

Condensed Matter Physics I

Prof. Dr. Ir. Paul H.M. van Loosdrecht
II Physikalisches Institut, Room 312
E-mail: pvl@ph2.uni-koeln.de

Website: <http://www.loosdrecht.net/>

Lectures and Working Class

Lectures & problem session on

Tuesday 10:00 : 11:30

Thursday 12:00 : 13:30

Problem sessions on Thursday alternating with lectures

Problem session teacher:

Dr. Matteo Montagnese, II P.I. room 318

m.montagnese@ph2.uni-koeln.de

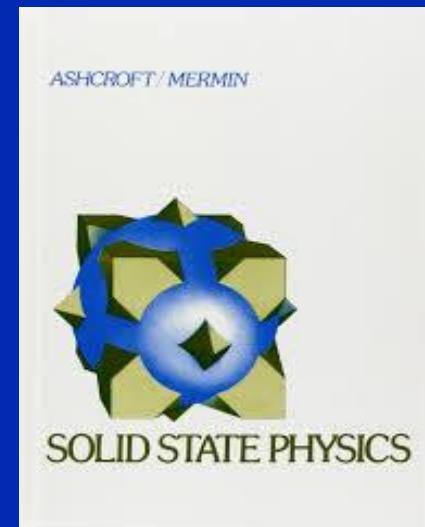
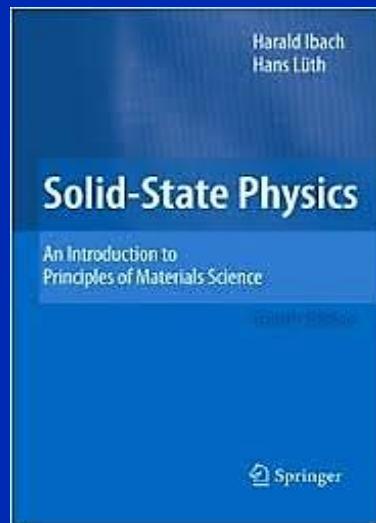
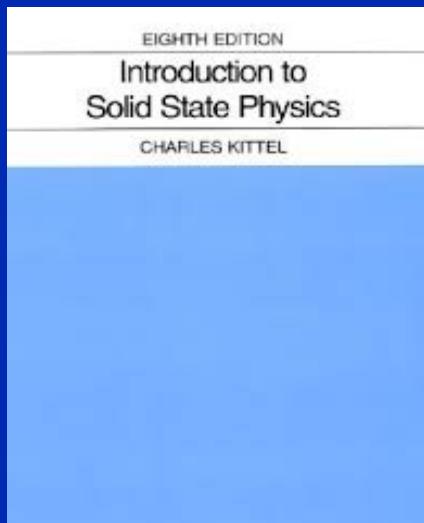
Hand in problem sets on Tuesday's (room 318).

Literature

Introduction to Solid State Physics, Charles Kittel

Solid State Physics, Ibach & Lüth

Solid State Physics, Ashcroft & Mermin



Quantum theory of solids, Kittel

Principles of condensed matter physics, Chaikin & Lubensky

Many, many more

States of matter

Plasma

Gas

Liquid

Liquid crystals

glasses

Amorphous solids

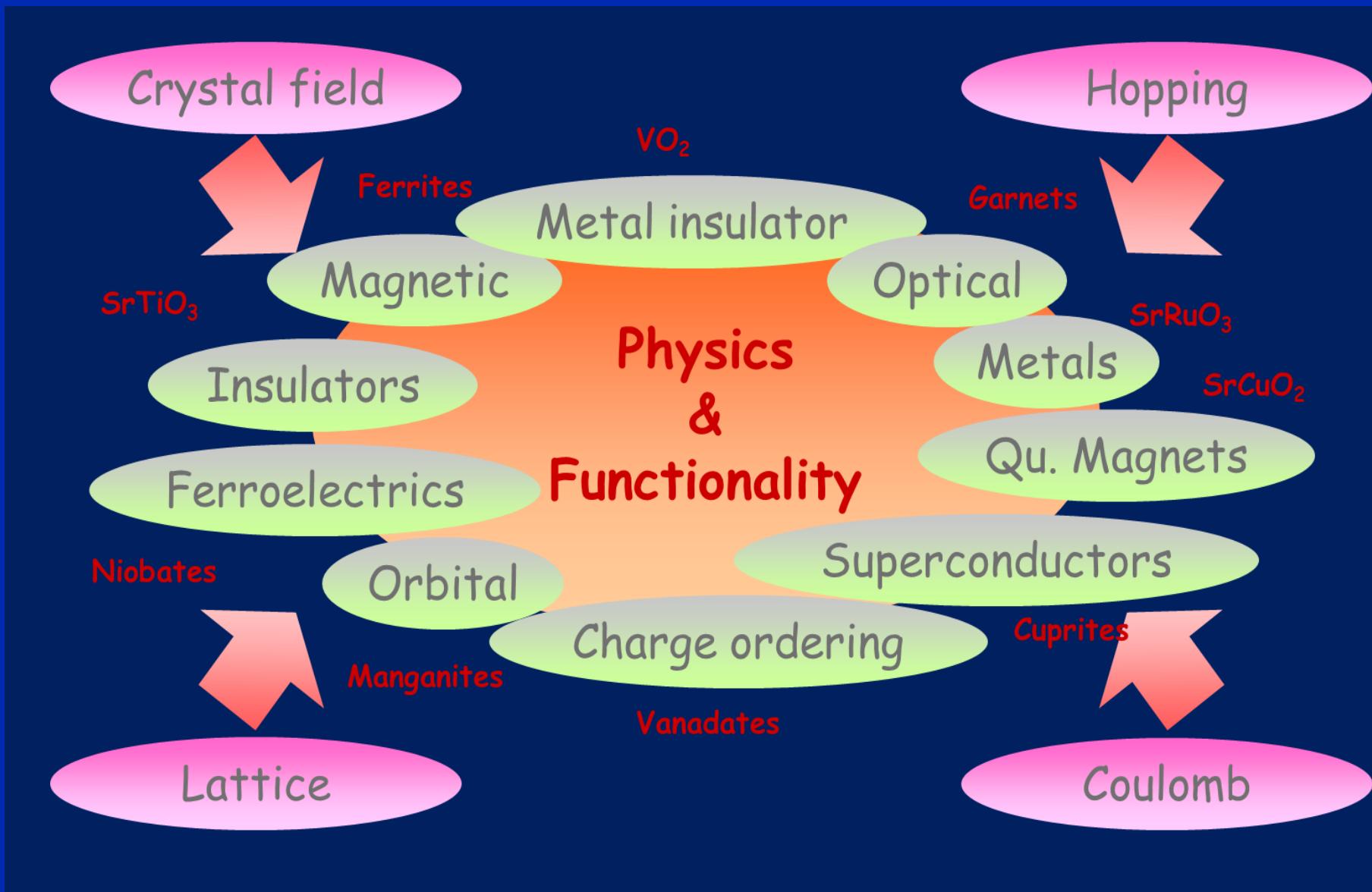
quasi crystals

nearly periodic crystals

PERIODIC CRYSTALS

- Structure and symmetry
- Binding
- Diffraction -> reciprocal space
- Vibrations & thermal properties, density of states
- Electrons, Fermi Gas, Energy bands, semiconductors
- Collective excitations (plasmons, excitons, etc.)
- Dielectrical properties
- Spins: magnetism & magnetic excitations

