

Experimental Study and Modeling of CO<sub>2</sub> Solubility in Aqueous Solution of DIPA and MDEA Blended Amine Using Electrolyte-UNQUAC model

최봉근, 김승모, 이종섭<sup>1</sup>, 신현용<sup>2</sup>, 민병무<sup>3</sup>, 문종호<sup>†</sup>

충북대학교; <sup>1</sup>한국에너지기술연구원; <sup>2</sup>서울과학기술대학교; <sup>3</sup>에너지토피아

(moonjongho@chungbuk.ac.kr<sup>†</sup>)

The technology of carbon capture and utilization (CCU) is a critical issue in various industries. Some kinds of amines have been used for CO<sub>2</sub> absorbent: MEA, DIPA, MDEA, and AMP. In this study, equilibrium solubility of CO<sub>2</sub> in DIPA and MDEA mixtures were evaluated by using experimental data and thermodynamic models. The solubility of CO<sub>2</sub> was measured at 50, 110 °C and at various amine mass fractions. To consider the non-ideality, activity coefficient model, electrolyte-UNQUAC model, is used to predicted interactions between chemical species in liquid phase. Calculations of CO<sub>2</sub> solubility and optimizations (parameter regression) were conducted by in-house model in MATLAB® 2020a version. Effect of temperature and blending ratio of DIPA, MDEA, and H<sub>2</sub>O about CO<sub>2</sub> solubility were estimated. Also, mole fraction of chemical species, pH, heat of absorption, cyclic capacity were predicted successfully.