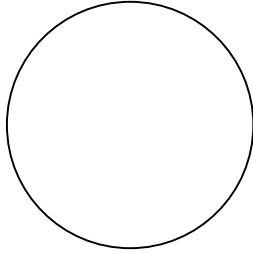


Name (pinyin please) _____

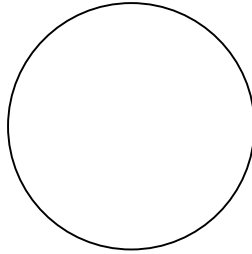
Homework #2. If you get your work to me by Monday evening, we can discuss it in class Tuesday morning.....before the exam on Wednesday.

It will be OK if you want to email your work to me at rx7@psu.edu

1. Construct the two point group diagrams for the point group 32.



symmetry elements



equivalent points

2. CaF_2 is $Fm\bar{3}m$, $a = 5.462 \text{ \AA}$, Ca in 4a, F in 8c. Make a drawing of the crystal structure projected down $[001]$. The equipoints for space group $Fm\bar{3}m$ are listed on the next page.

The F atoms have how many Ca nearest neighbors (i.e., how many Ca atoms touch each F atom)?

Since $a = 5.462 \text{ \AA}$, what is the Ca-F bond distance? (A simple geometry calculation.)

Positions		Coordinates					
		(0,0,0)+	(0, $\frac{1}{2}$, $\frac{1}{2}$)+	($\frac{1}{2}$,0, $\frac{1}{2}$)+	($\frac{1}{2}$, $\frac{1}{2}$,0)+		
Multiplicity,	Wyckoff letter,						
Site symmetry							
192	<i>l</i> 1	(1) x, y, z (5) z, x, y (9) y, z, x (13) y, x, \bar{z} (17) x, z, \bar{y} (21) z, y, \bar{x} (25) $\bar{x}, \bar{y}, \bar{z}$ (29) $\bar{z}, \bar{x}, \bar{y}$ (33) $\bar{y}, \bar{z}, \bar{x}$ (37) \bar{y}, \bar{x}, z (41) \bar{x}, \bar{z}, y (45) \bar{z}, \bar{y}, x	(2) \bar{x}, \bar{y}, z (6) z, \bar{x}, \bar{y} (10) \bar{y}, z, \bar{x} (14) $\bar{y}, \bar{x}, \bar{z}$ (18) \bar{x}, z, y (22) z, \bar{y}, x (26) x, y, \bar{z} (30) \bar{z}, x, y (34) y, \bar{z}, x (38) y, x, z (42) x, \bar{z}, \bar{y} (46) \bar{z}, y, \bar{x}	(3) \bar{x}, y, \bar{z} (7) \bar{z}, \bar{x}, y (11) y, \bar{z}, \bar{x} (15) y, \bar{x}, z (19) $\bar{x}, \bar{z}, \bar{y}$ (23) \bar{z}, y, x (27) x, \bar{y}, z (31) z, x, \bar{y} (35) \bar{y}, z, x (39) \bar{y}, x, \bar{z} (43) x, z, y (47) z, \bar{y}, \bar{x}	(4) x, \bar{y}, \bar{z} (8) \bar{z}, x, \bar{y} (12) \bar{y}, \bar{z}, x (16) \bar{y}, x, z (20) x, \bar{z}, y (24) $\bar{z}, \bar{y}, \bar{x}$ (28) \bar{x}, y, z (32) z, \bar{x}, y (36) y, z, \bar{x} (40) y, \bar{x}, \bar{z} (44) \bar{x}, z, \bar{y} (48) z, y, x		
96	<i>k</i> $\dots m$	x, x, z \bar{z}, \bar{x}, x x, x, \bar{z} $\bar{x}, \bar{z}, \bar{x}$	\bar{x}, \bar{x}, z \bar{z}, x, \bar{x} $\bar{x}, \bar{x}, \bar{z}$ x, \bar{z}, x	\bar{x}, x, \bar{z} x, z, x x, \bar{x}, z z, x, \bar{x}	x, \bar{x}, \bar{z} \bar{x}, z, \bar{x} \bar{x}, x, z z, \bar{x}, x	z, x, x x, \bar{z}, \bar{x} x, z, \bar{x} \bar{z}, x, x	z, \bar{x}, \bar{x} \bar{x}, \bar{z}, x \bar{x}, z, x $\bar{z}, \bar{x}, \bar{x}$
96	<i>j</i> $m \dots$	$0, y, z$ $\bar{z}, 0, y$ $y, 0, \bar{z}$ $0, \bar{z}, \bar{y}$	$0, \bar{y}, z$ $\bar{z}, 0, \bar{y}$ $\bar{y}, 0, \bar{z}$ $0, \bar{z}, y$	$0, y, \bar{z}$ $y, z, 0$ $y, 0, z$ $z, y, 0$	$0, \bar{y}, \bar{z}$ $\bar{y}, z, 0$ $\bar{y}, 0, z$ $z, \bar{y}, 0$	$z, 0, y$ $y, \bar{z}, 0$ $0, z, \bar{y}$ $\bar{z}, y, 0$	$z, 0, \bar{y}$ $\bar{y}, \bar{z}, 0$ $0, z, y$ $\bar{z}, \bar{y}, 0$
48	<i>i</i> $m \dots m 2$	$\frac{1}{2}, y, y$ $\bar{y}, \frac{1}{2}, y$	$\frac{1}{2}, \bar{y}, y$ $\bar{y}, \frac{1}{2}, \bar{y}$	$\frac{1}{2}, y, \bar{y}$ $y, y, \frac{1}{2}$	$\frac{1}{2}, \bar{y}, \bar{y}$ $\bar{y}, y, \frac{1}{2}$	$y, \frac{1}{2}, y$ $y, \bar{y}, \frac{1}{2}$	$y, \frac{1}{2}, \bar{y}$ $\bar{y}, \bar{y}, \frac{1}{2}$
48	<i>h</i> $m \dots m 2$	$0, y, y$ $\bar{y}, 0, y$	$0, \bar{y}, y$ $\bar{y}, 0, \bar{y}$	$0, y, \bar{y}$ $y, y, 0$	$0, \bar{y}, \bar{y}$ $\bar{y}, y, 0$	$y, 0, y$ $y, \bar{y}, 0$	$y, 0, \bar{y}$ $\bar{y}, \bar{y}, 0$
48	<i>g</i> $2 \dots m m$	$x, \frac{1}{2}, \frac{1}{2}$ $\frac{1}{2}, x, \frac{1}{2}$	$\bar{x}, \frac{1}{2}, \frac{1}{2}$ $\frac{1}{2}, \bar{x}, \frac{1}{2}$	$\frac{1}{2}, x, \frac{1}{2}$ $x, \frac{1}{2}, \frac{1}{2}$	$\frac{1}{2}, \bar{x}, \frac{1}{2}$ $\bar{x}, \frac{1}{2}, \frac{1}{2}$	$\frac{1}{2}, \frac{1}{2}, x$ $\frac{1}{2}, \frac{1}{2}, \bar{x}$	$\frac{1}{2}, \frac{1}{2}, \bar{x}$ $\frac{1}{2}, \frac{1}{2}, x$
32	<i>f</i> $\dots 3m$	x, x, x x, x, \bar{x}	\bar{x}, \bar{x}, x $\bar{x}, \bar{x}, \bar{x}$	\bar{x}, x, \bar{x} x, \bar{x}, x	x, \bar{x}, \bar{x} \bar{x}, x, x		
24	<i>e</i> $4m \dots m$	$x, 0, 0$	$\bar{x}, 0, 0$	$0, x, 0$	$0, \bar{x}, 0$	$0, 0, x$	$0, 0, \bar{x}$
24	<i>d</i> $m \dots m m$	$0, \frac{1}{2}, \frac{1}{2}$	$0, \frac{1}{2}, \frac{1}{2}$	$\frac{1}{2}, 0, \frac{1}{2}$	$\frac{1}{2}, 0, \frac{1}{2}$	$\frac{1}{2}, \frac{1}{2}, 0$	$\frac{1}{2}, \frac{1}{2}, 0$
8	<i>c</i> $\bar{4}3m$	$\frac{1}{2}, \frac{1}{2}, \frac{1}{2}$	$\frac{1}{2}, \frac{1}{2}, \frac{1}{2}$				
4	<i>b</i> $m \bar{3}m$	$\frac{1}{2}, \frac{1}{2}, \frac{1}{2}$					
4	<i>a</i> $m \bar{3}m$	$0, 0, 0$					

