

Day 2 pg 1

Prayer

Survey - put notes on website?

Comment - reading assignments

handouts for Yangping

Vocabulary review (call on students)

demo request
Stokes crystal models

crystal lattice (14 Bravais lattices)

lattice vectors

primitive lattice vectors

unit cell

primitive unit cell

Wigner-Seitz cell

conventional unit cell

→ I forgot to mention this (see day 1 pg 3 notes)

basis

symmetry operation

translation

point operations, or pt symmetries

rotation

→ forgot to mention notation

reflection

ex. σ_{xy}

inversion

ex. S_6

I

C_{4x}^+ , C_{2y}^- , etc

space group (230 space groups)

crystal class (32 classes)

sc

bcc

fcc

2D lattices, cont from last time

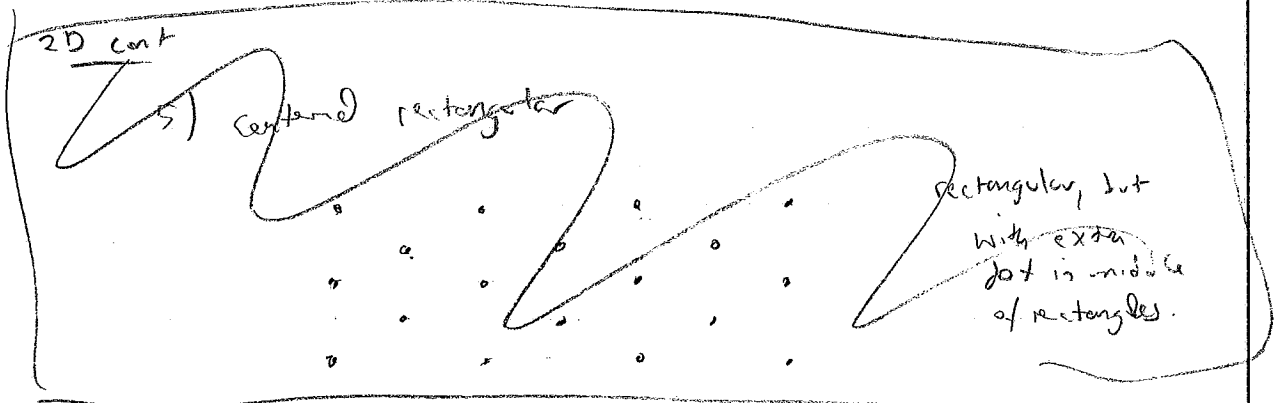
- why no centered square lattice?

(see next page)

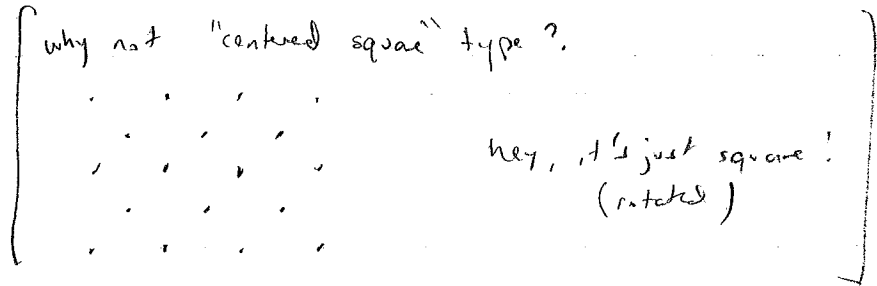
- why no pentagonal lattice?

