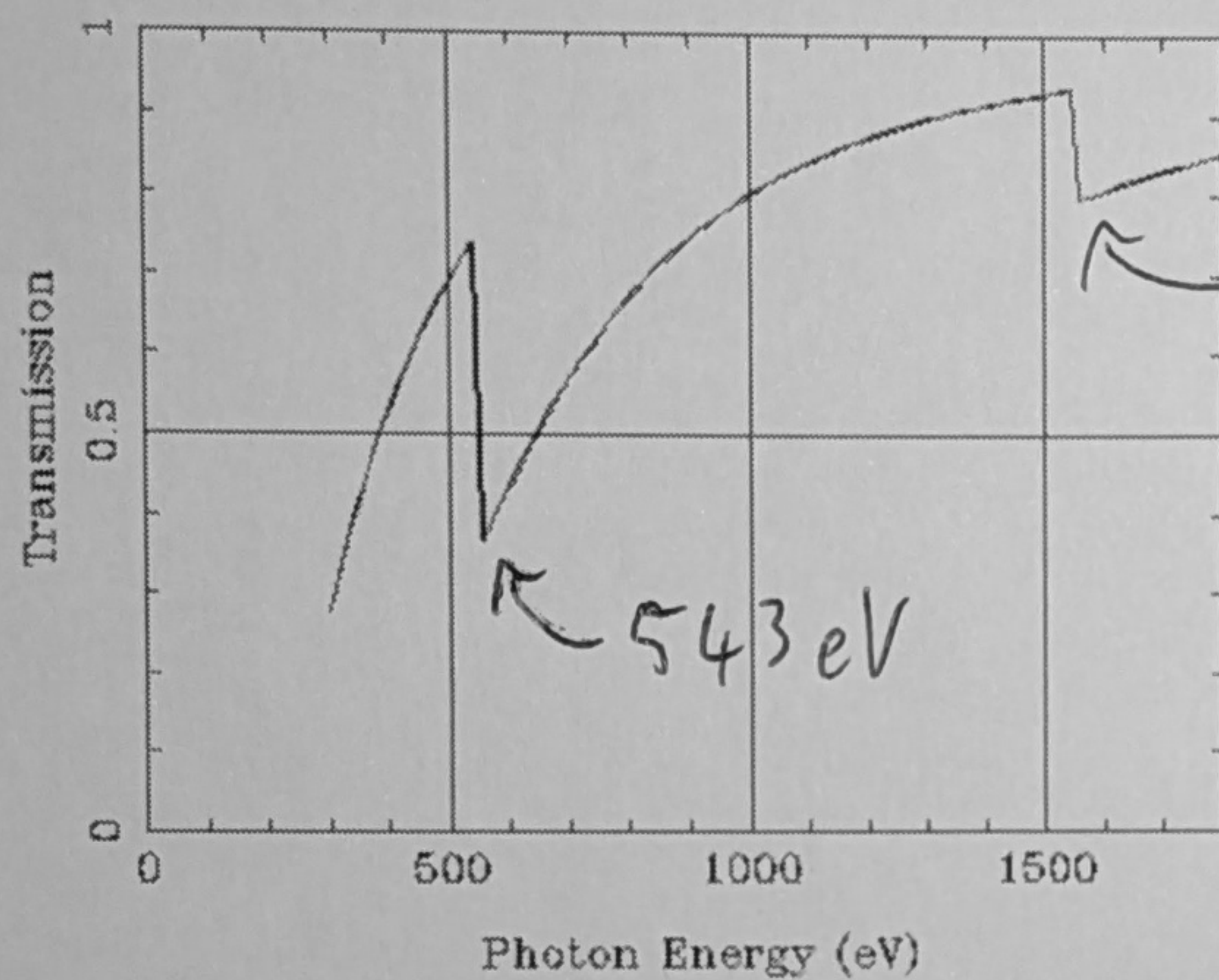


CH417 – X-ray exercises

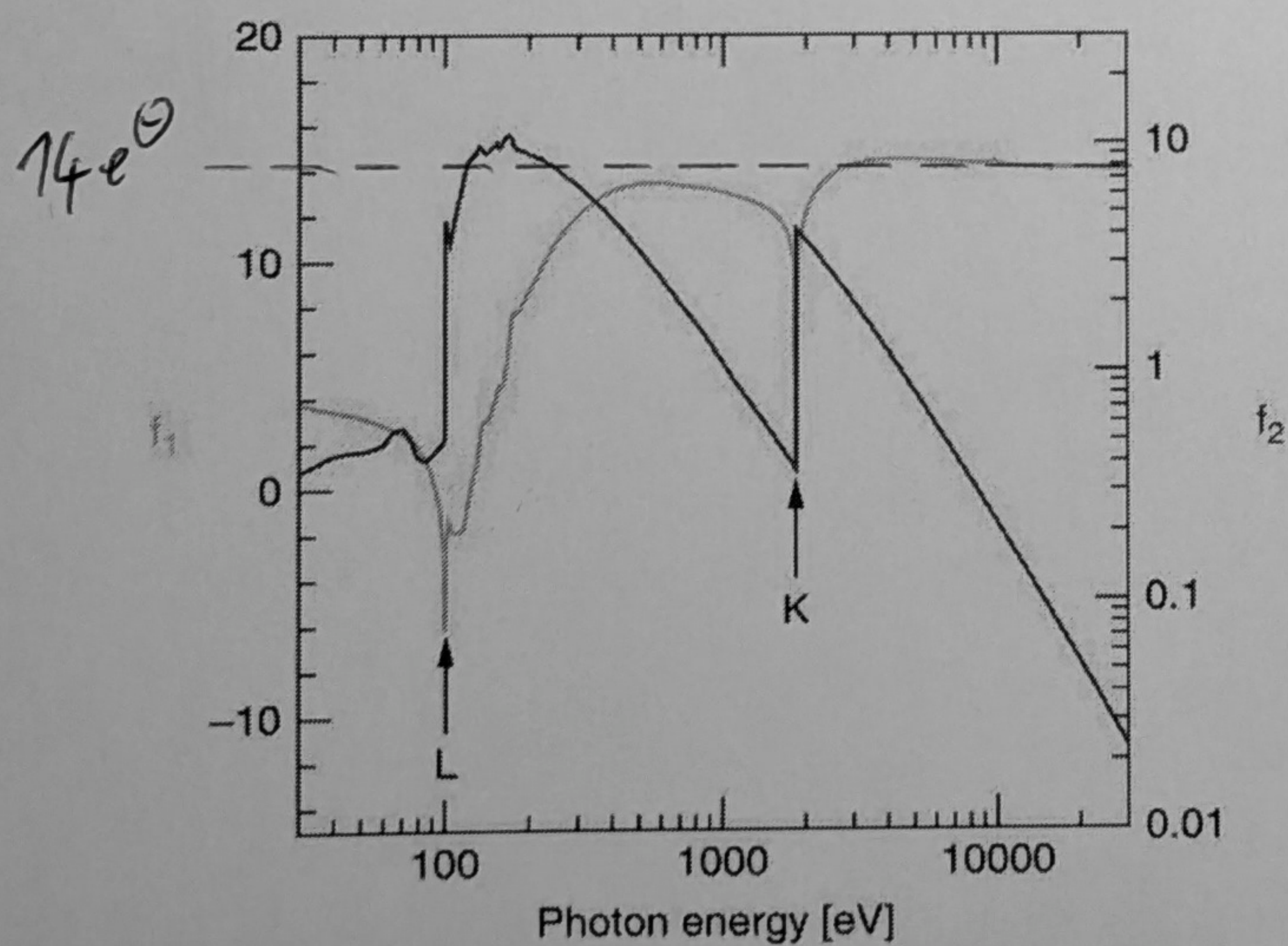
The Phoenix beamline at the Swiss Light Source can record absorption scans in the soft X-ray regime. Figure out the elements in the following sample. Use the electron binding energy tables from the x-ray data booklet on moodle or table on slide 18. Can you guess the material?