II. Physikalisches Institut



STRUCTURE

Kittel Ch.1



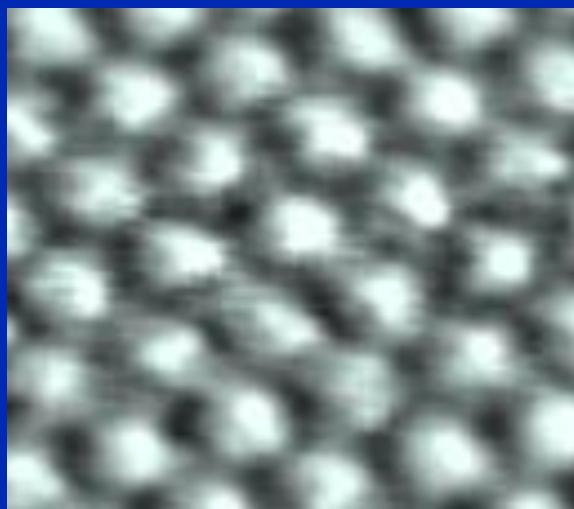
Quartz



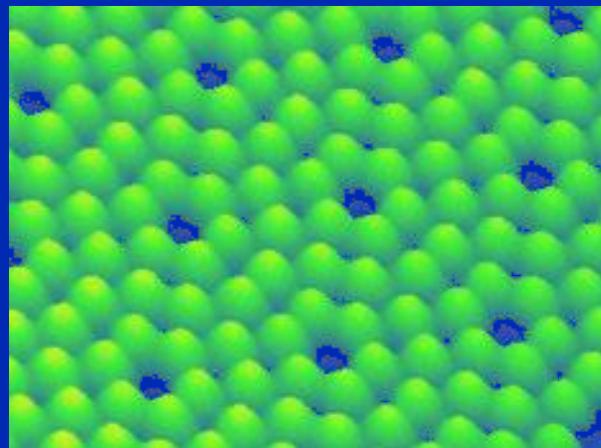
Protein



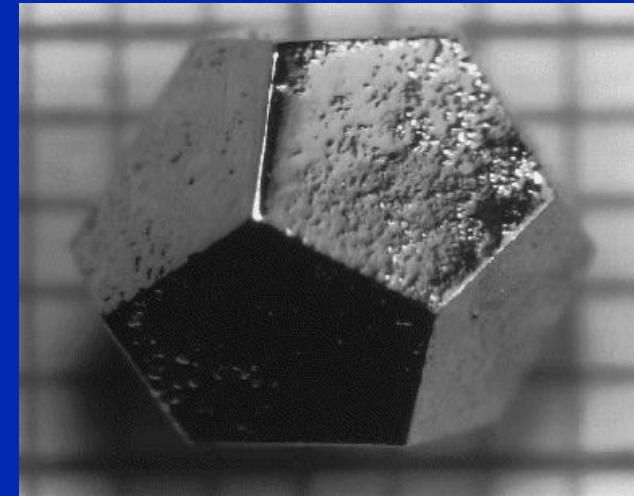
Cu_2O



Graphite



Silicon



Quasi-crystal

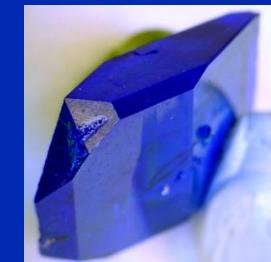
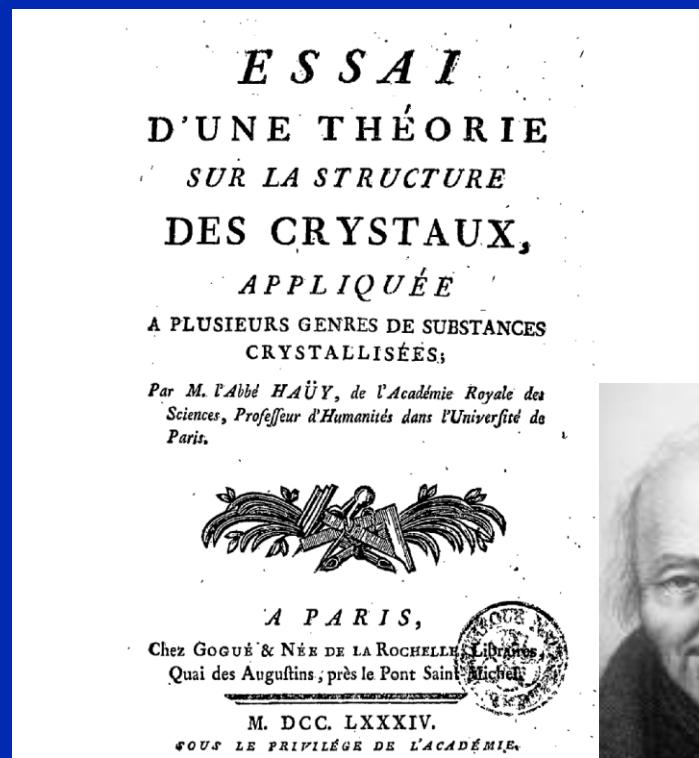
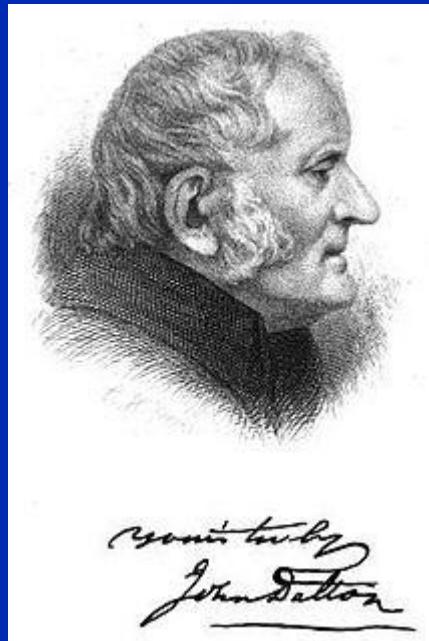
Structure

