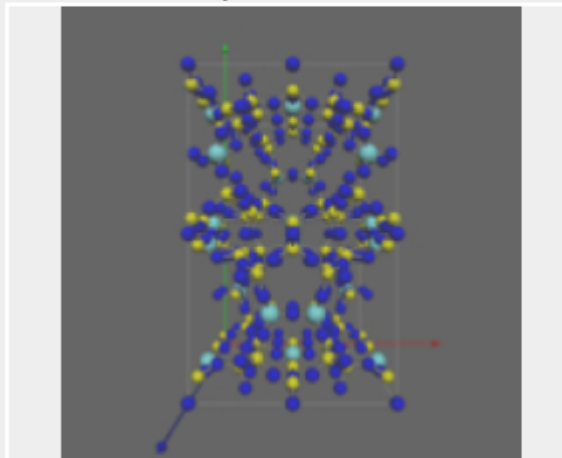


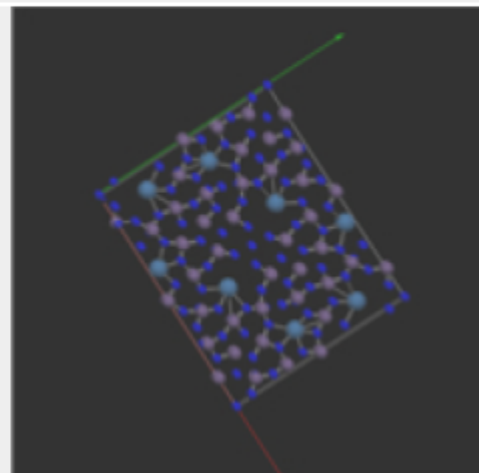


Cubic Hexagonal Trigonal Tetragonal Orthorhombic Monoclinic Triclinic Other Mineral Si3N4 HTC Super-cell