543 eV: Oxygen
1560 eV: Aluminium

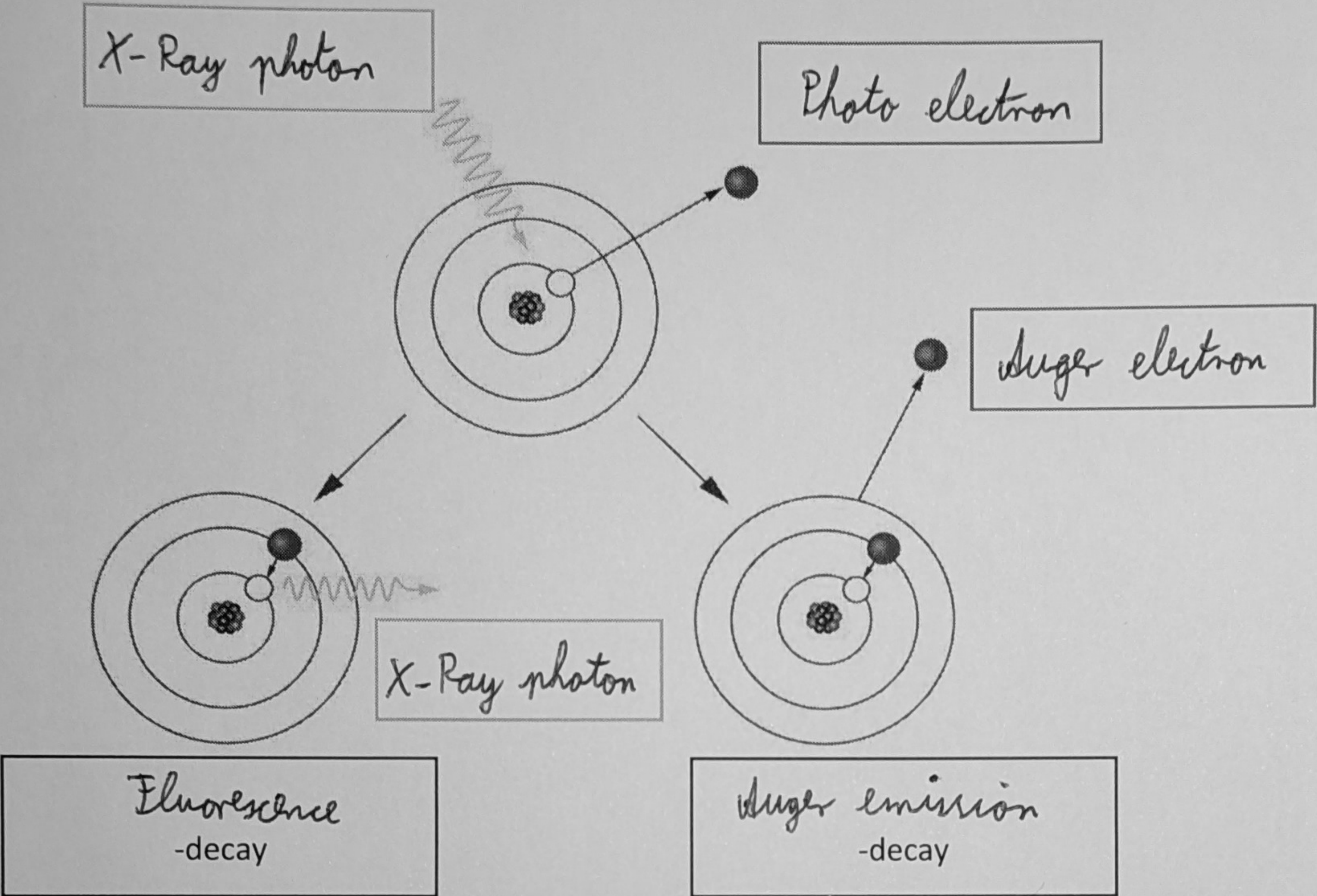
→ Al_2O_3

In the next figure the atomic scattering factors f_1 and f_2 of a single and simple element are plotted. How many electrons does each atom contain? What is the material?

