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SPRING COLLEGE IN MATERIALS SCIENCE

ON

"METALLIC MATERIALS"

(11 May - 19 June 1987)

AUGER MICROSCOPY AND SPECTROSCOPY,

PHOTOELECTRON SPECTROSCOPY

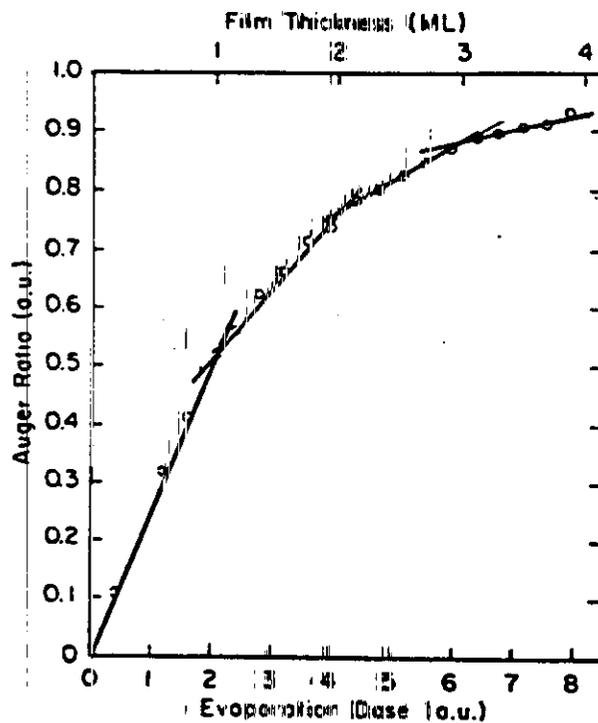
(Lecture III)

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These are preliminary lecture notes, intended only for distribution to participants.

Growth of Co overlayer on Cu(001)



$$\text{Auger Ratio} = \frac{C_{\text{Co}}(656) + C_{\text{Co}}(716)}{C_{\text{Co}}(656) + C_{\text{Co}}(716) + C_{\text{Cu}}(845) + C_{\text{Cu}}(920)}$$

Pescia et al. P.R.L. 58, 933, 1987

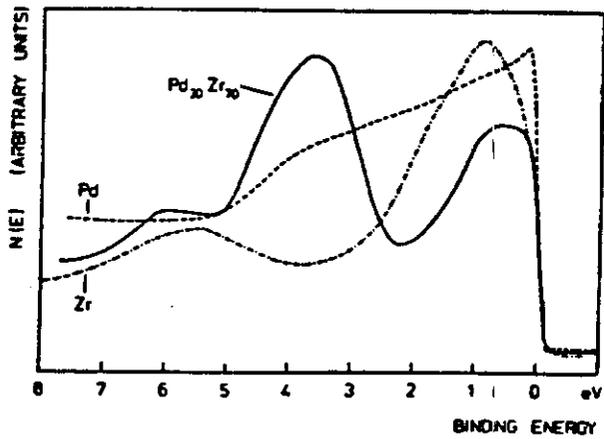


FIG. 1. Uncorrected (without background subtraction) UPS ($h\nu = 40.8$ eV) spectra of the metallic glass $\text{Pd}_{30}\text{Zr}_{70}$, polycrystalline Pd and Zr.

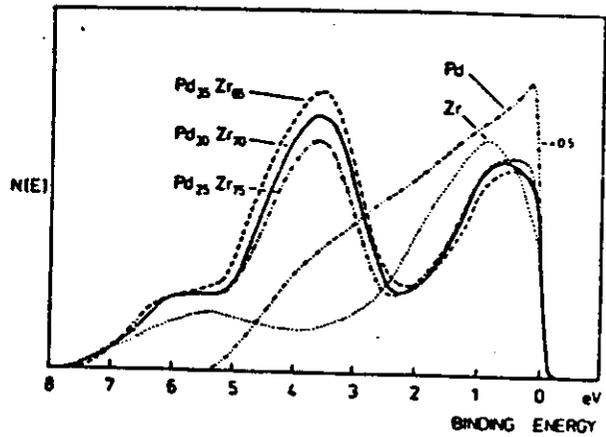


FIG. 2. UPS ($h\nu = 40.8$ eV) spectra of three glassy Pd-Zr alloy compositions, polycrystalline Pd and Zr. The spectra are normalized (see text).