460 BC Demokritus: idea that matter is made of 'atoms'

1611 J. Keppler: Snow Xtals made of spherical water molecules

1784 R.J. Haüy: Law of integral indices

1802 J. Dalton: atomic theory



Structure

1784 R.J. Haüy: Law of integral indices



1850 A. Bravais: Space lattices



1895 W. Röntgen: X-rays (1st Nobel prize 1901,
Work on cathode rays: Lenard 1905)

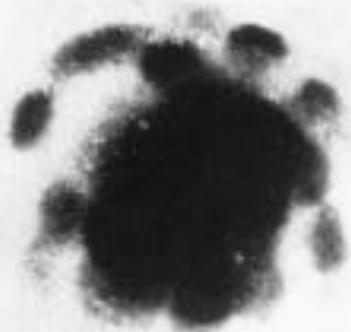


1912 M. van Laue: Diffraction (Nobel prize 1914)

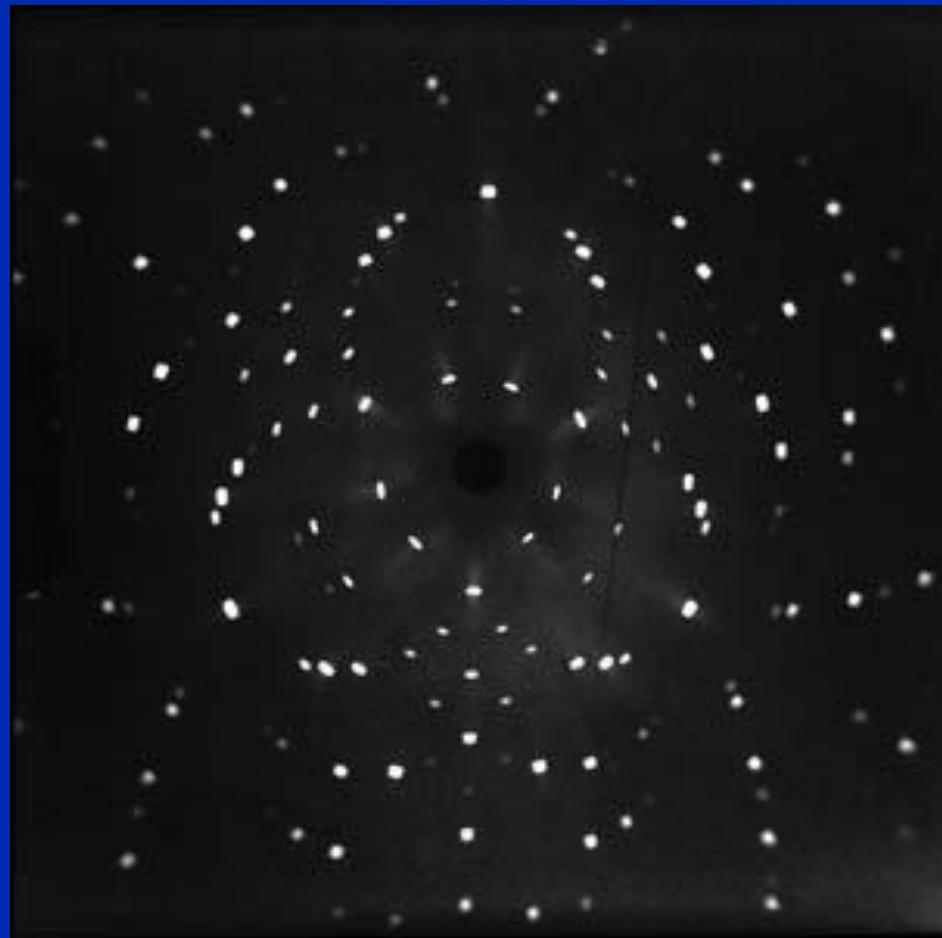


Beugungsexperimente

Die erste Röntgen-Durchbeleuchtung eines
Kristalls.



M.v. Lahn

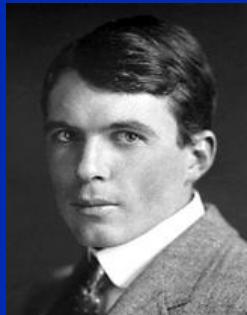
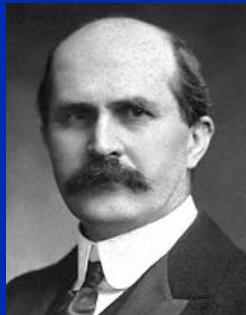


Si (FCC) 111 axis

Using hard X-rays ($\lambda \sim a \sim 10 \text{ \AA}$)

Structure

1913/14 W.&L. Bragg: Diffraction (Nobel prize 1915)



1984 D. Shechtman: Quasicrystals (Chem. Nobel prize 2011)



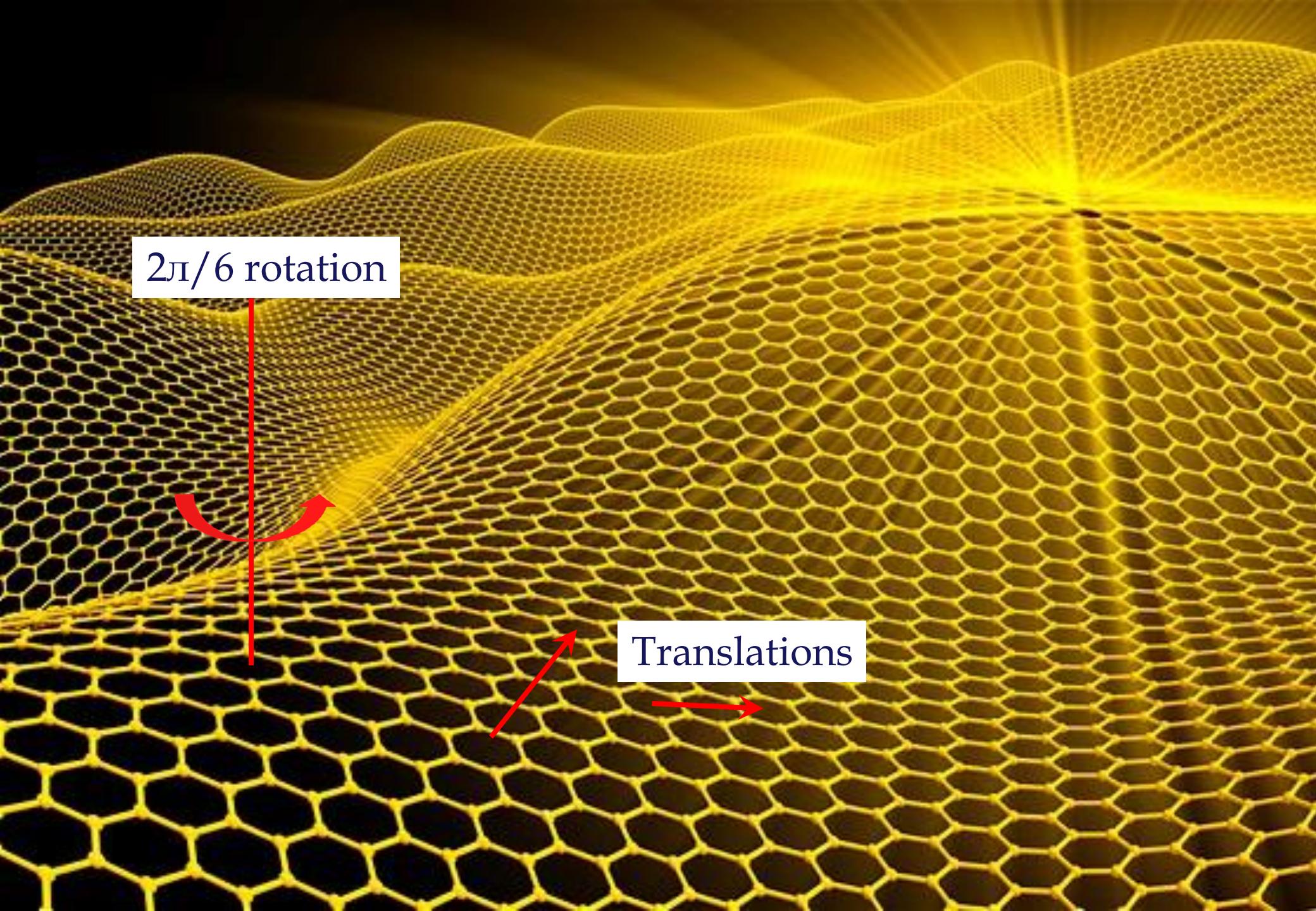
Symmetry

Translation symmetry:

vectors representing translations which transform a structure into itself

Rotation symmetry:

Operations (rotations, mirrors) which transform a structure into itself



$2\pi/6$ rotation

Translations

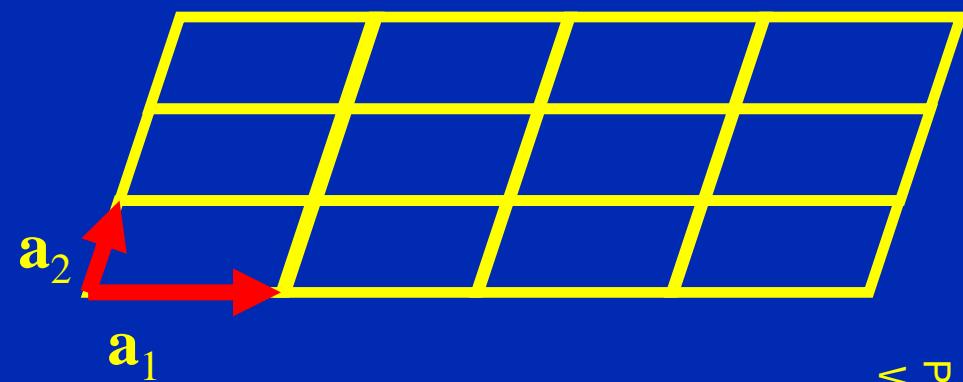
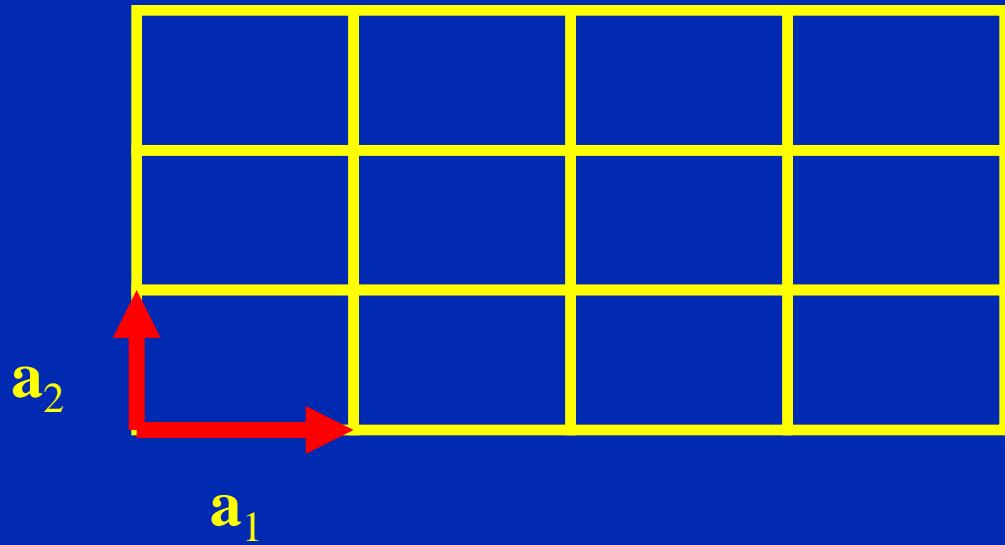
Lattice

Periodic Arrays of points

or

Lattice = set of fundamental translation vectors:

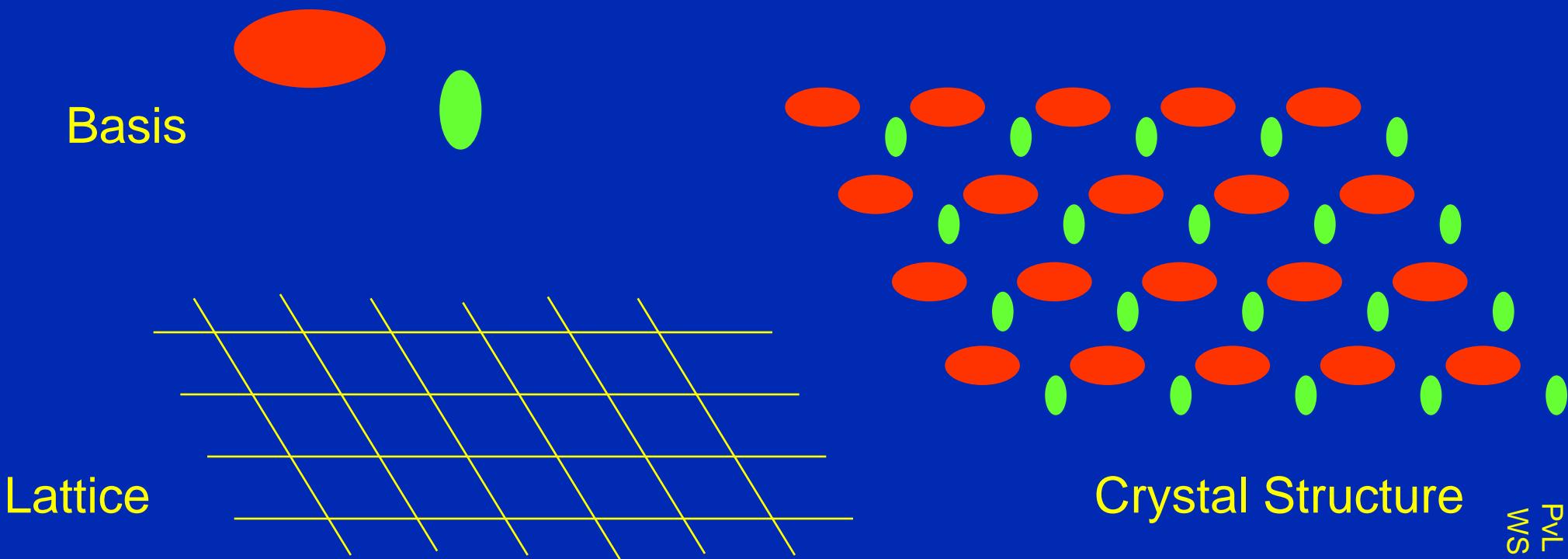
$$\mathbf{T} = u_1 \mathbf{a}_1 + u_2 \mathbf{a}_2 + u_3 \mathbf{a}_3$$



Crystal structure

Lattice + Basis = Crystal structure

Basis = group of atoms forming the unit cell:

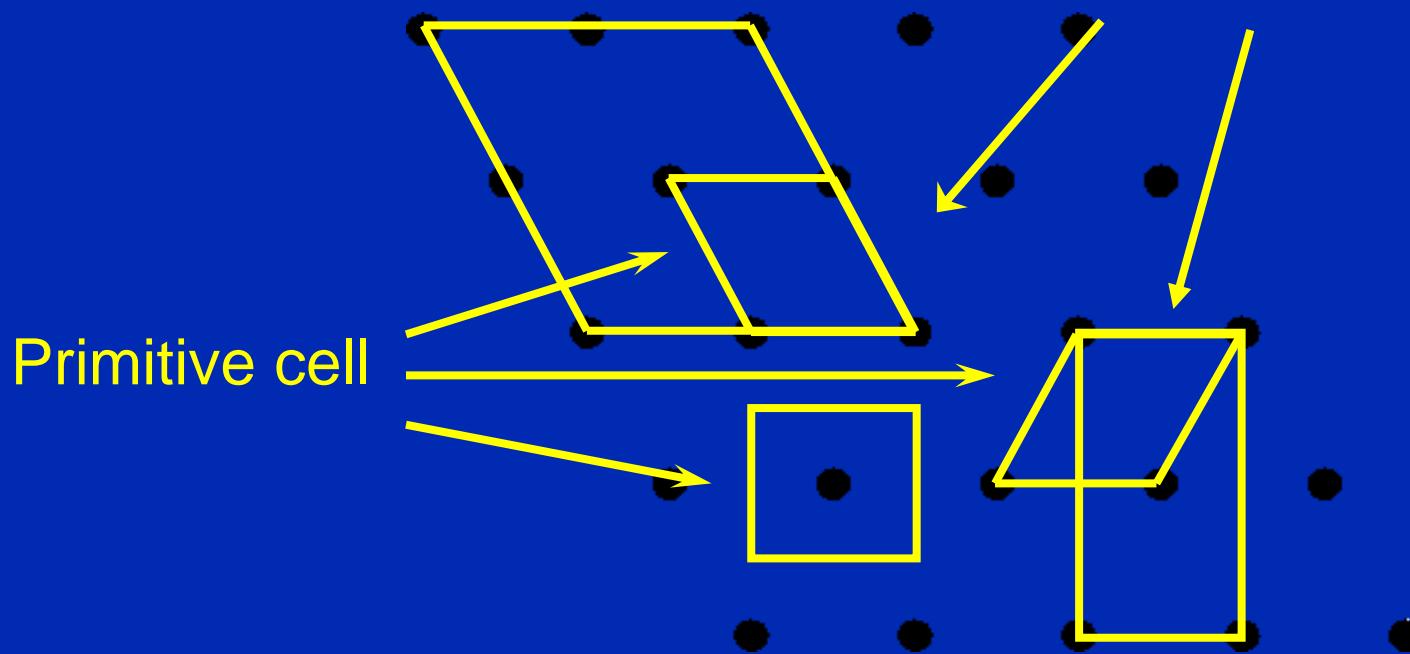


Primitive cell

Minimal volume unit cell

“Conventional” unit cell

A unit cell is not necessarily primitive!



