

## Influence of capillary pressure on CO<sub>2</sub> storage and monitoring

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### SUMMARY

Sequestration of CO<sub>2</sub> in geological formations is one of the solutions to mitigate the greenhouse effect. We are interested in analyzing the influence of capillary pressure on CO<sub>2</sub> injection, storage and monitoring in saline aquifers. To do so, we present a methodology integrating numerical simulation of CO<sub>2</sub>- brine flow and seismic wave propagation. Besides, we build a suitable geological model that includes mudstone layers and fractures. The simultaneous flow of CO<sub>2</sub> and brine in porous media is described by the Black-Oil formulation, which applies a simplified thermodynamic model. Capillary pressure is represented as a potential function of CO<sub>2</sub> saturation. The wave propagation is simulated using a viscoelastic model that includes attenuation and dispersion effects due to mesoscopic scale heterogeneities. The fluid simulator properly models the CO<sub>2</sub> injection, obtaining accumulations below the mudstone layers as injection proceeds. We are able to identify the time-lapse distribution of CO<sub>2</sub> from the synthetic seismograms, which show the typical pushdown effect. When capillary pressure is higher, CO<sub>2</sub> upward migration is slower and thicker zones of CO<sub>2</sub> accumulations are obtained. Numerical examples show the effectiveness of this methodology to detect the spatio-temporal distribution of CO<sub>2</sub> and to make long term predictions.

### INTRODUCTION

Capture and storage of CO<sub>2</sub> in deep saline aquifers is a valid alternative for reducing the amount of greenhouse gases in the atmosphere (Arts et al., 2008). The first industrial CO<sub>2</sub> injection project is the Sleipner gas field, where CO<sub>2</sub> separated from natural gas is being injected in the Utsira formation, a highly permeable porous sandstone (Arts et al., 2008)-(Chadwick et al., 2005) with several mudstone layers which act as barriers to the vertical flow of the CO<sub>2</sub>. Numerical modeling of CO<sub>2</sub> injection and seismic monitoring are important tools to understand the long term behavior after injection and to test the effectiveness of CO<sub>2</sub> sequestration. Recent papers (Carcione and Picotti, 2006)-(Carcione et al., 2012) successfully apply seismic modeling for monitoring the spatio-temporal distribution of CO<sub>2</sub> using assumed saturation maps. Instead, we introduce a methodology to model the gas flow and monitor the storage. We also build a petrophysical model of the Utsira formation based on fractal porosity and clay content, taking into account the variation of properties with pore pressure and saturation (Carcione et al., 2003). The simultaneous flow of brine and CO<sub>2</sub> is modeled with the Black-Oil formulation for two-phase flow in porous media (Aziz and Settari, 1985), which uses the PVT data as a simplified thermodynamic model (Has-

sanzadeh et al., 2008). The wave propagation is based on an isotropic viscoelastic model that considers dispersion and attenuation effects. The complex P-wave and S-wave moduli are determined using the Zener model in the brine saturated mudstone layers (Carcione, 2007); and the White's theory (White et al., 1975) in zones saturated with brine and CO<sub>2</sub>. The results of the flow simulator, allows us to calculate the phase velocities and attenuation coefficients of the P and S waves in order to compute the synthetic seismograms. The methodology is used to model CO<sub>2</sub> injection and flow and compute time-lapse seismograms corresponding to the Utsira aquifer at Sleipner field applying different capillary pressure curves.

### CO<sub>2</sub> - BRINE FLOW MODEL IN POROUS MEDIA

The well-known Black-Oil formulation applied to two-phase, two component fluid flow (Aziz and Settari, 1985) is used to simulate the simultaneous flow of brine (subindex *b*) and CO<sub>2</sub> (subindex *g*), considering that CO<sub>2</sub> may dissolve in the brine but the brine is not allowed to vaporize into the CO<sub>2</sub> phase. This formulation uses the PVT data ( $R_s$ : CO<sub>2</sub> solubility in brine;  $B_g$  and  $B_b$ : CO<sub>2</sub> and brine formation volume factors) as a simplified thermodynamic model, computed with the Hassanzadeh correlations (Hassanzadeh et al., 2008). Combining the mass conservation equations with Darcy's empirical Law, the following nonlinear system of partial differential equations is obtained,

$$\nabla \cdot \left( \underline{\kappa} \left( \frac{\kappa_{rg}}{B_g \eta_g} (\nabla p_g - \rho_g g \nabla D) + \frac{R_s \kappa_{rb}}{B_b \eta_b} (\nabla p_b - \rho_b g \nabla D) \right) \right) + \frac{q_g}{\rho_g^{SC}} = \frac{\partial \left[ \phi \left( \frac{S_g}{B_g} + \frac{R_s S_b}{B_b} \right) \right]}{\partial t}, \quad (1)$$

$$\nabla \cdot \left( \underline{\kappa} \frac{\kappa_{rb}}{B_b \eta_b} (\nabla p_b - \rho_b g \nabla D) \right) + \frac{q_b}{\rho_b^{SC}} = \frac{\partial \left[ \phi \frac{S_b}{B_b} \right]}{\partial t}. \quad (2)$$