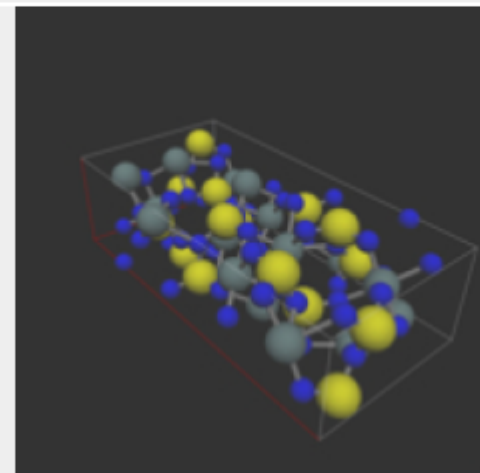
Orthorhombic crystals



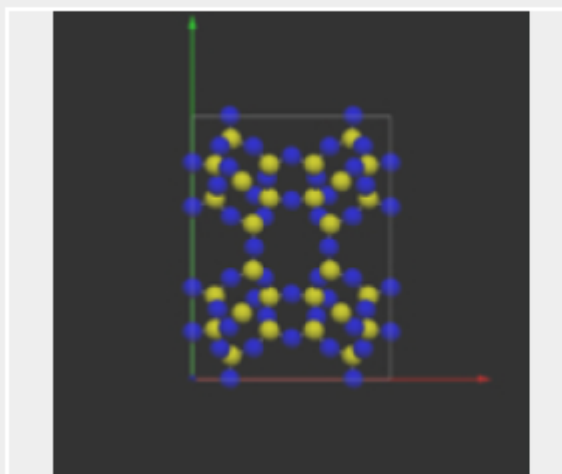
Al₅Si₁₀₇O₂₂₄ - C 2/m 2/m 2/e



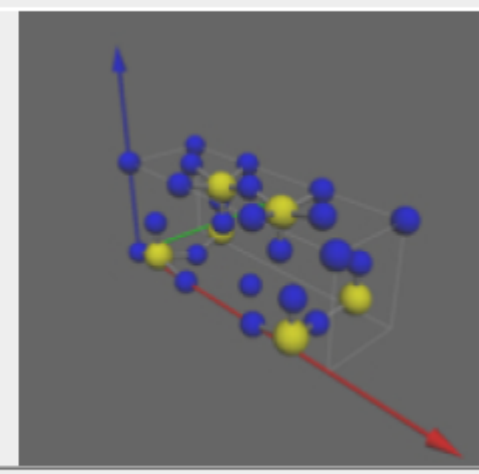
CsNbWO₁₄ - P 21/b 21/a 2/m



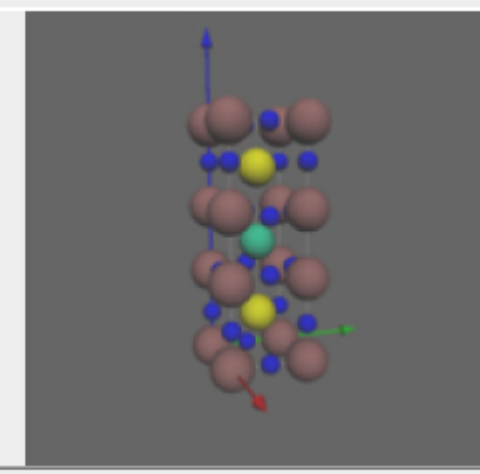
MgSiO₄ - P 21/b 21/c 21/a



Si₇₂O₁₄₄ - F 2/m 2/m 2/m



V₂O₅ - P 21/m 21/m 2/n



YBa₂Cu₃O₇ - P 2/m 2/m 2/m

- Help on JEMS
 - View a crystal structure
 - Diffraction pattern
 - Transfer function
 - Multislice simulation
-
- Import CIF files

