

Controlling Orientation of phthalocyanine molecules by use of PTCDA ultra-thin template layer

Takeaki SAKURAI¹, Shunsuke KAWAI¹, Jo SHIBATA¹, Ryousuke FUKASAWA¹,
Katsuhiko AKIMOTO¹, Takahiro MARUYAMA², and Hiroo KATO³

¹Institute of Applied Physics, University of Tsukuba, Tsukuba, Ibaraki 305-8573, Japan

²Material Science and Engineering, Meijo University, Tenpaku-ku, Nagoya 468-8502, Japan

³Faculty of Science and Technology, Hirosaki University, Hirosaki 036-8561, Japan

Introduction

Phthalocyanine (H_2Pc) and metal phthalocyanine (MPc) have been widely applied to the optoelectronic devices due to their high carrier mobility and high photoabsorbance in visible region. However, single crystalline phthalocyanine shows the anisotropic electric conductivity because the stacking direction of π orbitals mainly contributes to the carrier transport. Therefore, the development of a technique of controlling molecular orientation is of importance for fabricating high performance devices.

It was recently reported that over 75 Å thick perylene tetracarboxylic dianhydride (PTCDA) template layers controlled the stacking direction of H_2Pc films on quartz [1]. However, definite orientation of H_2Pc grown on various thick PTCDA layer, particularly in the initial stage of growth, have not been fully understood so far. In the present study, we applied angle resolved ultraviolet photoemission spectroscopy (ARUPS) to determine orientation of H_2Pc on PTCDA template layers.

Experiment

Highly oriented pyrolytic graphite (HOPG) was cleaved in air and cleaned by annealing at 700 °C in an ultra high vacuum (UHV) chamber. H_2Pc and PTCDA were purified by temperature-gradient sublimation method. Thin film samples were prepared by vacuum deposition in a UHV chamber. The deposition was performed at a rate of 1-5 Å/min with the substrate held at room temperature. ARUPS was carried out at the station BL-11C of photon factory in the institute of Material Structure Research.

Results and Discussions

Figure 1 shows the ARUPS spectra of H_2Pc on PTCDA template layer with the incidence angle (α) = 45 ° and $h\nu = 22$ eV. The intensity was normalized by the photon flux. With increasing in the detection angle (θ), the π band labeled A in Fig.1 increased up to near 55 °, and then decreased, while peak position was almost independent of θ .

Figure 2 shows the angular distribution of the peak A intensity. The solid curve and the open square indicate the thickness of PTCDA template layer of 7.5 Å and 30 Å, respectively. The observed angular distribution is very clear and simple giving a maximum intensity of about 55 ° for both samples. In order to analyze the molecular orientation from the ARUPS spectrum, the result of the

independent atomic center (IAC) approximation in which molecules are assumed to stay parallel to the substrate surface is shown by solid curve [2]. The observed polar angles of the maximum intensity agree well with the calculated ones. It is thus concluded that H_2Pc molecules are arranged on PTCDA with their molecular plane parallel to the substrate surface.

References

- [1] S. Heutz et al., Appl. Phys. Lett. 77, 3928 (2000).
[2] T. Kawaguchi et al., J. Appl. Phys. 75, 1486 (1994).

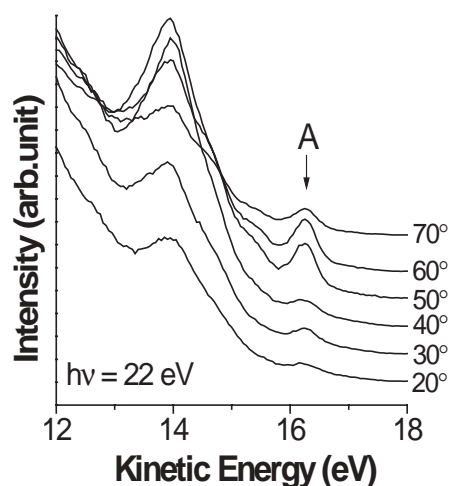


Fig.1 ARUPS spectra of the 90 Å H_2Pc film grown on the 7.5 Å PTCDA template layer.

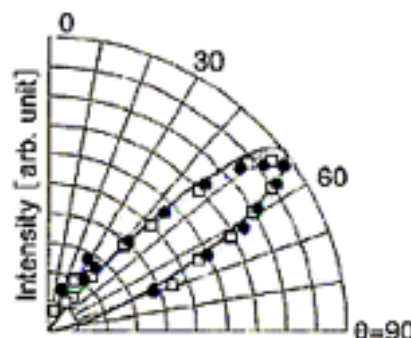


Fig.2 Angular distribution of the peak A intensity. The solid circle and open square indicate the thickness of PTCDA template layer of 7.5 Å and 30 Å, respectively. The thickness of H_2Pc films is 90 Å for both samples. The solid curve shows the result of IAC approximation [2].