27p-PA1-5

Stability of Mesogenic - Phthalocyanine - Based Bulk Heterojunction Solar Cell 阪大院工¹, 産総研ユビキタス² [°]Dao Quang Duy¹, 堀 哲郎¹, 増田 哲也¹, 福村 薫¹, 上門敏也¹, 藤井 彰彦¹, 清水 洋², 尾崎 雅則¹

Osaka Univ.¹, AIST², [°]Dao Quang Duy¹, Tetsuro Hori¹, Tetsuya Masuda¹, Kaoru Fukumura¹, Toshiya Kamikado¹, Akihiko Fujii¹, Yo Shimizu², and Masanori Ozaki¹

E-mail: duy@opal.eei.eng.osaka-u.ac.jp

<u>はじめに</u>: A mesogenic phthalocyanine derivative, 1,4,8,11,15,18,22,25-octahexylphthalocyanine (C6PcH₂) has been demonstrated as a promising small molecule for use in bulk heterojunction (BHJ) solar cells.^[1] C6PcH₂ exhibits not only excellent processability for thin films but also appropriate electronic characteristics for a solar cell, such as a deep highest occupied molecular orbital energy level, a relatively small band gap, strong optical absorption, and high hole and electron drift mobilities exceeding 1.4 $cm^2V^{-1}s^{-1}$ and 0.5 $cm^2V^{-1}s^{-1}$ in the crystalline phase, respectively. In the simple structure, solar cells, the active which layer of was composed of C6PcH₂ and a fullerene derivative, 1-(3-methoxy-carbonyl)-propyl-1-1-phenyl-(6,6)C61 (PCBM), showed a high performance and high stability.^[2] Herein, we clarify mechanism of stability on mesogenic-phthalocyanine-based BHJ Solar Cell.

<u>実験</u>: MoO_x films were thermally evaporated onto ITO substrates. A solution containing a mixture of C6PcH₂:PCBM (2:1) in toluene or chloroform was spin-cast onto a MoO_x layer. A similar process was performed to fabricate the P3HT-based devices. Finally, a LiF buffer layer and an aluminum layer were deposited through a shadow mask by thermal evaporation.

<u>結果</u>: Fig. 1 shows the cell characteristics of the based devices stored in the air as a function of time, when illuminated by a solar simulator. Obviously, the efficiency is correspondingly fallen to the short – circuit current and decreased to 70% after 12-hour illuminating. It exhibits the high stability of the C6PcH₂ based devices, compared with polymers based ones which decreased to 60% after 9-hour illuminating by solar simulator. ^[2] The C6PcH₂ and PCBM composite thin films have also been characterized by XPS and XRD. It exhibited that the bonds of nitrogen atoms at twofold-coordinated nitrogens: the two pyrrole aza nitrogens, denoted as N2 in Fig. 2, and the four meso-bridging aza nitrogens, denoted as N3 with neighboring carbons were broken, or the C6PcH₂ rings were opened.

謝辞: 本研究は JST 先端的低炭素化技術開発(ALCA)の援助の基に行われた。

[1] Y. Miyake et al., Appl. Phys. Express, 4 (2011) 021604. [2] Q.D. Dao et al., Jpn. J. Appl. Phys., 52 (2013) 012301



Fig. 1: Characteristics, normalized by their initial values, as a function of illumination time,

Fig. 2: Molecular structure of C6PcH₂

of solar cells fabricated utilizing (a) P3HT and (b) C6PcH₂