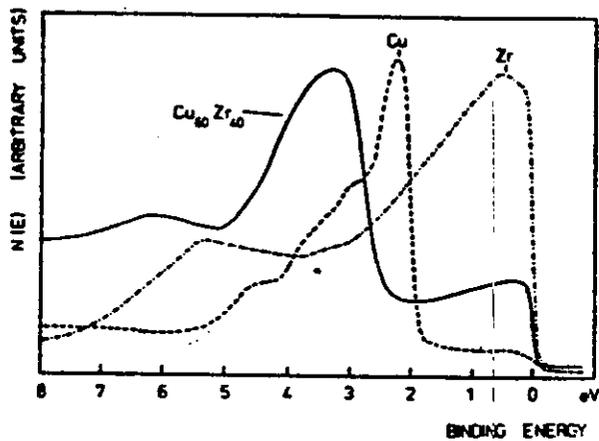


FIG. 3. Uncorrected (without background subtraction) UPS ($h\nu = 21.2$ eV) spectra of glassy $\text{Cu}_{80}\text{Zr}_{20}$, polycrystalline Cu and Zr.

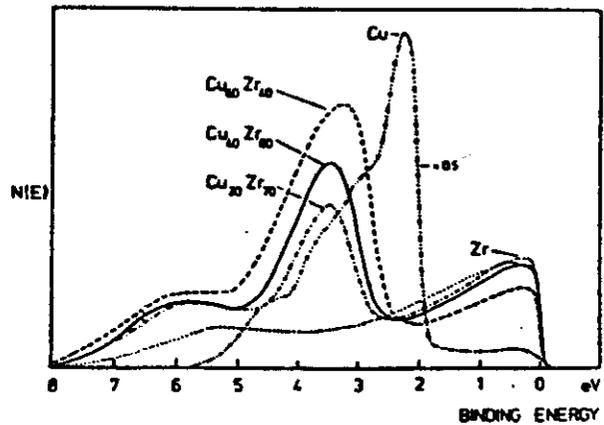


FIG. 4. UPS ($h\nu = 21.2$ eV) spectra of three Cu-Zr samples, polycrystalline Cu and Zr. The spectra are normalized (see text).

Gelhofen et al P.R.L. 43, 1134, 1979.

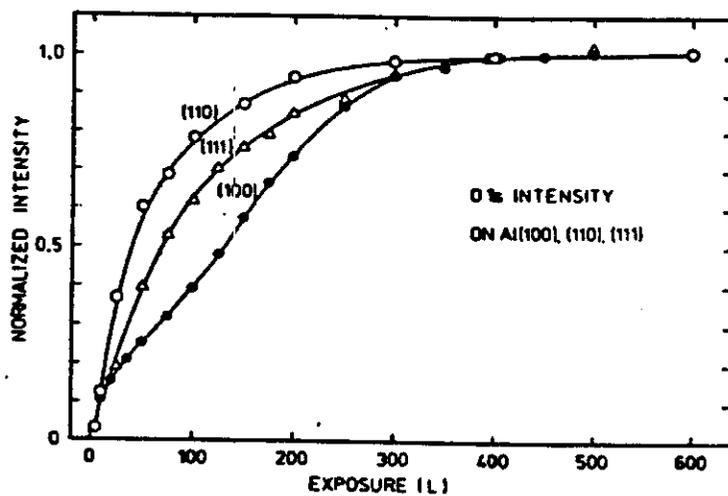


Fig. 2. Intensity change for the O_{1s} peak on the (100), (110) and (111) crystal faces as a function of oxygen exposure.