(see overhead

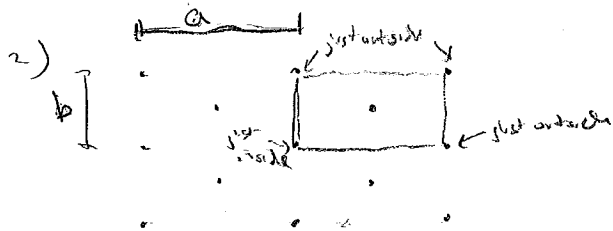
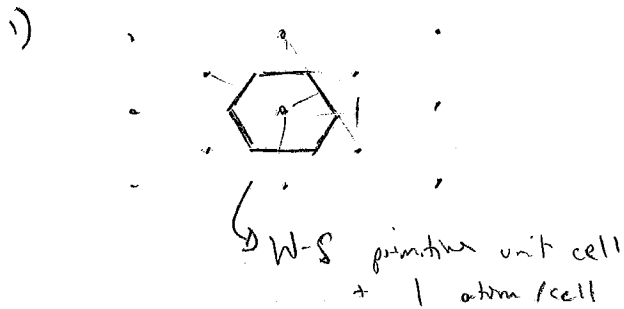
No translational symmetry!)



Ans. Day 2



Two ways of looking at centered rectangular w/ 1 atom/cell



$$\vec{r}_1 = (0, 0)$$

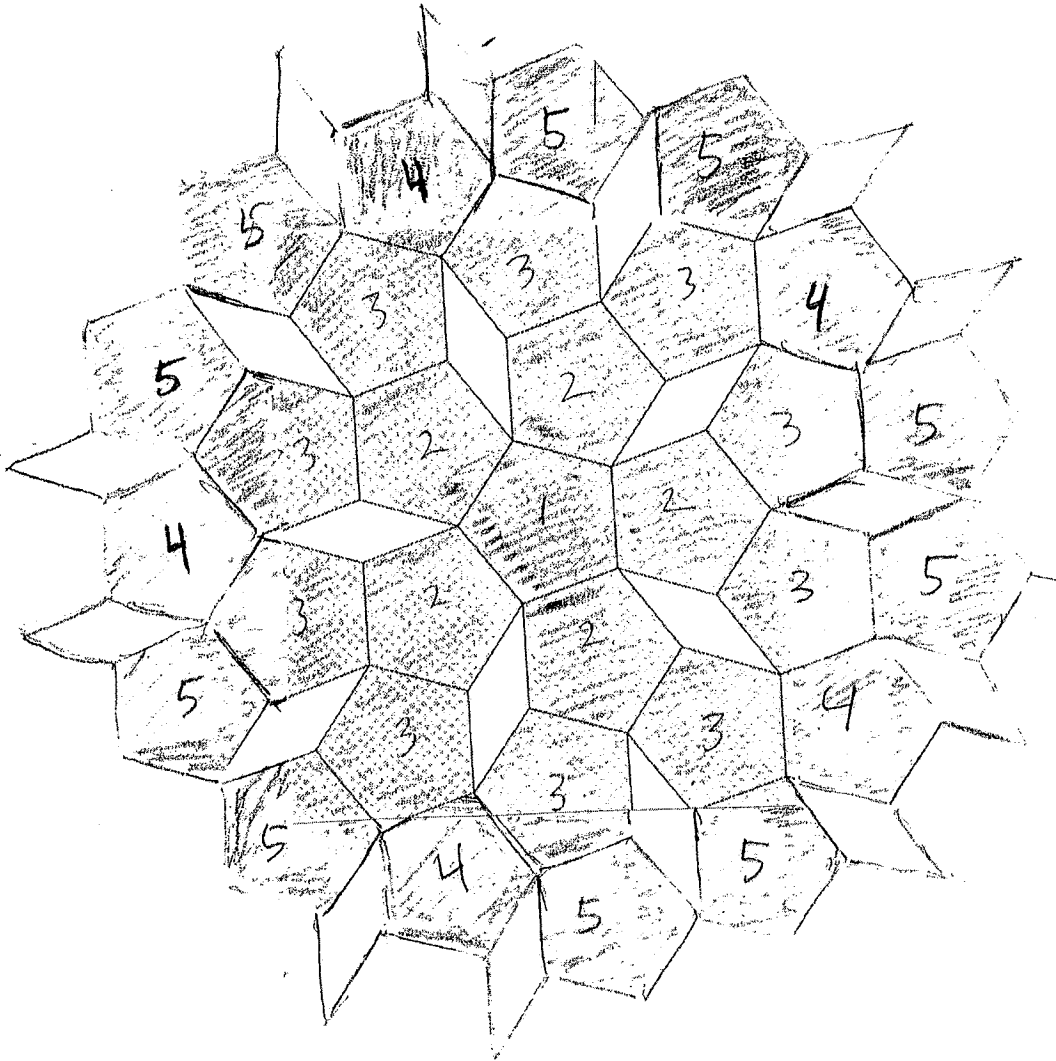
$$\vec{r}_2 = (a/2, b/2)$$

in this example.

which is easier?

Pentagonal Tiling of 2D Plane
Physics 581 Colton

From Kittel Fig. 1.5, pg 7



One more notation note - [] vs () vs { }

[] a direction. Eg $[100] = \hat{x}$
 $[110] = \frac{\hat{x} + \hat{y}}{\sqrt{2}}$

() a single plane or set of parallel planes "Miller indices"

Defined by intercepts on primitive l.v. lines

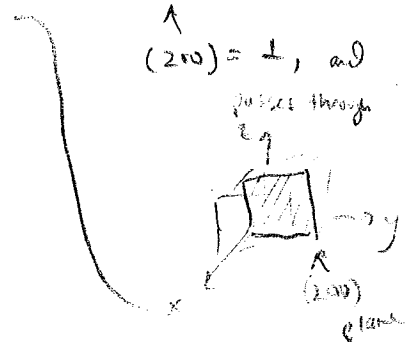
If plane intersects axes at: $A_1 \hat{a}_1$
 $A_2 \hat{a}_2$
 $A_3 \hat{a}_3$

take inverse $(\frac{1}{A_1}, \frac{1}{A_2}, \frac{1}{A_3})$

multiply by LCD
 $\rightarrow (h, k, l)$

careful - for (simple) cubic the planes are exactly what you'd expect: $(100) = \perp$ to $[100]$ direction