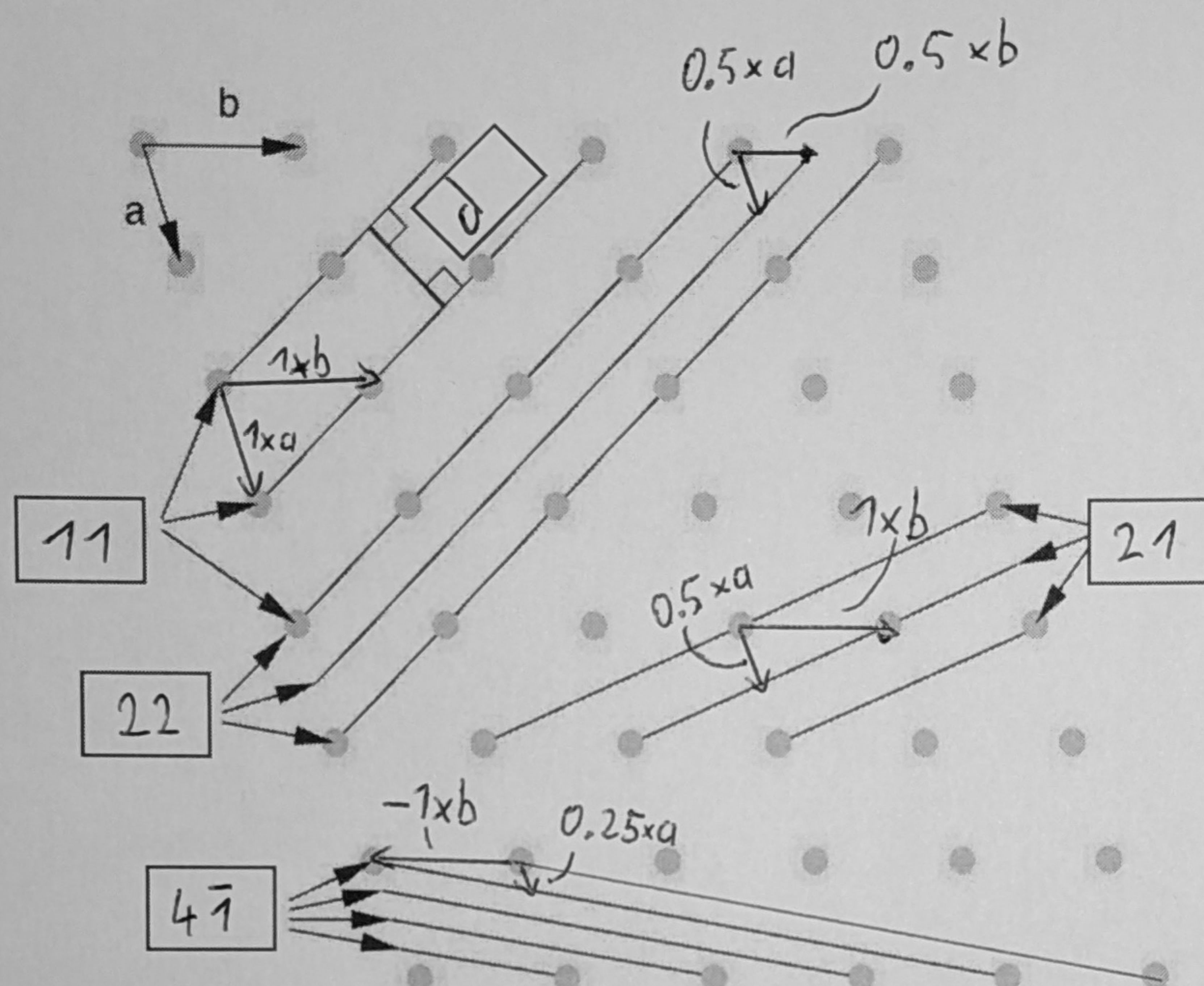


14 electrons: Silicon

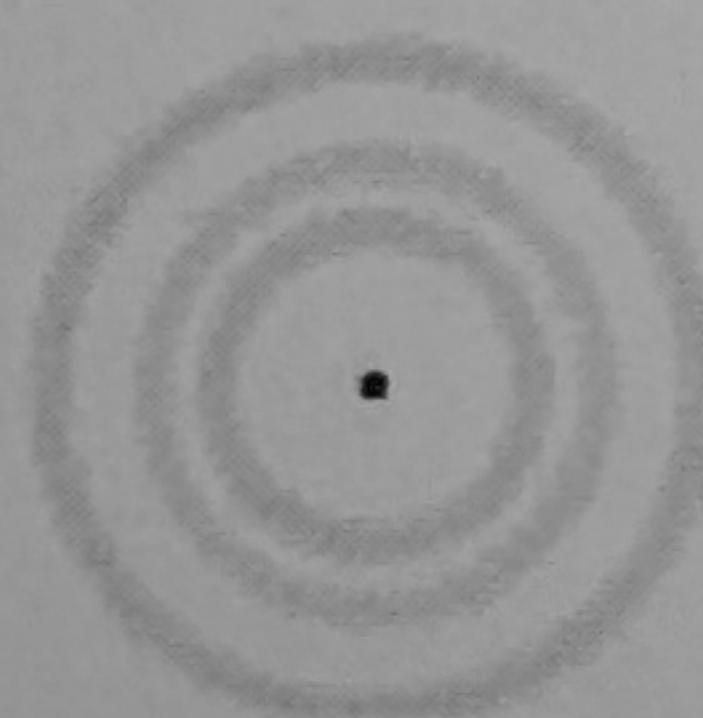
Label the following figure



Identify the lattice planes



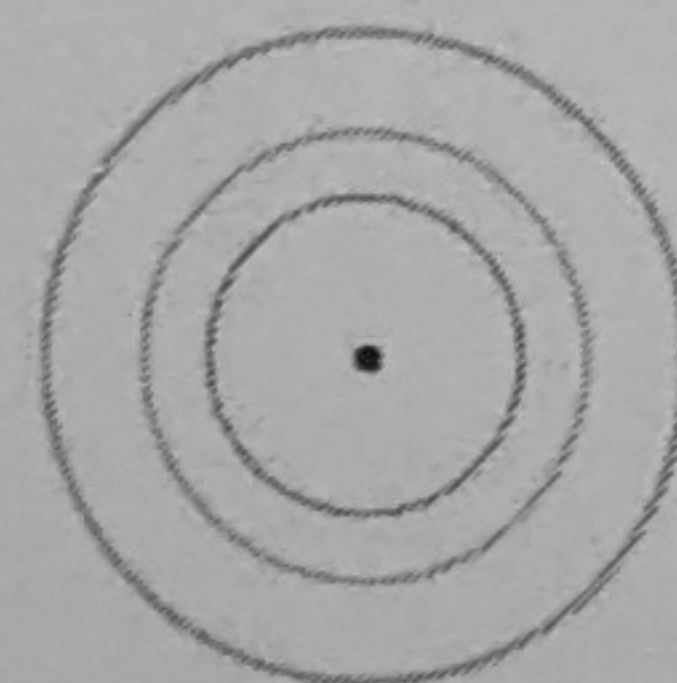
Which diffraction pattern do you attribute to a) single crystal b) powder c) nanocrystalline powder



c)



a)



b)

CH314 – Structural Analysis

Part III : X-ray Tools

Exercise 3, Dec 7

1) Lets pick up the problem from last week again. The data below shows diffraction patterns from a solution of colloidal nanoparticles taken with an x-ray tube using the Cu K α line with a wavelength of 1.54 Angstrom