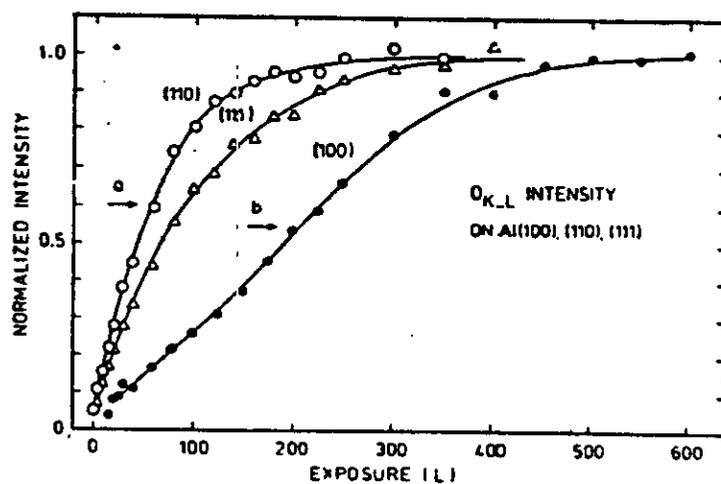
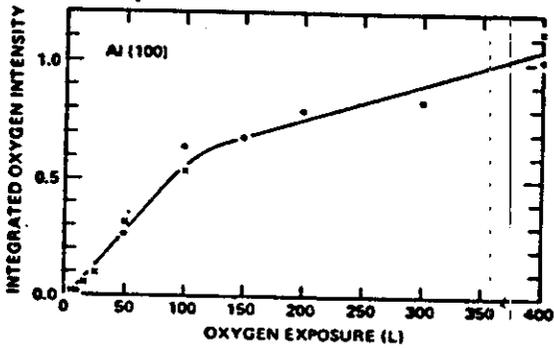
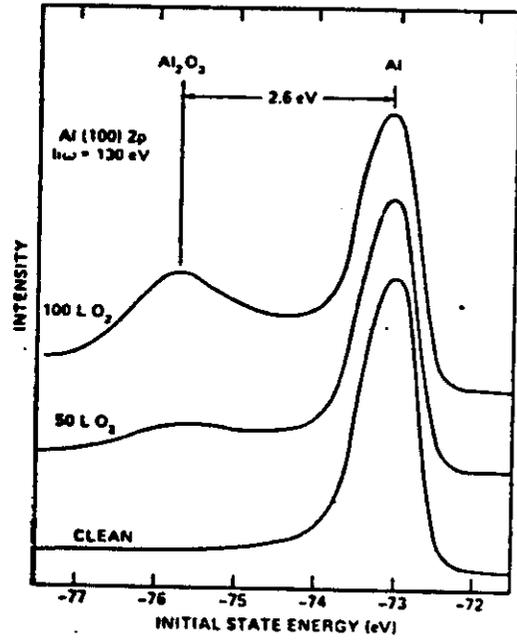


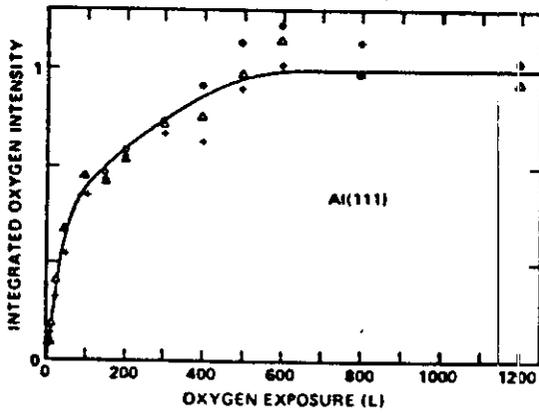
Fig. 1. Intensity change for the O_{KLL} peak on the (100), (110) and (111) crystal faces as a function of oxygen exposure. The disappearance of the LEED pattern from the (110) and (100) faces is indicated by arrows, a and b respectively.



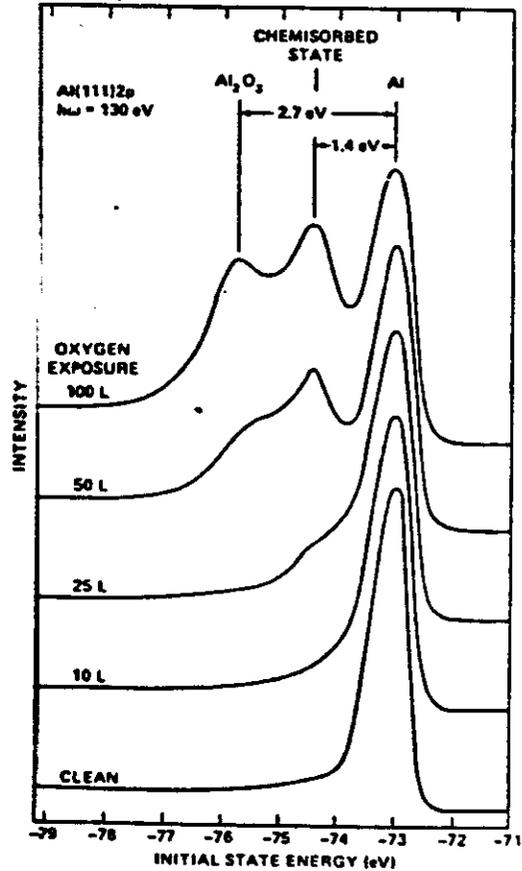
Integrated $2p$ O_2 intensity
Versus exposure for Al(100) surface.