 but not for other systems

Convention ~~is~~
 is to use abc for
 use cubic planes



If representing a set of planes, $(200) =$



{ } planes equivalent by symmetry
 Ex. in sc, $\{100\} = (100), (010), \text{ and } (001)$ planes

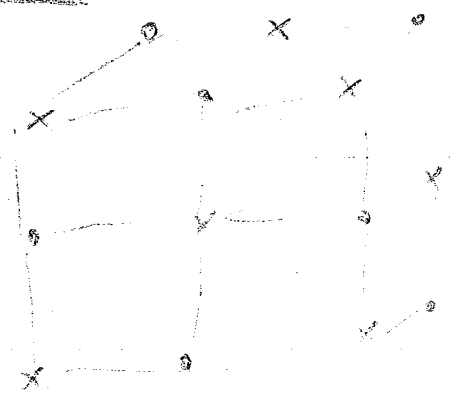
"bcc structure" vs "bcc lattice"

Some real materials! (f.c.c.)

Elements pg 20*

- * bcc: Li, Na, K, V, Cr, Fe, Rb, Nb, Mo, Cs, Ba, Ta, W, (Eu)
- * fcc: Ne, Al, Ar, Ca, Ni, Cu, Kr, Sr, Rh, Pd, Ag, Xe, Ir, Pt, Au, Pb, Ac, Ce, Yb, Th

Sodium chloride



~~lattice~~

~~more for discussion~~
~~the cause of~~
~~relaxation~~

<http://cst-www.nrl.navy.mil/lattice>

also Stille's website

↳ links at bottom of class website

Lattice = ? fcc, not sc

basis: Cl⁻ ion at 000
Na⁺ ion at 1/2 1/2 1/2

8 atoms in conventional unit cell

"NaCl structure" = name of this structure & also used by other crystals

but list:	LiH	AgBr
	MgO	PbS
	MnO	KCl
	(NaCl)	KBr

Note 1:1 stoichiometry