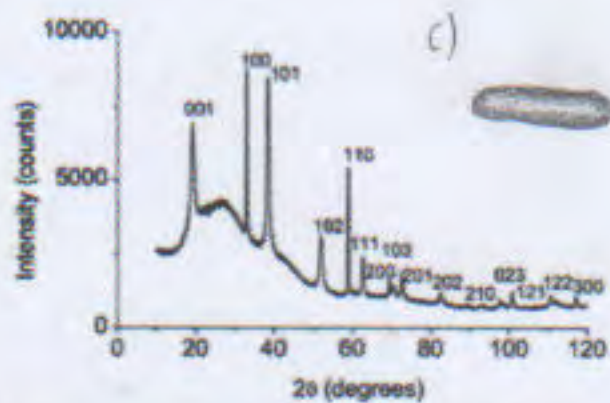


Fig. 1. X-ray diffraction pattern from Ni(OH) $_2$ dispersions ($a = 3.1273(2)$ Å, $c = 4.610(3)$ Å) with diffraction peaks labeled with hkl indices. The 001 peak is

Table 1
Results from the pseudo-Voigt function fit to the Ni(OH) $_2$ sample. 2θ is the peak position for the Cu K α radiation wavelength (0.154056 nm).

hkl	Bragg angle 2θ (°)	Peak width β (°)	Instrument resolution (°)	Corrected β (°)	Diameter (nm)
001	19.25	1.140	0.121	1.02	
100	33.08	0.197	0.115	0.08	
101	38.52	0.593	0.114	0.48	
102	52.06	0.876	0.117	0.76	
110	58.05	0.348	0.120	0.13	
111	62.67	0.550	0.122	0.43	

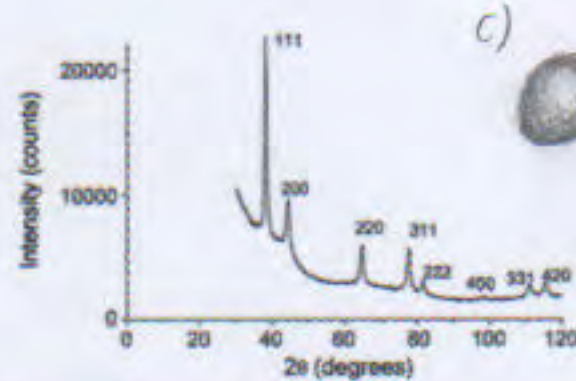


Fig. 2. X-ray diffraction patterns from dispersed Au particles ($a = 4.067(3)$ Å) with diffraction peaks are labeled with hkl indices.

Table 2
Results from the pseudo-Voigt fit to the Au particles. 2θ is the peak position for the Cu K α radiation wavelength (0.154056 nm).

hkl	Bragg angle 2θ (°)	Peak width β (°)	Instrument resolution (°)	Corrected β (°)	Diameter (nm)
111	38.21	0.580	0.114	0.47	
200	44.34	0.822	0.115	0.71	
220	64.69	0.810	0.123	0.68	
311	77.63	0.905	0.135	0.77	
420	115.34	1.194	0.119	1.07	

- What can you tell about the Ni(OH) $_2$ and the Au nanoparticles by just looking at the data?
- Calculate the nanoparticle dimensions for the 001 and 100 reflections for the Ni(OH) $_2$ and 111, 200 reflections for the Au particles.
- Sketch the dimensionality of each of the particles.
- Which lattice do the particles exhibit? Work on the Au nanoparticles first and guess for the Ni(OH) $_2$ particles.

b)
$$\text{Size} = \frac{k \cdot a}{\text{width} \cdot \cos \theta} \quad (\text{see exercise 2})$$

Ni(OH) $_2$

Gold

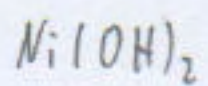
001	7,9 nm	111	17,9 nm
100	103,5 nm	200	12,7
101	17,5	220	13,8
102	17,6	311	13,2
110	70,2	420	13,9
111	27,6		

4) For Gold the peaks have a somewhat similar width $\rightarrow$ almost spherical particle

For the Ni(OH) $_2$ particles the peaks on perpendicular planes (100 and 001) differ severely $\rightarrow$ cylindrical particle

d) Gold: fcc

Ni(OH) $_2$: sc



$\Rightarrow$ Broader peak
indicates fewer
layers

Plane	2θ	β	Diameter
001	19	1.02	8 nm
100	33	0.08	104 nm

