Equilibrium and Kinetic H₂ Separation in Organic Templating Membrane by Molecular Properties

문종호, 정종태¹, <u>이창하</u>* 연세대학교 화학공학과; ¹한국가스공사 (leech@yonsei.ac.kr*)

Hydrogen separation on micro-porous methyltriethoxy- silane templating silica composite/ α -alumina membrane (MTES membrane) was studied using six pure CO $_2$, CO, N $_2$, CH $_4$, He and H $_2$ gases , four binary mixtures – H $_2$ /N $_2$, H $_2$ /CO $_2$, H $_2$ /CH $_4$, H $_2$ /CO and one quaternary mixture – H $_2$ /CH $_4$ /CO/CO $_2$. The equilibirum and kinetic characteristics of an MTES membrane were studied using unsteady and steady-state permeation. The separation mechanisms could be primarily ascribed to the molecular size, and structure of each gas which likely contributed to steric hindrance or molecular sieving within the membrane pore. In addition, the adsorption affinity of each molecule on the membrane surface acted as a key factor in separation performance because it highly contributed to the surface diffusion. The generalized Maxwell–Stefan model incorporating the dust gas model and the Langmuir model could successfully predicted the transient and steady-state permeation/separation.