Here  $\underline{\kappa}$  is the absolute permeability tensor,  $p_g, p_b$  are the fluid pressures and for  $\beta = g, b$ ,  $k_{r\beta}$  and  $\eta_\beta$  are the relative permeability and viscosity of the  $\beta$ -phase, respectively. Two algebraic equations relating the saturations and pressures, complete the system:

$$S_b + S_g = 1, \quad p_g - p_b = PC(S_b), \quad (3)$$

where  $PC$  is the capillary pressure. The relationship between capillary pressure and CO<sub>2</sub> saturation may be represented by the following potential function:

$$PC(S_g) = PC_{max} \left( \frac{S_g - S_{gc}}{1 - S_{gc} - S_{bc}} \right)^{n_c}, \quad (4)$$

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where  $S_{gc}$  and  $S_{bc}$  are the saturations at which CO<sub>2</sub> and brine phases become mobile, respectively.

The solution of the Black-Oil model is obtained employing the public domain software BOAST (Fanchi, 1997) which solves the differential equations using the IMPES algorithm (Implicit Pressure Explicit Saturation), based on a finite difference technique (Aziz and Settari, 1985).

### SEISMIC MODELING

#### Mesoscopic attenuation effects

An important mechanism of P-wave attenuation and dispersion at seismic frequencies is known as mesoscopic loss, due to heterogeneities larger than the pore size but much smaller than the predominant wavelengths (mesoscopic-scale heterogeneities). Due to the extremely fine meshes needed to properly represent these type of media, numerical simulation at the macroscale is very expensive or even not feasible. Our approach to solve this problem consists in applying an upscaling procedure to include the mesoscale effects in the macroscale (Santos et al., 2008). Within the formation, in zones where CO<sub>2</sub> is present, we determine complex and frequency dependent P-wave modulus  $E(\omega) = \lambda(\omega) + 2\mu(\omega)$  at the mesoscale using White's theory for patchy saturation (White et al., 1975), where  $\lambda(\omega)$  and  $\mu(\omega)$  are the Lamé coefficients and  $\omega$  is the angular frequency. Shear wave attenuation is taken into account using another relaxation mechanism, related to the P-wave White mechanism, to make the shear modulus  $\mu(\omega)$  complex and frequency dependent. Outside the formation and in zones where only brine is present, the complex bulk and shear moduli as function of frequency are determined using a Zener model. These complex moduli define an equivalent viscoelastic model at the macroscale that takes into account dispersion and attenuation effects occurring at the mesoscale.

#### Constitutive Relations

Let  $u = u(\omega) = (u_x(\omega), u_z(\omega))$  be the time-Fourier transform of the displacement vector. Therefore, the stress-strain relations in the space-frequency domain are:

$$\sigma_{jk}(u) = \lambda(\omega)\nabla \cdot u \delta_{jk} + 2\mu(\omega)\varepsilon_{jk}(u), \quad (5)$$

where  $\sigma_{jk}(u)$  is the stress tensor,  $\varepsilon_{jk}(u)$  is the strain tensor,  $\delta_{jk}$  is the Kroenecker delta, and  $\lambda(\omega), \mu(\omega)$  are the complex Lamé coefficients determined using White's theory.

#### Phase velocities and attenuation coefficients

For isotropic viscoelastic solids, the frequency dependent phase velocities  $v_t(\omega)$  and quality factors  $Q_t(\omega)$  ( $t = p, s$ ) are defined by the relations

$$v_t(\omega) = \left[ \text{Re} \left( \frac{1}{v_{ct}(\omega)} \right) \right]^{-1}, \quad Q_t(\omega) = \frac{\text{Re}(v_{ct}(\omega)^2)}{\text{Im}(v_{ct}(\omega)^2)}, \quad (6)$$

where  $v_{cp}(\omega), v_{cs}(\omega)$  are the complex and frequency dependent compressional and shear velocities defined as

$$v_{cp}(\omega) = \sqrt{\frac{E(\omega)}{\rho}}, \quad v_{cs}(\omega) = \sqrt{\frac{\mu(\omega)}{\rho}} \quad (7)$$

with  $\rho$  the bulk density.