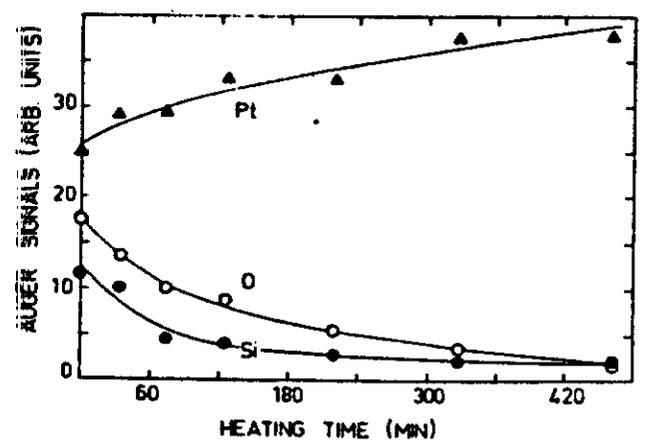
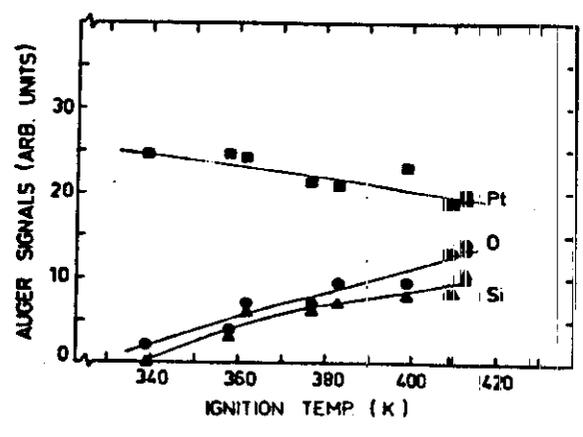
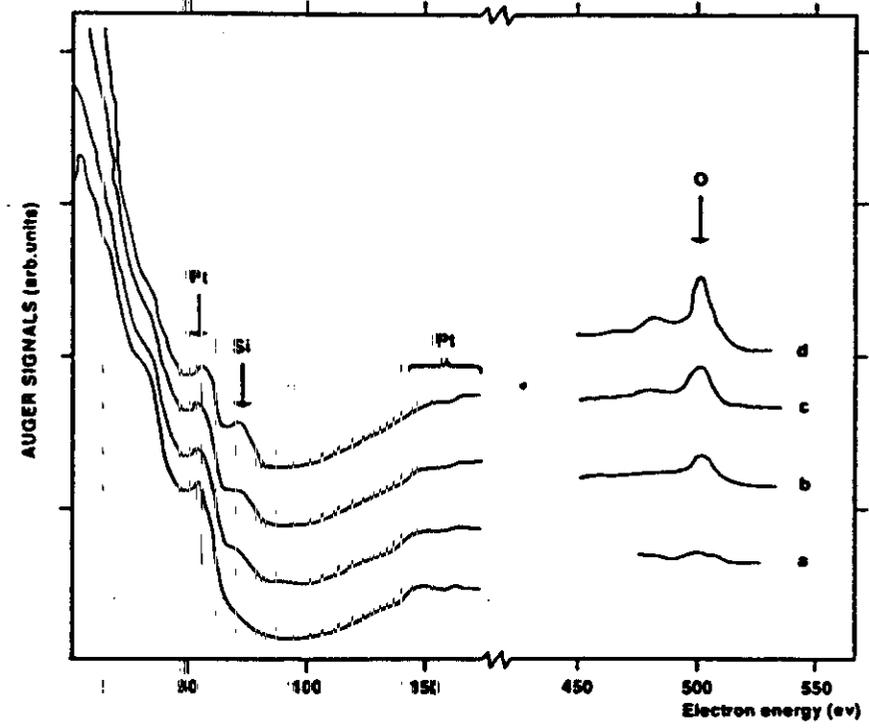
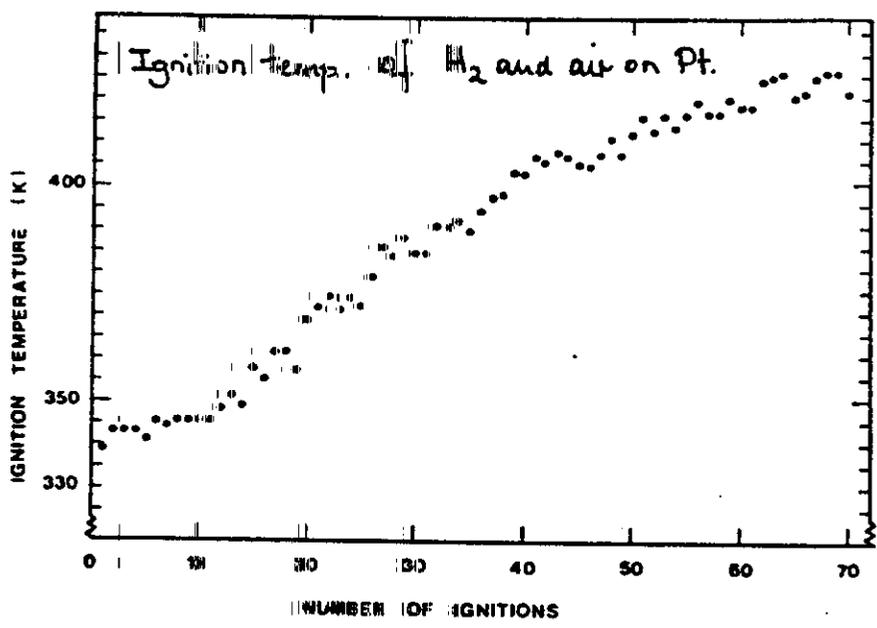