Primitive cell

Minimal volume unit cell

Primitive Translation vectors: $\mathbf{a}_1, \mathbf{a}_2, \mathbf{a}_3$

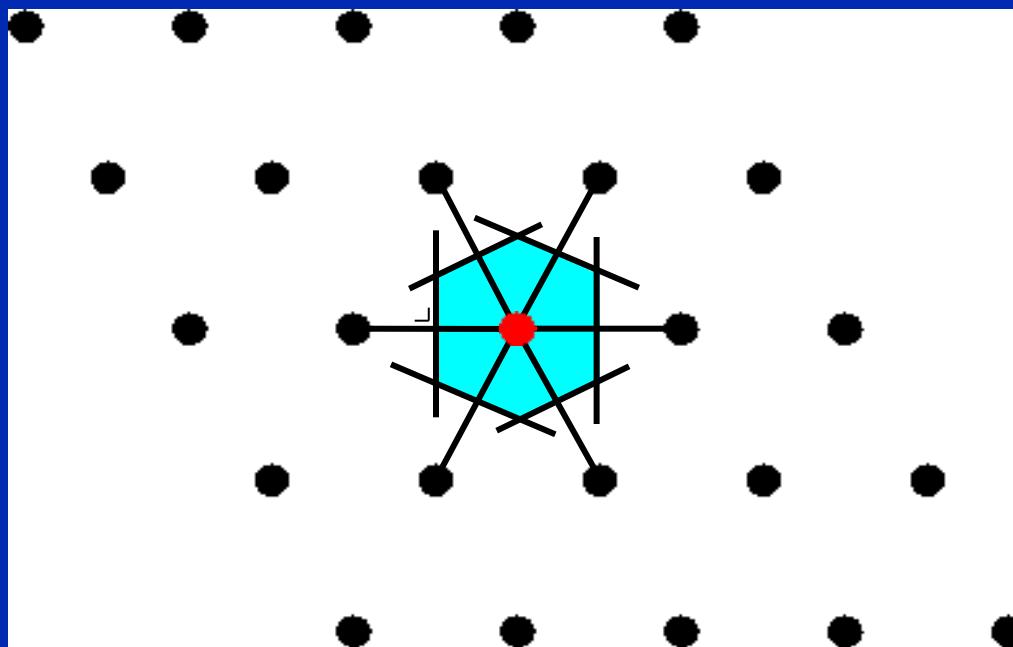
Primitive Lattice Cell = parallelepiped of $\mathbf{a}_1, \mathbf{a}_2, \mathbf{a}_3$

Many equivalent options to choose from,

Volume of Primitive Lattice cell: $V = \mathbf{a}_1 \cdot (\mathbf{a}_2 \times \mathbf{a}_3)$

Wigner-Seitz Cell: Most symmetric choice of Primitive Cell

Wigner-Seitz cell



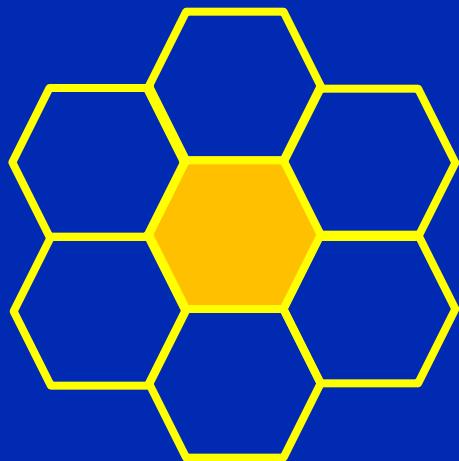
Symmetry elements

Translations: $\mathbf{T} = u_1\mathbf{a}_1 + u_2\mathbf{a}_2 + u_3\mathbf{a}_3$

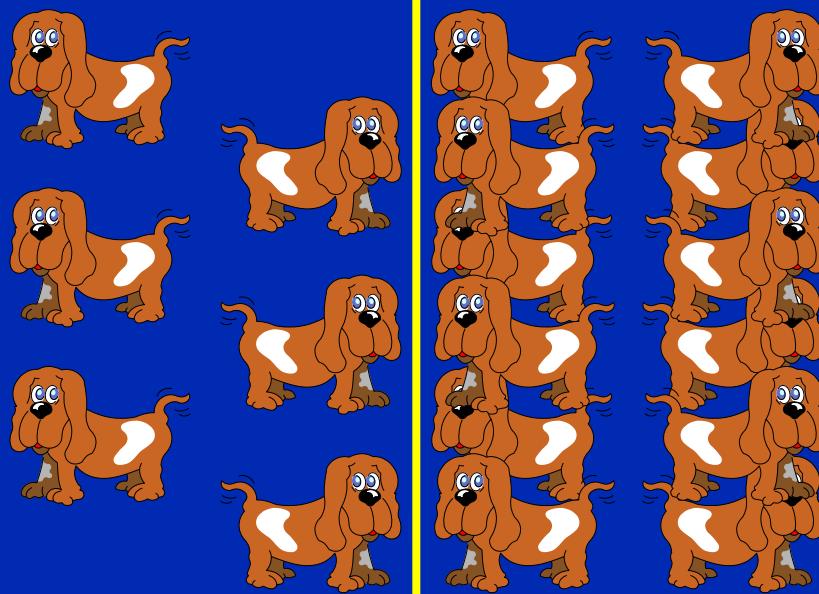
Point group operations: Rotations and Mirrors