3. A monoclinic crystal has $a = 6$, $b = 4$, $c = 10 \text{ \AA}$, $\beta = 110^\circ$. Make a rough drawing of a portion of the $k = 0$ level of the reciprocal lattice. Show and name the reciprocal lattice vectors and angle for the reciprocal lattice unit cell.

4. In a powder diffractometer, give the function (tell what it does) of the:

divergence slit

receiving slit

Soller slits

monochromator

5. Index this cubic powder diffraction pattern on the next page.

Give the Bravais lattice.

Calculate the lattice parameter for this material. Show your calculations.

$\lambda = 1.54184 \text{ \AA}$.

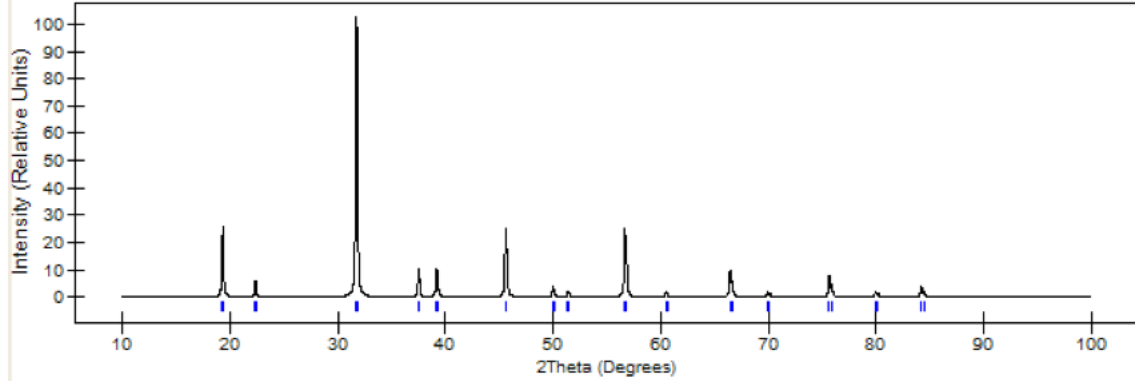
91.214, -7.940

Approximate Digitized Diffraction Pattern

PDF# 00-016-0547. Lead Magnesium Strontium Tellurate. $Pb_2MgSrTeO_6$

00-016-0547 (Approximate) 100%

Peak Markers



2θ	d (Å)
19.279	4.6000
22.319	3.9800
31.703	2.8200
37.473	2.3980
39.169	2.2980
45.569	1.9890
49.960	1.8240
51.345	1.7780
56.667	1.6230
60.457	1.5300
66.440	1.4060
69.937	1.3440
75.654	1.2560
79.947	1.1990
84.195	1.1490