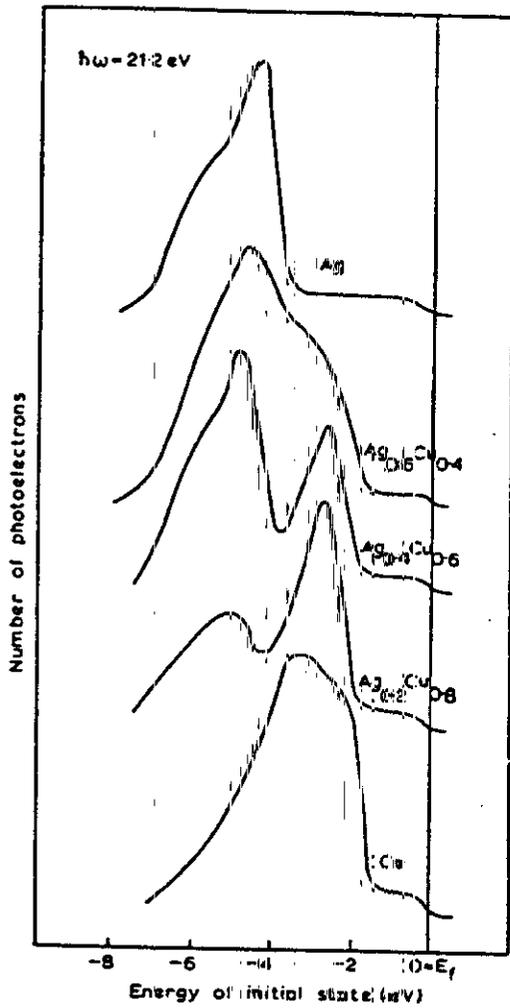
Al $2p$ spectra.



Al(111) surface

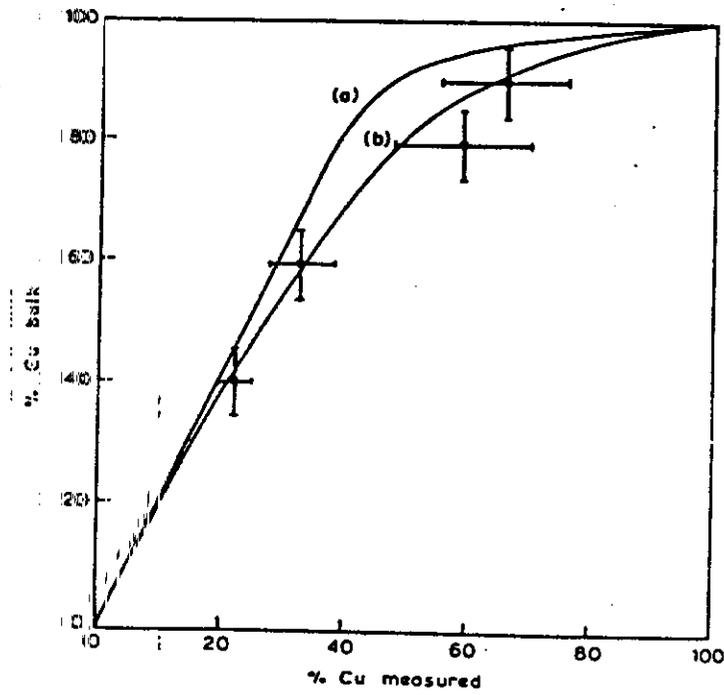






Surface Enrichment
in
Ag-Cu liquid

Williams & Norris Phil. Mag.



Surface enrichment in Ag-Cu alloys. The experimental data are compared with a thermodynamics calculation using (a) no surface enthalpy relaxation, (b) surface enthalpy relaxation parameter $\delta = 0.2$ (see text).

