

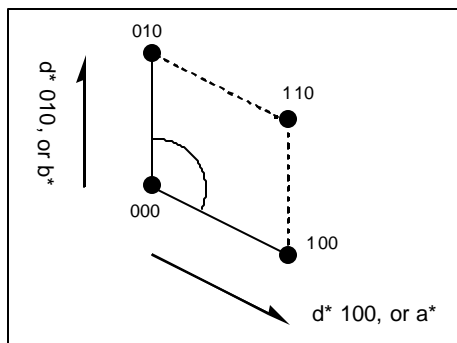
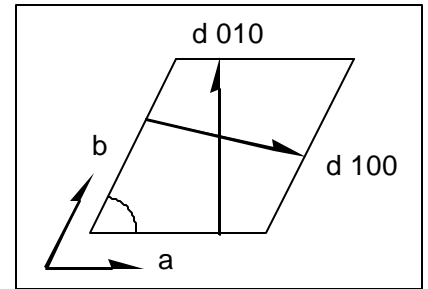
The exercise yields a repeating pattern of reciprocal lattice points that is generated by

- (a) constructing normals to families of planes that pass through one unit cell,
- (b) calculating the reciprocal of the length of each normal - and from a common origin,
- (c) marking the end of the reciprocal length after
- (d) positioning it in the same orientation as the original normal.

In real space, suppose the angle between two cell axes (say a and b) is not 90 degrees, then, the normal between parallel cell edges is given by $a \sin g$ and $b \sin g$ respectively.

These particular normals are labeled as d_{100} and d_{010} respectively.

In a real space, cell limits are marked by an imaginary boundary selected so it contains the unique chemical/crystallographic information of its contents.



In reciprocal space, a unit cell may also be recognized. It has edge lengths that are the reciprocals of the normals of the real space cell, and they are labeled as d^*_{100} and d^*_{010} , or a^* and b^* respectively.

In reciprocal space, cell limits are marked by lattice points which represent families of planes. Each lattice point can be uniquely labeled with a set of indices.

Accordingly, the relationship between so-called real and reciprocal space for this 2-D examples is as follows:

real cell	normals	reciprocal cell
edge lengths		edge lengths
a, b	$d_{100} = a \sin g$	$a^* = d^*_{100} = \frac{1}{d_{100}}, \quad b^* = d^*_{010} = \frac{1}{d_{010}}$
	$d_{010} = b \sin g$	
cell area	$A = ab \sin g$	$A^* = a^* b^* \sin g^* \quad g^* = 180 - g$

for a monoclinic unit cell in three dimensional space:

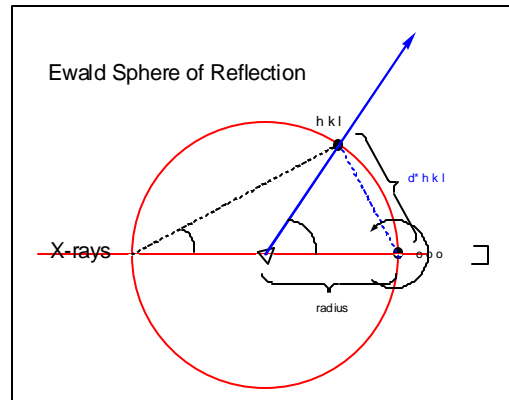
$$a \neq b \neq c, \quad a = g = 90 \neq b \quad a^* = 1/a \sin g, \quad b^* = 1/b, \quad c^* = 1/c \sin g$$

$$\text{cell volume } V = abc \sin g \quad \frac{1}{V} = V^* = a^* b^* c^* \sin g^*$$

Suffice it to say the reciprocal lattice is observed as a diffraction pattern. A two-dimensional example of a reciprocal lattice was prepared for today's class. Of course a three-dimensional crystal with corresponding 3-D unit cell will produce a 3-D diffraction pattern associated with a 3-d reciprocal lattice. It is convenient to examine the 3-D lattice in terms of separate "levels" or "slices" in which one index is held constant. So, for example, reflections of the type $hk0$ for which the l index is zero could be grouped and considered. This yields a 2-D pattern similar to the one prepared for today's class. The complete set of reflections would be obtained by collecting additional levels having $l = 1, 2, 3, \dots$

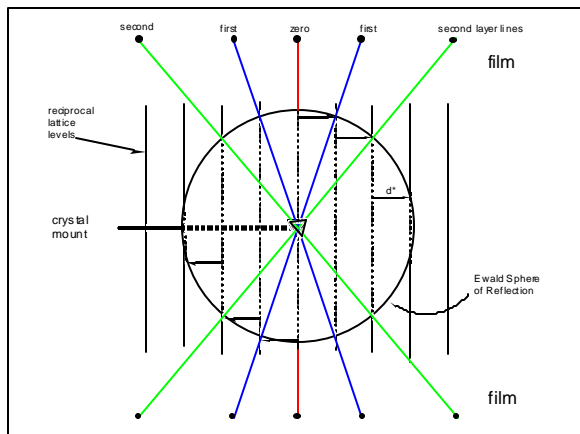
Diffraction: in order to "reflect" x-rays of wavelength λ a family of planes (hkl) separated by a normal (d_{hkl}) must be oriented at an angle q w/r to the x-ray beam, so that Bragg's Law is satisfied: $2d_{hkl} \sin q = \lambda$.

This is accomplished in reciprocal space by invoking a model known as Ewald's sphere of reflection. The origin of the reciprocal lattice (000) is fixed at the point where x-rays exit the sphere. The radius of the sphere is $1/\lambda$. (A triangle whose hypotenuse is a diameter is a right triangle.)



As the crystal is moved in the x-ray beam its associated reciprocal lattice also moves correspondingly. Consider the situation when a reciprocal lattice point is in contact with the sphere... It was demonstrated that Bragg's law is obeyed in such situations.

Rotation Method: is one of the first methods used to detect the reciprocal lattice. DEMO.



A real axis is perpendicular to parallel reciprocal lattice layers.

