

Improvements in the Measurement and Modeling of CO₂-Brine Properties

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Carbon capture and storage (CCS) and geothermal energy extraction are alternative approaches to fossil fuels exploitation considered to effectively reduce CO₂ emissions in the atmosphere. One possibility is to store CO₂ in depleted petroleum reservoirs or saline aquifers. Another approach is to use CO₂ as a working fluid to enhance heat transfer from depth to surface in geothermal reservoirs. The correct prediction of the phase equilibrium conditions of CO₂ + brine mixtures and the long-term trapping of CO₂ in the saline aquifer require the accurate knowledge of the *P-T-x* properties of these fluids at reservoir conditions.

While the majority of experimental and modeling works have used NaCl as the only salt component in the brine, the presence of other salts can affect the brine properties differently. Therefore, it is important to obtain new datasets for the thermophysical properties of CO₂ + brines containing multiple salts. The objective of this work is to review new improvements for the experimental determination of the volumetric and transport properties of CO₂-brine solutions at conditions of temperature and pressure relevant to carbon storage and geothermal energy production.

The solubility of CO₂ in brine solutions containing NaCl, CaCl₂, MgCl₂, SrCl₂, KCl, NaHCO₃, BaCl₂ and Na₂SO₄ were measured by using an experimental gravimetric method, at 303, 333 and 373 K, from 5 to 40 MPa, and for salinities up to 2 mol·kg⁻¹. A vibrating U-tube densimeter and a capillary viscometer were used to measure the density and viscosity of brines loaded with controlled amounts of CO₂ at the same conditions. The geochemical simulator PHREEQC was used to predict the solubility of CO₂ in multi-component brines. In addition, attempts were made to generalize the Søreide-Whitson model to calculate the solubility of CO₂ in brines containing NaCl, KCl, MgCl₂, CaCl₂ and Na₂SO₄ for temperatures between 273 and 473 K, pressures up to 200 MPa and salinities up to 6 mol·L⁻¹. The results of our study showed that the modified model considerably improves the description of the individual salt brines and calculates the CO₂ solubility in a large variety of brine solutions, including our measurements, with good accuracy when compared with previous parameterizations. The apparent molar volume approach was used for the density modeling.