#### A viscoelastic model for wave propagation

The equation of motion in a 2D isotropic viscoelastic domain  $\Omega$  with boundary  $\partial\Omega$  is (Santos et al., 2008) :

$$\omega^2 \rho u + \nabla \cdot \sigma(u) = f(x, \omega), \quad \Omega, \quad (8)$$

with first-order absorbing boundary condition:

$$-\sigma(u)\nu = i\omega \mathcal{D}u, \quad \partial\Omega. \quad (9)$$

Here  $f(x, \omega)$  is the external source and  $\mathcal{D}$  is a positive definite matrix which definition is given in (Ha et al., 2002).

The numerical solution is computed at a selected number of frequencies in the range of interest using an iterative finite element domain decomposition procedure. The time domain solution is obtained using a discrete inverse Fourier transform (Ha et al., 2002). To approximate each component of the solid displacement vector we employ a nonconforming finite element space which generates less numerical dispersion than the standard bilinear elements.

### PETROPHYSICAL MODEL OF A SHALY SANDSTONE

The pressure dependence of properties is based on the following relationship between porosity and pore pressure  $p(t) = S_b p_b(t) + S_g p_g(t)$ ,

$$\frac{(1 - \phi_c)}{K_s} (p(t) - p_H) = \phi_0 - \phi(t) + \phi_c \ln \frac{\phi(t)}{\phi_0}, \quad (10)$$

where  $\phi_c$  is a critical porosity,  $\phi_0$  is the initial porosity at hydrostatic pore pressure  $p_H$  and  $K_s$  is the bulk modulus of the solid grains (Carcione et al., 2003). The rock is formed with quartz (bulk modulus of 40 GPa) and clay (bulk modulus of 15 GPa).  $K_s$  is computed as the arithmetic average of the Hashin Shtrikman upper and lower bounds. The relationship among horizontal permeability, porosity and clay content  $C$  is (Carcione et al., 2003),

$$\frac{1}{\kappa_x(t)} = \frac{45(1 - \phi(t))^2}{\phi(t)^3} \left( \frac{(1 - C)^2}{R_q^2} + \frac{C^2}{R_c^2} \right), \quad (11)$$

where  $R_q$  and  $R_c$  are the average radii of the sand and clay grains. Also, as permeability is anisotropic, we assume the following relationship between horizontal and vertical permeability  $\kappa_z$  (Carcione et al., 2003)

$$\frac{\kappa_x(t)}{\kappa_z(t)} = \frac{1 - (1 - 0.3a) \sin \pi S_b}{a(1 - 0.5 \sin \pi S_b)}, \quad (12)$$

where  $a$  is the permeability-anisotropy parameter. The bulk modulus of the dry matrix,  $K_m$ , is computed using the Krief equation (Krief et al., 1990) as follows:

$$K_m(t) = K_s (1 - \phi(t))^{A/(1 - \phi(t))}. \quad (13)$$

Assuming a relation  $K_m/\mu_m = K_s/\mu_s$ , we set

$$\mu_m(t) = \mu_s (1 - \phi(t))^{A/(1 - \phi(t))}. \quad (14)$$

Using the moduli  $K_s, K_m, \mu_m$ , the porosity  $\phi$  and permeabilities  $\kappa_x, \kappa_z$ , as well as the fluids bulk moduli and viscosities (computed using the Peng-Robinson model (Peng and Robinson, 1976)), we determine the complex and frequency dependent Lamé coefficients  $\lambda(\omega), \mu(\omega)$  as explained before.

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### EXAMPLES

To test the proposed methodology, we consider a model of the Utsira formation having 1.2 km in the  $x$ -direction, 10 km in the  $y$ -direction and 0.4 km in the  $z$ -direction (top at 0.77 km and bottom at 1.17 km b.s.l.). The pressure-temperature conditions are  $T = 31.7z + 3.4$ , where  $T$  is the temperature (in °C) and  $z$  is the depth (in km b.s.l.);  $p_H = \rho_b g z$  is the hydrostatic pressure, with  $\rho_b = 1040 \text{ kg/m}^3$  the density of brine and  $g$  the gravity constant. Within the formation, there are several mudstone layers which act as barriers to the vertical motion of the CO<sub>2</sub>. The initial porosity  $\phi_0$  (at hydrostatic pore pressure) for the Utsira sandstone is assumed to have a fractal spatial distribution, based on the so-called von Karman self-similar correlation functions (Frankel and Clayton, 1986). Horizontal and vertical permeabilities were determined by using equations (11) and (12), considering an anisotropy parameter  $a = 0.1$  and a fixed clay content  $C = 6 \%$ . We describe the results obtained using two different capillary pressure curves, that are shown in Figure 1.

