

TIME-LAPSE MODELING AND INVERSION OF CO<sub>2</sub> SATURATION FOR  
SEQUESTRATION AND ENHANCED OIL RECOVERY

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## ABSTRACT

In the second quarter of this DOE NETL project, we have continued Phase I and Phase II activities for researching and developing new technology to quantitatively model the rock physics effects of CO<sub>2</sub>-oil-water systems. These activities included completing a literature search of currently available equation-of-state methods, initiating work in molecular dynamics modeling, and building a prototype seismic modeling code for predicting time-lapse CO<sub>2</sub> changes in well-log models. We have also received permission to use the Sleipner time-lapse CO<sub>2</sub> data set from the North Sea for the current NETL project, and have issued a formal request for the data.

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## I. EXECUTIVE SUMMARY

In the second quarter of this DOE NETL project, we have continued Phase I and Phase II activities for quantitatively modeling the rock physics effects of CO<sub>2</sub> injection (both miscible and immiscible) in porous reservoir/aquifer rock systems containing oil-water phases. These activities have included completing a literature search of currently available equation-of-state methods, initiating work in molecular dynamics modeling, and building a prototype seismic modeling code for predicting time-lapse changes from well-log data during CO<sub>2</sub> injection. We have also received permission to use the Sleipner time-lapse CO<sub>2</sub> data from the North Sea for our project work, and have formally requested the data.

## II. EXPERIMENTAL

No experimental methods were used during this reporting period.

## III. RESULTS AND DISCUSSION

We have completed our literature search for calculating CO<sub>2</sub>-oil and CO<sub>2</sub>-brine properties via equation-of-state (EOS) methods. We also initiated research into applying molecular dynamics modeling for calculating properties (such as bulk modulus and density) of CO<sub>2</sub>-oil-brine mixtures. Molecular dynamics modeling is a numerically intensive modeling approach that is able to generate macroscopic properties of a specific mixture of molecular components at equilibrium. The two-pronged approach of using both equation-of-state methods and molecular dynamics modeling in our fluid modeling effort should result in novel methodologies that improve upon current, state-of-the-art techniques, particularly for problematic zones near phase transitions in supercritical CO<sub>2</sub> solutions.

Other work accomplished during this quarter involved the prototyping and refining of a seismic modeling code for predicting time-lapse changes in well-log data ( $V_p$ ,  $V_s$ , and density) during CO<sub>2</sub> injection (see Figure 1). This code will eventually use our fluid modeling methodology developed in Phase I to generate time-lapse changes in seismic traces resulting from CO<sub>2</sub> saturation and pressure changes.

Lastly, we obtained permission to use the Sleipner CO<sub>2</sub> time-lapse data set for this NETL project, and have already requested the data. This industry-standard data set consists of four vintages of 3D seismic data recorded over a shallow saline aquifer during CO<sub>2</sub> injection, and should prove very useful for testing our inversion methods in Phase III.

## IV. CONCLUSIONS

In the next quarter, we will complete our analysis of the most promising equation-of-state methods found in the literature, and will use this information to work towards a derivation of the best theoretical approach for modeling oil-water-CO<sub>2</sub> mixtures. This effort, coupled with continued work in molecular dynamics modeling, should result in a novel, accurate, and robust

fluid modeling capability that can be used for seismic data modeling in Phase II, as well as for inversion of CO<sub>2</sub> pressure and saturation changes in Phase III.

## V. FIGURES

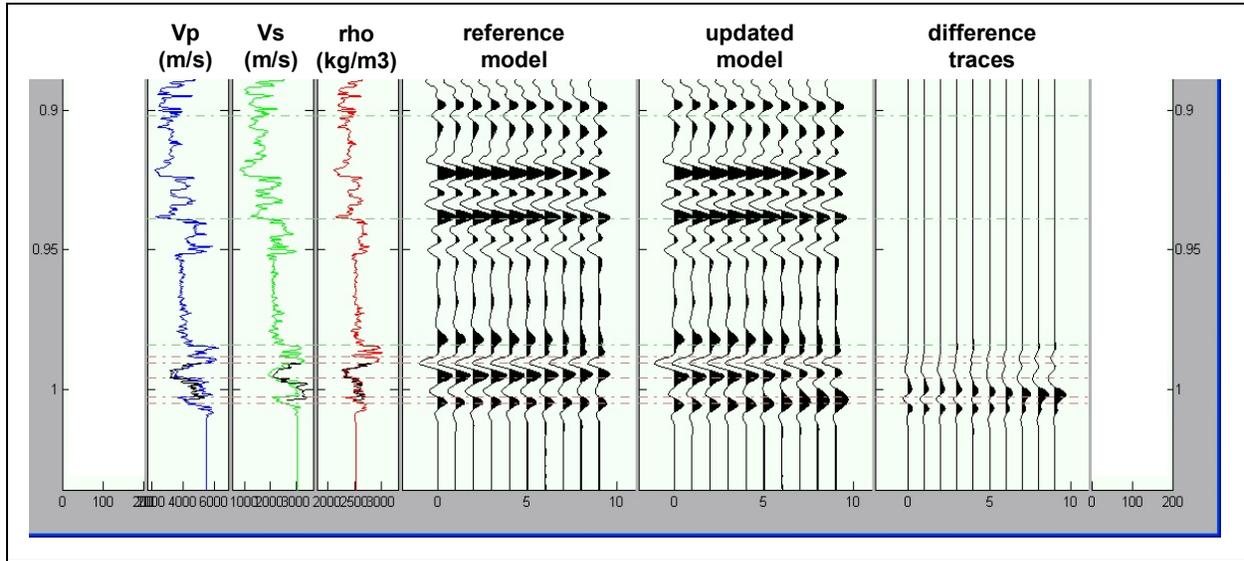


Figure 1 Example of time-lapse seismic modeling of well-log velocities and densities during CO<sub>2</sub> injection.