Combinations: Glide planes and Screw axes

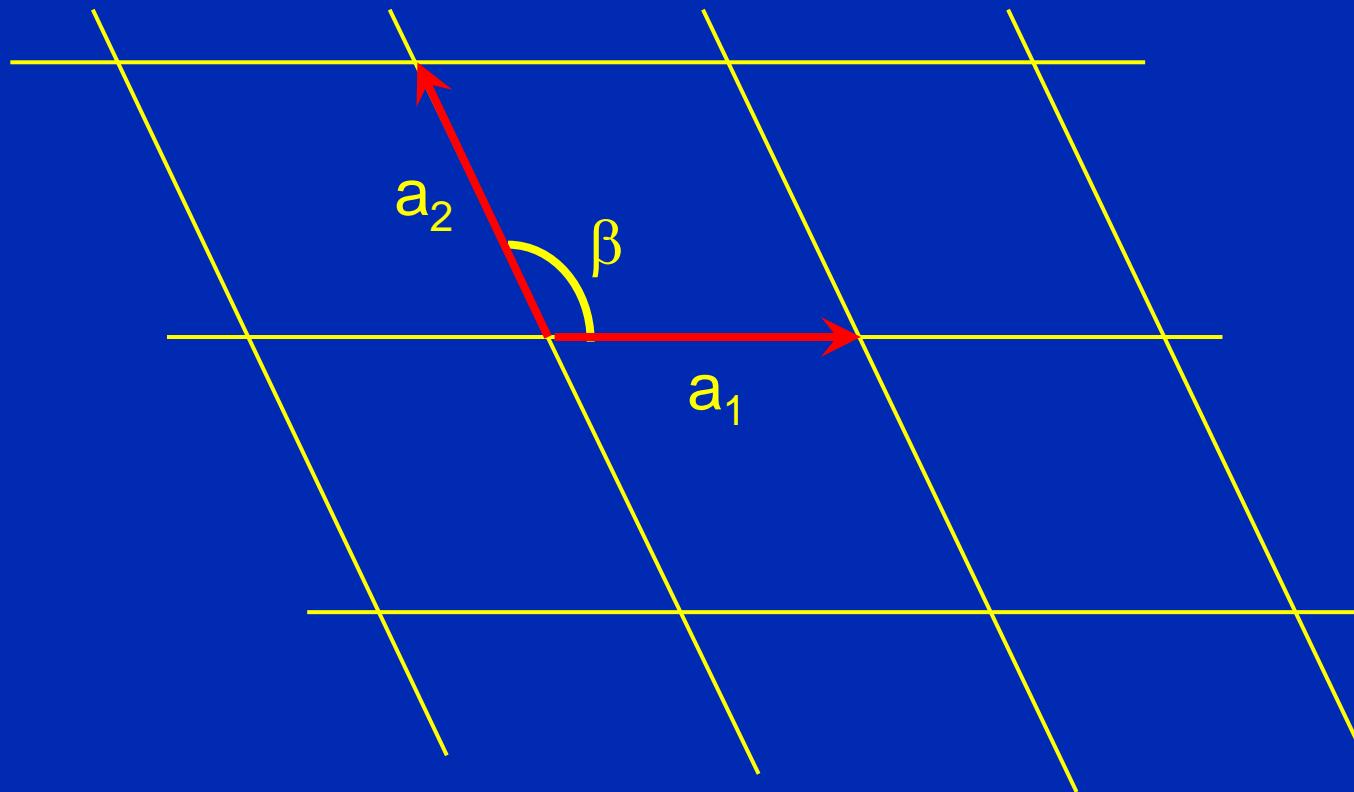
6 fold rotation



Glide plane



Two dimensional lattices (Net)



Ratio $|a_1/a_2|$
Angle β

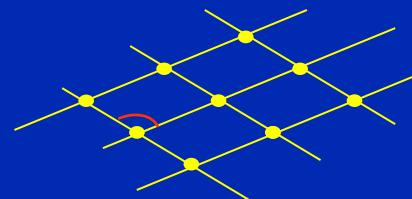
Two dimensional lattices (Net)



2D lattice types

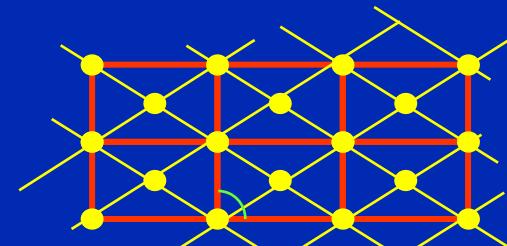
1-fold: Oblique

$$a_1 \neq a_2; \beta$$



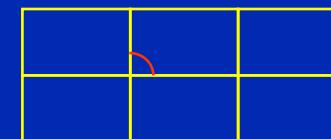
1-fold + mirror:
Centered rectangular

$$a_1 \neq a_2; \alpha = 90^\circ$$



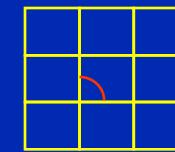
2-fold: Rectangular

$$a_1 \neq a_2; \beta = 90^\circ$$



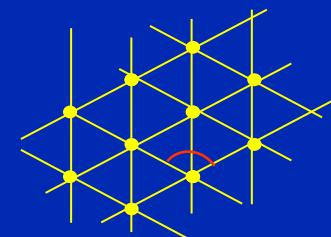
4-fold: Square

$$a_1 = a_2; \beta = 90^\circ$$



6-fold: Hexagonal

$$a_1 = a_2; \beta = 120^\circ$$



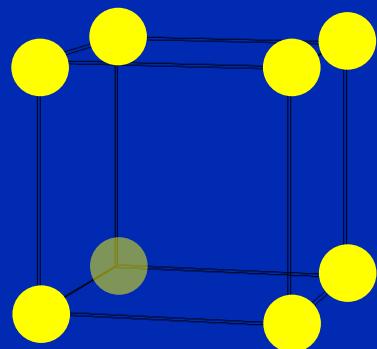
3D lattice types

The 3D lattice types (14)

Triclinic	1	No symmetries
Monoclinic	2	Two right angles
Orthorhombic	4	Three right angles
Tetragonal	2	Three right angles + 4 fold
Cubic (SC, FCC, BCC)	3	Three right angles + 4 fold + 3 fold
Trigonal	1	Three equal angles (not 90°) + 3 fold
Hexagonal	1	Two right, one 120° + 6 fold