- $PC_{max}=30 \text{ kPa}$ ,  $n_c=4$
- $PC_{max}=100 \text{ kPa}$ ,  $n_c=2$

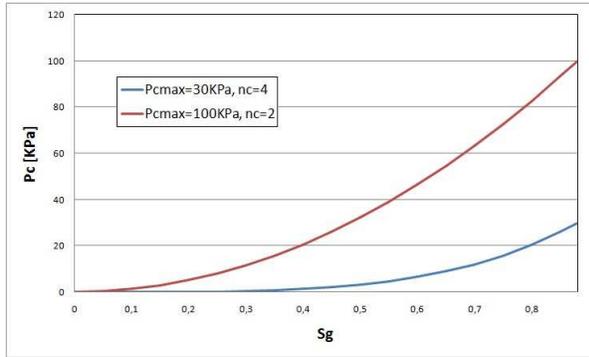
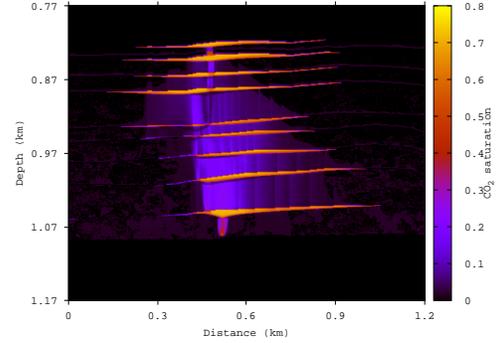


Figure 1: Capillary pressure curves.

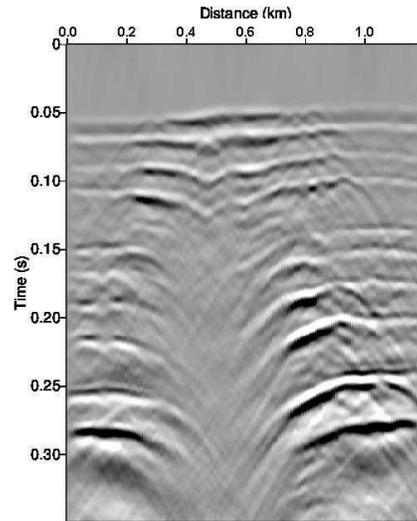
### CO<sub>2</sub> injection

CO<sub>2</sub> is injected during seven years in the Utsira formation at a constant flow rate of one million tons per year. The injection point is located at the bottom of the formation:  $x = 0.6 \text{ km}$ ,  $z = 1.082 \text{ km}$ . The simulation uses a mesh with equally-spaced blocks in each direction:  $n_x = 300$  in the  $x$ -direction,  $n_y = 5$  in the  $y$ -direction and  $n_z = 400$  in the  $z$ -direction. Actually the model is 2.5D since the properties are uniform along the  $y$ -direction, which has an extension of 10 km. The source is located at the third grid point along this direction. To satisfy the CFL stability condition due to IMPES formulation (Savioli and Bidner, 2005), the time step is 0.05 d. Recall that the petrophysical properties of the formation are time dependent due to the CO<sub>2</sub> injection and consequent increase in pore fluid pressure, but they change at a much slower rate than pressure and saturations. As a consequence, we have two time scales, and we use a much larger time step to update petrophysical properties than to run the flow simulator. In this work, the petrophysical properties are updated every year.

Figures 2(a) and 3(a) show 2D vertical slices (corresponding to  $n_y = 3$ ) of the CO<sub>2</sub> saturation fields after three years of CO<sub>2</sub> injection, computed using the lower and higher  $PC$  curves, respectively. In both cases, CO<sub>2</sub> accumulations below the mudstone layers can be observed. As injection proceeds, part of the injected fluid migrates upwards due to the openings in the mudstone layers that generate chimneys. Comparing both figures, when capillary pressure is higher, CO<sub>2</sub> upward migration is slower and, consequently, we can observe thicker zones of accumulations below the layers and high CO<sub>2</sub> saturations levels between layers.



