

# Small-Angle Neutron Scattering Study on the Miscibility and Concentration Fluctuation of Hydrogen Bonding Polymer Blends

Osamu Urakawa, Hiroyoshi Ikuta, Toshiyuki Shikata

*Dept Macromolecular Sci. Osaka Univ.*

The effects of hydrogen bonding (H-bonding) on the composition fluctuation in miscible polymer blends were examined by utilizing small-angle neutron scattering (SANS) method. Specifically, we used the blends composed of poly(vinyl acetate-co-vinyl alcohol) (P(VA-VOH)) with various VOH content ( $f_{OH}$ ) and deuterated poly(ethylene oxide) (dPEO), since the formation of the H-bond between the hydroxyl group of P(VA-VOH) and the ether oxygen of dPEO is expected. Furthermore, the advantageous point of this system is that the degree of H-bonding can be changed by the variation of  $f_{OH}$ . In this study, we prepared P(VA-VOH) samples having different  $f_{OH}$  (molar fraction) ( $= 0, 0.10, 0.18, 0.28, 0.35$ ) by the saponification reaction of PVA. The blend composition of P(VA-VOH):dPEO = 80:20 was mainly examined for the SANS measurements.

For all the blends, the SANS intensities increased with increasing temperature  $T$  implying the existence of lower critical solution temperature. For the P(VA-VOH)/dPEO (80/20) blends with  $f_{OH} < 0.28$ , the scattering profiles  $S(q)$  could be well represented by the random phase approximation (RPA) theory. Based on this theory, the interaction parameter  $\chi$  was determined as a function of  $T$ . Figure 1 shows the  $f_{OH}$  dependence of  $\chi$  at  $T=343K$ . As seen in this figure, the  $\chi$  shows a minimum at around  $f_{OH}=0.15$ , meaning that the blend at this particular  $f_{OH}$  is the most miscible. Furthermore, we found that the random copolymer theory (solid curve in the figure) could quantitatively explain the  $f_{OH}$  dependence of  $\chi$ .

For the blend of  $f_{OH}=0.35$ , on the other hand, the RPA theory did not fit the  $S(q)$  well, especially at the small  $q$  re-

gion. The experimental data at the small  $q$  were found to be more suppressed than the theoretical prediction. This behavior suggests that the formation of larger number of H-bonds might suppress the composition fluctuation.

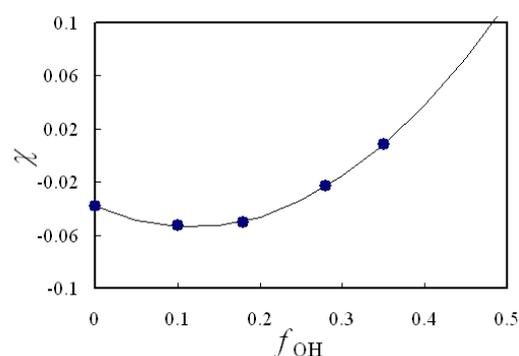


Fig. 1.  $f_{OH}$  dependence of the interaction parameter  $\chi$  for P(VA-VOH)/dPEO (80/20) blends at  $T=343K$ . Solid curve represents the fitted result with the random copolymer theory.