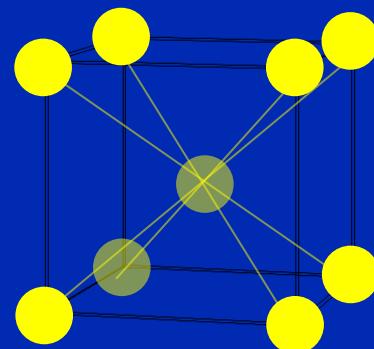
Cubic lattices

Simple cubic
SC



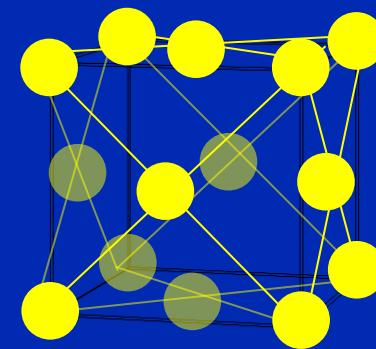
Po

Body centered
BCC



Na, K, Rb, Cs,
Fe, V, Cr, Mo, W

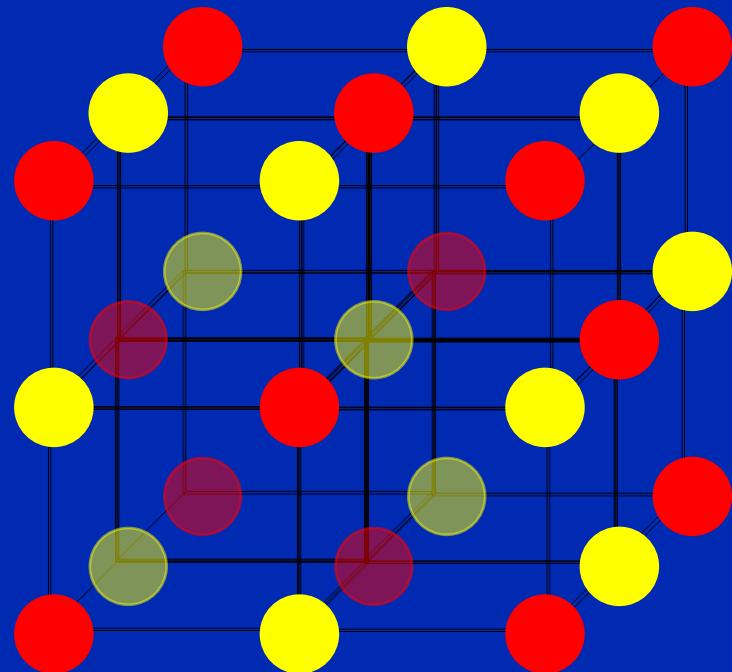
Face centered
FCC



Ni, Cu, Ag, Au, Pt,
Ne, Ar, Kr, Xe, Rn

NaCl Structure

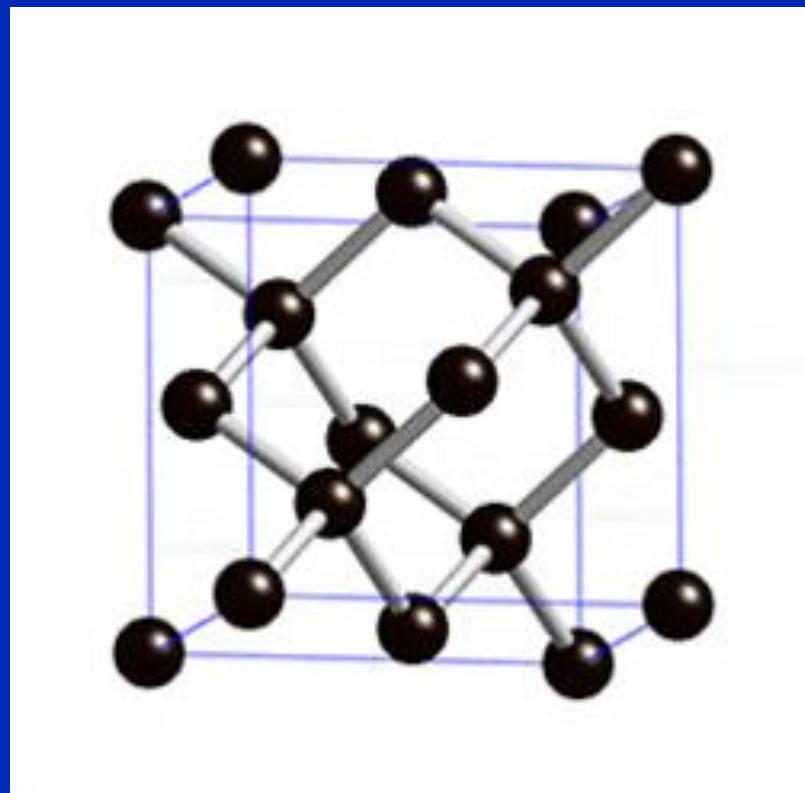
Rocksalt structure



NaCl
SnTe
NaF
MgO

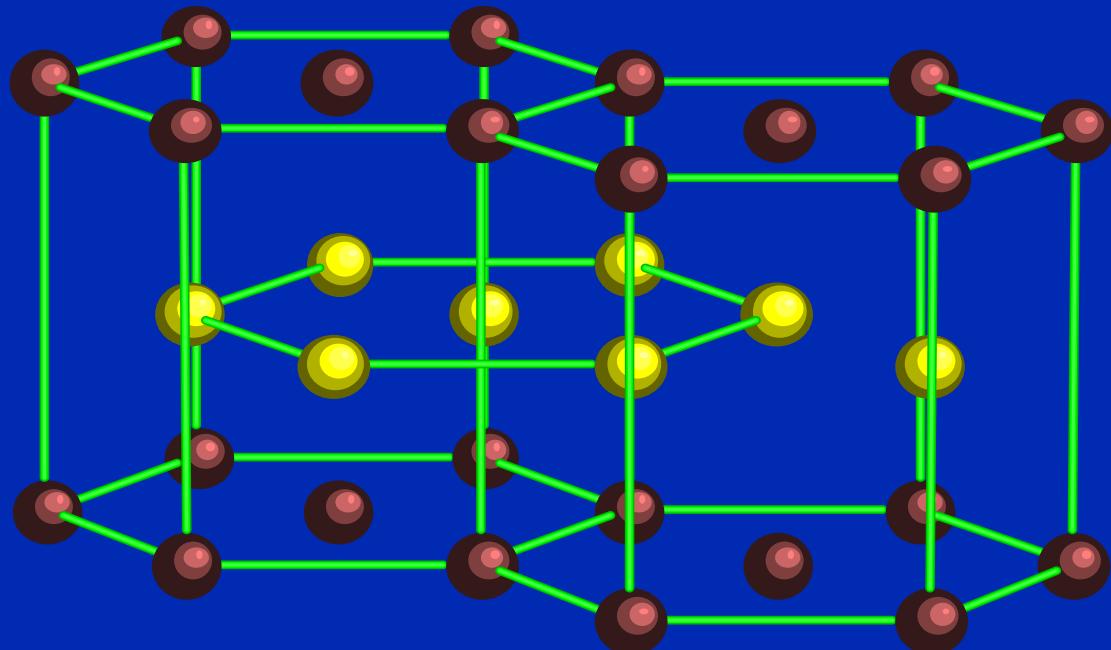


Diamond structure



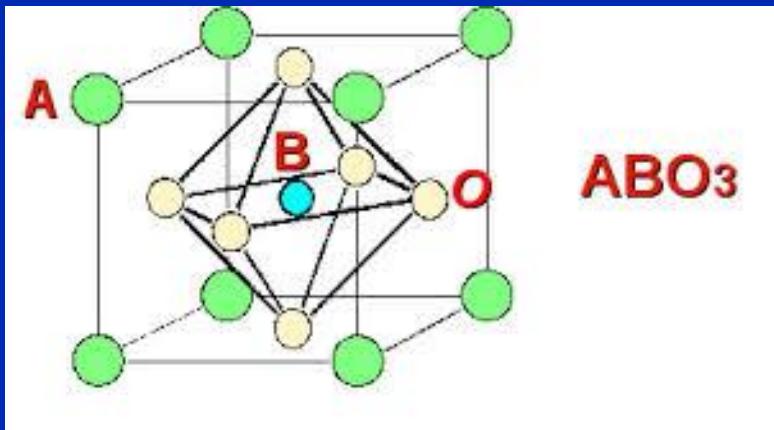
C,Si,Ge

Hexagonal closed packed



H, He, N
Be, Mg
Co, Zn, Sc, Ti
Y, Zr, Tc, Ru, Cd
Hf, Re, Os, Tl,
Pr, Gd, Tb, Dy, Ho, Er, Lu

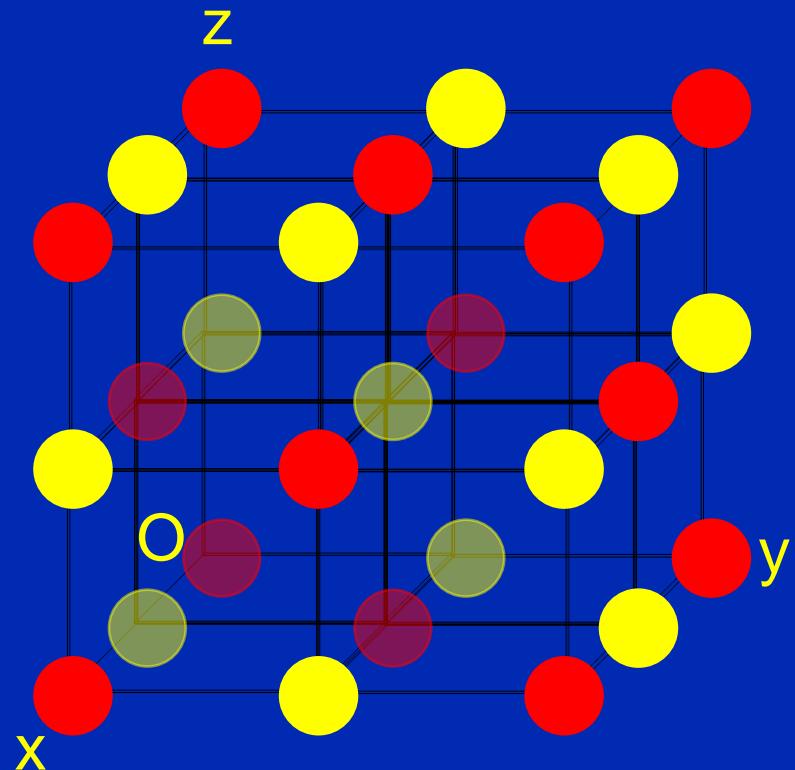
Perovskite structure



CaTiO₃



Planes and directions



Miller indices:

1. Find intercepts with axes of unit cell (in cell units).
2. Take reciprocals
3. Find (smallest) integers with the same ratio

() Planes

{ } Equivalent planes

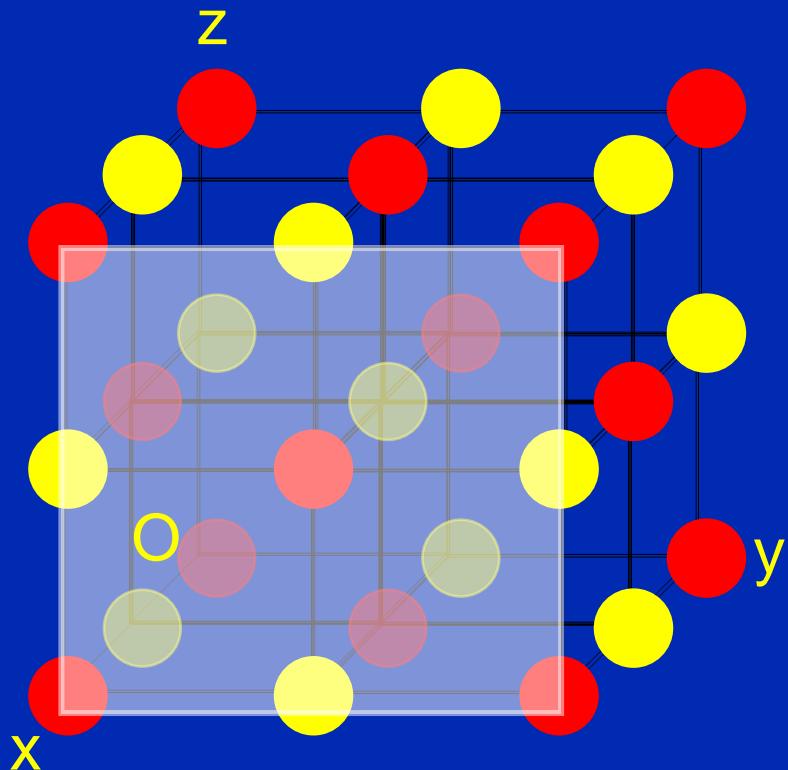
Cubic

{100}: (100),(200),(001)

[] directions:

x-direction [100]

Planes and directions

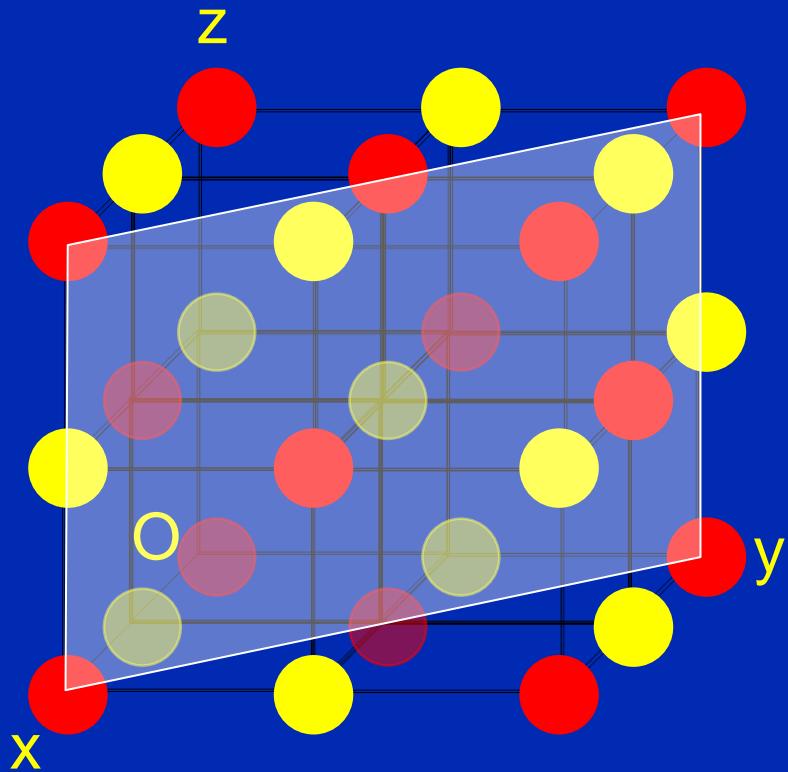


Miller indices:

1. Find intercepts with axes of unit cell (in cell units).
2. Take reciprocals
3. Find (smallest) integers with the same ratio

Intersections at $(1, \infty, \infty)$
Miller indices (100)

Planes and directions

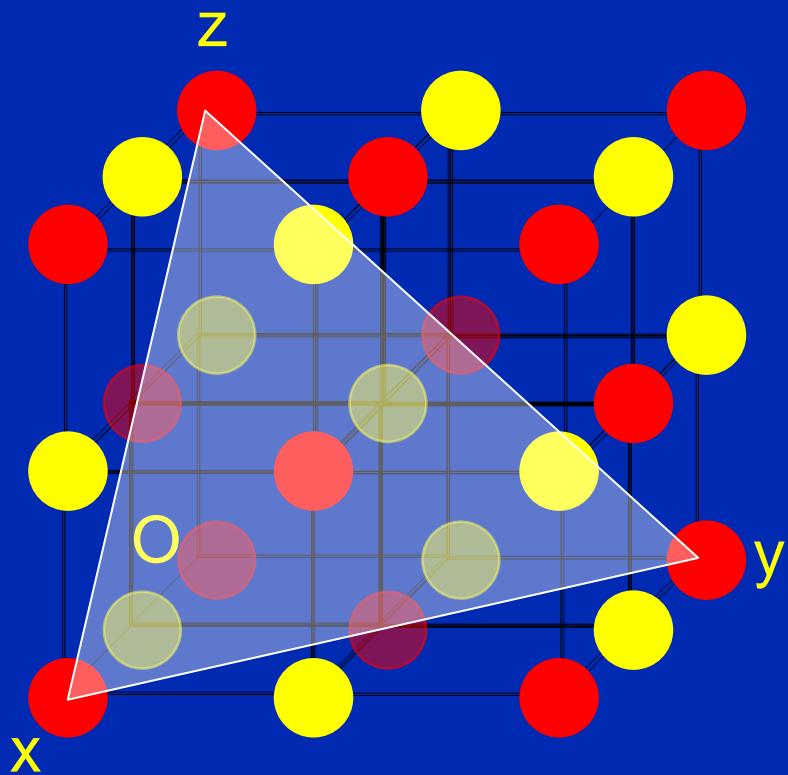


Miller indices:

1. Find intercepts with axes of unit cell (in cell units).
2. Take reciprocals
3. Find (smallest) integers with the same ratio

Intersections at $(1, 1, \infty)$
Miller indices (110)

Planes and directions



Miller indices:

1. Find intercepts with axes of unit cell (in cell units).
2. Take reciprocals
3. Find (smallest) integers with the same ratio

Intersections at (1,1,1)
Miller indices (111)