(a)



(b)

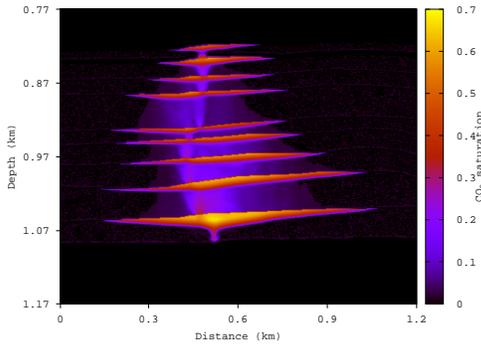
Figure 2: CO<sub>2</sub> saturation distribution and synthetic seismogram after three years of CO<sub>2</sub> injection. Lower capillary pressure,  $PC_{max}=30 \text{ kPa}$ ,  $n_c=4$ .

### Seismic Monitoring

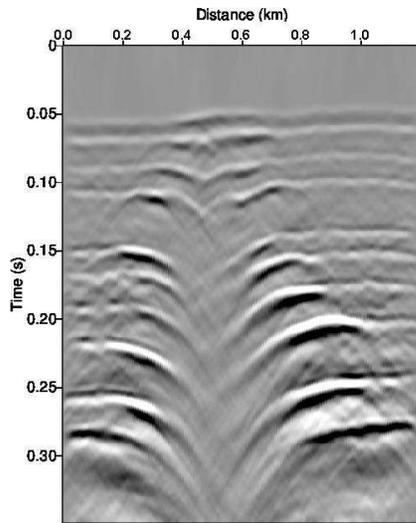
In this section, we analyze the capability of seismic monitoring to identify zones of CO<sub>2</sub> accumulation and migration. With this purpose, we use 2D slices of CO<sub>2</sub> saturation and fluid pressure obtained from the flow simulator to construct a 2D model of the Utsira formation. The seismic source is a spatially lo-

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calized plane wave of main frequency 60 Hz. Figures 2(b) and 3(b) display seismograms after three years of CO<sub>2</sub> injection associated with the CO<sub>2</sub> saturations shown in Figures 2(a) and 3(a), respectively. A standard  $f$ - $k$  filter is applied to the seismic sections. The reflections seen in those seismograms show the CO<sub>2</sub> accumulations below the mudstone layers as the injection proceeds. In particular, the pushdown effect observed in the real seismograms (Chadwick et al., 2009) due to CO<sub>2</sub> accumulations is clearly observed. But the seismograms obtained from the two capillary pressure curves are quite different. The difference in P-wave phase velocities and attenuation coefficients is clearly observed in Figure 4. Therefore, the influence of capillary forces is very important in CO<sub>2</sub> storage and a precise estimation of this curve is needed to perform realistic long term predictions.

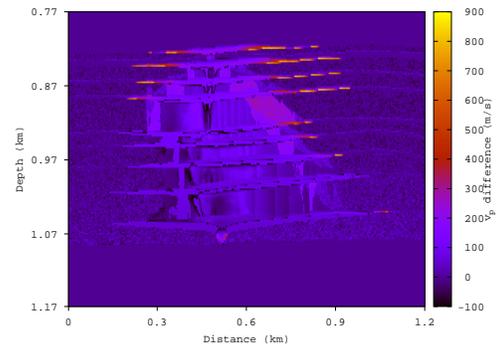


(a)

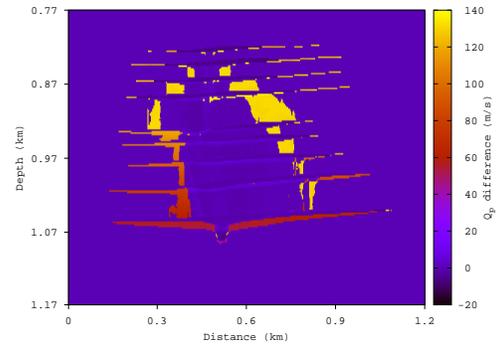


(b)

Figure 3: CO<sub>2</sub> saturation distribution and synthetic seismogram after three years of CO<sub>2</sub> injection. Higher capillary pressure,  $PC_{max}=100$  kPa,  $n_c=2$ .



(a)



(b)

Figure 4: Difference between P-wave phase velocities and attenuation coefficients obtained with higher and lower capillary pressure

## CONCLUSIONS

The fluid-flow simulator yields CO<sub>2</sub> accumulations below the mudstone layers and the corresponding synthetic seismograms resemble the real seismic data, where the pushdown effect is clearly observed. Capillary forces affect the migration and dispersal of the CO<sub>2</sub> plume; higher values of this curve cause a slower CO<sub>2</sub> upward migration and thicker zones of CO<sub>2</sub> accumulations. The influence of capillary forces is clearly observed in the synthetic seismograms.