File Crystal **Drawing** Imaging Indexing Measuring Parameters Window Help

Diffraction

Ctrl-D

Image 3-D

Ctrl-I

Perspective

Ctrl-P

Reciprocal space

Ctrl-R

Stereogram

Ctrl-S

Thon diagram

Ctrl-H

Transfer function

Ctrl-T

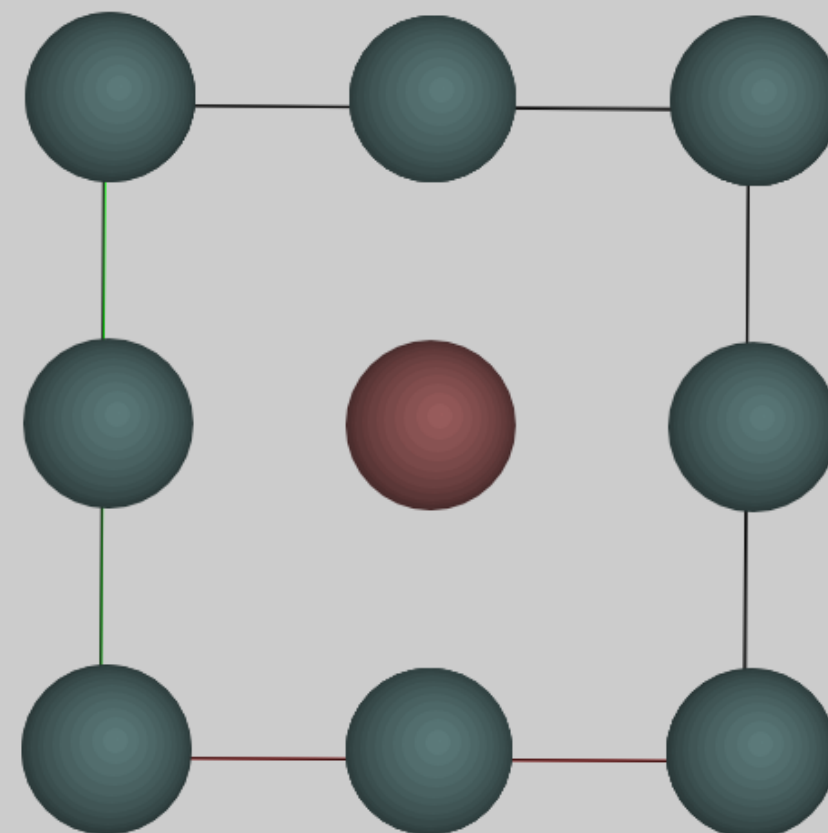
X-Ray powder lines

Ctrl-X



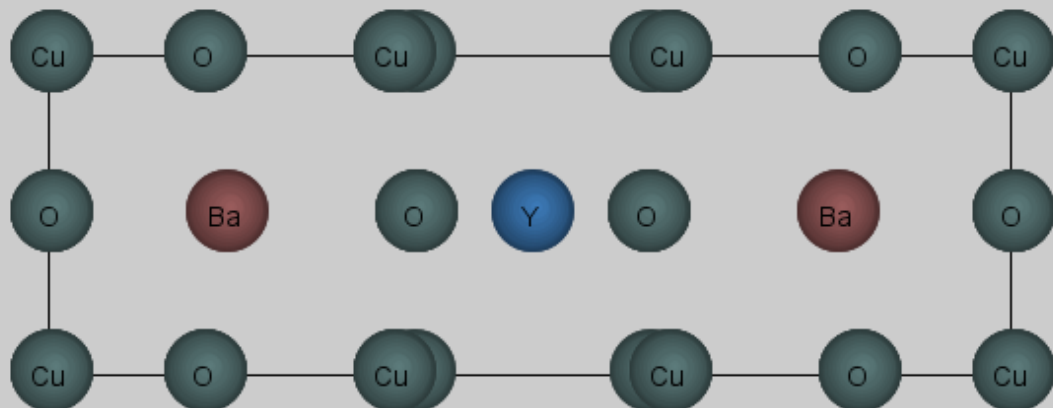
#	Atom	Perspective	Ctrl-P		z	D-W	Oc
0	Y	Reciprocal space	Ctrl-R	0	0.500000	0.0050	1.00
1	Ba	Stereogram	Ctrl-S	0	0.183000	0.0050	1.00
2	Ba	Thon diagram	Ctrl-H	0	0.817000	0.0050	1.00
3	Cu	Transfer function	Ctrl-T	0	0.000000	0.0050	1.00
4	Cu	X-Ray powder lines	Ctrl-X	0	0.355600	0.0050	1.00
5	Cu			0	0.644400	0.0050	1.00
6	O	e	0.000000	0.500000	0.000000	0.0050	1.00
7	O	s	0.500000	0.000000	0.377900	0.0050	1.00
8	O	s	0.500000	0.000000	0.622100	0.0050	1.00
9	O	r	0.000000	0.500000	0.379000	0.0050	1.00
10	O	r	0.000000	0.500000	0.621000	0.0050	1.00
11	O	r	0.000000	0.000000	0.159000	0.0050	1.00
12	O	r	0.000000	0.000000	0.841000	0.0050	1.00

Cu(0)::(x,y,z)=(0.0000,0.0000,0.0000)



YBa2Cu3O7 : [0, 0, 1]

$\text{Cu}(0)::(x,y,z)=(0.0000,0.0000,0.0000)$



Print

Save

Export to Mathematica

☒ Label atoms

☒ Show cell frame

Show powder pattern

Show projected potential image

Show HAADF image

Show SAED pattern

Show WPOA image

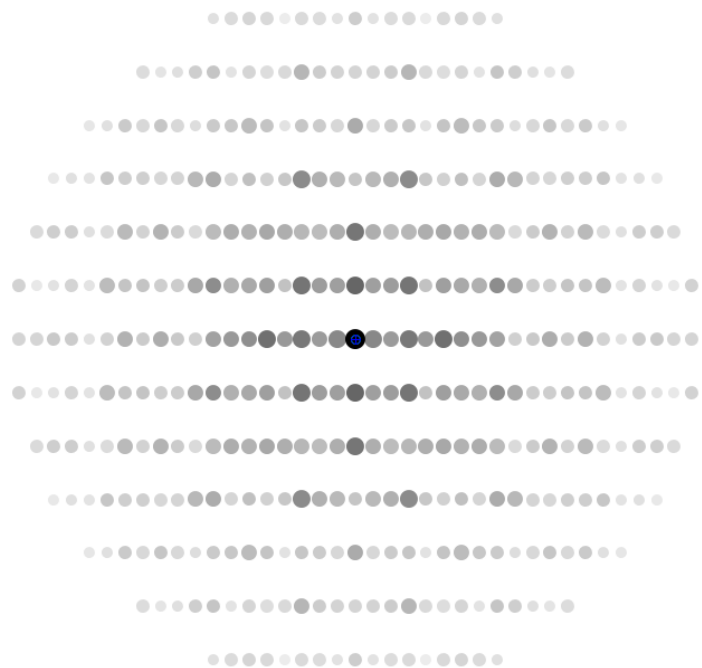
Transfer to clipboard

[UVW] orientation

YBa₂

20 nm⁻¹

Center of Laue circle : (0.0000, 0.0000, 0.0000)::Zone axis : [1, 0, 0]
Tilt angle / deg. : 0.00



YBa2Cu3O7 (407 reflections), thickness:10.0 [nm]
AV/kV:300.000, CL/mm:3292, ZA:[1, 0, 0], FN:[1, 0, 0]

Avalanche Crystal / matrix Diffraction Lattice Laue zones Options Orientation

Options Thresholds

☐ Background
 ☐ Color
 ☒ Dots
 ☐ Double diff.

☐ Fast
 ☐ HOLZ
 ☐ HOLZ shift
 ☐ Indices

☐ Kikuchi
 ☐ Laue circles
 ☐ LZ colors
 ☒ Spots

☐ ZOLZ def-lin.

Atomic Form Factors

☐ DTSB
 ☐ EJK
 ☐ PRDW
 ☐ WK
 ☐ WKc.
 ☒ XRay
 ☐ JEMS