

Examining How CO₂ Displaces Brine at the Pore Level

Dustin Crandall^{1,2}, Robert P. Warzinski³, and William K. O'Connor⁴

¹*URS Washington Division, Morgantown, West Virginia, USA*

²*National Energy Technology Laboratory, Morgantown, West Virginia, USA*

³*National Energy Technology Laboratory, Pittsburgh, Pennsylvania, USA*

⁴*National Energy Technology Laboratory, Albany, Oregon, USA*

ABSTRACT: Combining micro-computed tomography scanning and computational fluid dynamics this study examined how CO₂ displaces brine in a small volume of pore space. The pore volume ($4.41 \cdot 10^{-11} \text{ m}^3$) was obtained from an Xradia Micro XCT-400 scanner with a voxel resolution of 27.1 microns from a sub-core of Mt Simon sandstone, recovered from a depth of $\approx 1770\text{m}$. The complex micro-structure was preserved via careful three-dimensional reconstruction of the scanned data into a fully meshed volume, with over 3.5 million hexahedral cells. A series of simulations was performed with the commercial CFD code FLUENT to evaluate the immiscible flow behavior of CO₂ and brine within this volume, under a range of conditions that might be expected at a CO₂ storage site. This was done using the volume-of-fluid method to handle multiphase flow, and solving the full Navier-Stokes equations. Initial results show that an increase in the CO₂ viscosity, which could be obtained by adding a surfactant or by increasing depth of the injection, results in an increase in the residual volume of CO₂ within the pore space.