## Papers, etc.

The two Al 2p peaks characteristic of metallic Al, peaks and -b, were separated by 0.18 eV. To obtain information about the origin of these peaks, Al 2p spectrs of the clean surface at different electron emission singles were masured. Peak-b at normal emission was stronger than at 30 degree take-off single. Therefore, the origin of poak-b is located deeper below the surface and peaks is stributed to a top surface component. Such chemical shifts between top surface and belk atoms were observed also in NAM [100], where the separation of the two peaks was 0.13 eV.95 fixor peak-adisappeared at 512 L, the surface was completely control of the two peaks was 0.13 eV.95 fixor peak-adisappeared at 512 L, the surface was completely control of the state of the surface was completely control of the state of the surface was completely control of the state of the surface was completely control to the state of the surface was completely control to the state of the surface was completely control to the state of the surface was completely control to the state of the surface was completely control to the surface was contr

To obtain information about the origin of peak- (27.63 eV) and -d (74.90 eV), Al 29 poocts were measured at normal emission and 30 degrees take-off single fart 4000 L. exposure, Fig. 3. Peak-d is more pronounced than peak-fig 30 degrees take-off single. Therefore, peak-d is located most towards the surface after outlation. Since peak-b arises from the bulk, it is considered that peak-c originates at the interface between the substrate and outer layer (peak-d). The peak-d at 74.50 eV is attributed to Al in AbOs from the following two considerations: AlOs is the only oxide of aluminum in the phase diagram, and in Al 2p binding energy has been reported as approximately 75 eV, both for bulk outlier and AbOs. Such a chemical shift for a interface component is well-known for The SciOSS interface. <sup>30</sup> An Al 2p chemical shift for a interface component is well-known for the SciOSS interface. <sup>30</sup> An Al 2p chemical shift the to an interface component we well-known for the SciOSS interface. <sup>30</sup> An Al 2p chemical shift the to an interface component was reported for a

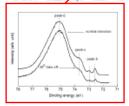


Fig. 3 The Al 2p photoelectron spectra obtained with 106 eV light at normal emission and 30 degrees take off angle after 4000 L exposure.

NiAl(110)<sup>47</sup>, where the interface was considered to be Al-ternizated. It has been reported that in the case of an oxygen terminated interface made by the deposition of pare Cu on alpha-AlOc(0001), no Al 2p peak due to an interface Al component was observed. The Therefore, it is believed that the interface is Al-terminated in

The thickness of the interface layer and alumina layer was estimated using the method described in reference 15. With this method, the thickness of the interface layer (di) is estimated from eq. (1) and that of the alumina layer  $(d_2)$ from eq. (2).



where  $P^{\mu\nu}_{amon}$ ,  $P^{\mu\nu}_{amon}$ , and  $P^{\mu\nu}_{cool}$  are Al. 2p intensity from alumnus (peak-d), the interface (peak-d) and Ca-Al alloy (peak-d). The following values of the inelastic mean free path of the Al. 2p photoelectron are used (kinetic energy of cs. 30 eV); alumnus,  $X^{\mu\nu}_{amon}$ ,  $X^{\mu\nu}_{amon}$ ,  $X^{\mu\nu}_{amon}$ ,  $X^{\mu\nu}_{amon}$ ,  $X^{\mu\nu}_{amon}$ ,  $X^{\mu\nu}_{amon}$ ,  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are atomic densities of Al in the Cu-Al alloy, in the interface layer at Al (60.2 x 107° /erm<sup>2</sup>) and alumnus (4.60 x 107° /erm<sup>2</sup>) was used. The values used for  $N^{\mu\nu}_{amon}$ ,  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  are  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  and  $X^{\mu\nu}_{amon}$  and  $X^{\mu$ 

