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# Validating compositional fluid flow simulations using 4D seismic interpretation and vice versa in the SECARB Early Test—A critical review

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## Abstract

This paper discusses strengths and weaknesses of 4D seismic interpretation as a technique for monitoring carbon dioxide that was injected as part of a large scale test associated with commercial enhanced oil recovery (EOR) at Cranfield Field, Mississippi, USA. The goals of the monitoring effort are 1) to make measurements to verify that the CO<sub>2</sub> is contained in the reservoir according to operational designs and model predictions, and 2) that if there are deviations, to provide data which can be used to update the earth models and determine if any mitigation is needed. The current work uses a compositional numerical simulation to model CO<sub>2</sub> flow in the reservoir and compares the results with estimates of changes in seismic properties between a pre-injection and a survey after more than 2 million metric tons injected. The complicated physics of the problem in Cranfield field present challenges to seismic interpretations. Our results show partial agreement between the results of the numerical simulation and the seismic interpretations. Possible causes of discrepancies among the fluid flow model and multiple interpretations of the same seismic data sets performed with different work flows illuminates the types of uncertainties that should be considered to achieve the goals of monitoring.

## Keywords

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CO<sub>2</sub> injection monitoring; 4D seismic; Compositional simulation; Enhanced oil recovery

## **1. Introduction**

An important aspect of underground CO<sub>2</sub> injection, as an enhanced oil recovery method or as part of carbon capture and storage, is to monitor and track the movement of the carbon dioxide. This will help in maintaining a more effective sweep during EOR operation or improved effective storage during CO<sub>2</sub> sequestration. An increasing amount of time-lapse surface seismic work is being used for CO<sub>2</sub> injection monitoring. Arts et al. (2004) presented the results of seismic interpretations to monitor CO<sub>2</sub> injection into a saline aquifer at the Sleipner site. Their pioneering work was one of the first to use repeated time-lapse seismic surveys to monitor a large CO<sub>2</sub> injection and storage operation. Several subsequent seismic monitoring efforts have been conducted, such as those at the Weyburn (Davis et al., 2003), Otway (Urosevic et al., 2011), and Ketzin (Ivanova et al., 2012) reservoirs.

Alfi and Hosseini (2016) analyzed CO<sub>2</sub> injection into the Lower Tuscaloosa Formation of Cranfield field by comparing the results of a black oil numerical simulation of CO<sub>2</sub> movement with several seismic interpretations (Carter, 2014; Ditekof et al., 2011; Zhang et al., 2014). Their analyses showed that the flow simulation results and the seismic interpretations agree in some locations while diverging in others. In this paper, we will investigate the possible reasons for the observed discrepancies. Seismic technology provides a key advantage in monitoring CO<sub>2</sub> injection sites by covering a large acreage in a single survey. However, the application of 4D seismic for CO<sub>2</sub> plume tracking has always been hampered by several different factors (Gendrin et al., 2013; Ivanova et al., 2012; Roach et al., 2016). One challenge is the depth of the formation of interest. Generally, the deeper the reservoir, the harder it becomes to detect the induced changes in seismic signals because of CO<sub>2</sub> injection. Formation geology and rock properties, e.g. high rock stiffness, are additional factors that can negatively affect the reliability of seismic interpretations. The sensitivity of seismic surveys to changes in acoustic impedance make them highly sensitive in some, but not all geological contexts. Such effects can be magnified in environments with higher noise levels (e.g. onshore fields). Formation heterogeneity is another factor that influence our seismic interpretations. Regions with low porosity rocks, the presence of low-density layers (e.g.

coal) in the overburden, or non-homogeneous mixing of CO<sub>2</sub> with the formation fluid leading to the formation of patches, are examples of the effect of heterogeneity. Properties of the reservoir pore fluids prior to CO<sub>2</sub> injection and the presence of residual gas or a low-density oil phase in the reservoir provides another challenge for seismic monitoring of CO<sub>2</sub> injection, as will be discussed in more detail in the following sections. In addition, thickness of the accumulated carbon dioxide itself can impact the interpretation results. Thin plumes, below about one quarter of a wavelength, will not be detected in seismic survey.

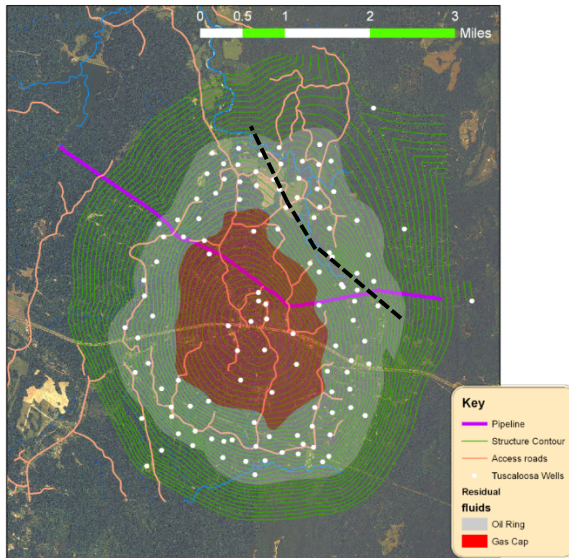
Alfi and Hosseini (2016) modeled the CO<sub>2</sub> injection process using a black oil simulator (CMG-IMEX). Such an approach is effective specifically for large models with a significant number of grid blocks as it reduces the computational cost. On the other hand, black oil simulators come with the cost of not accurately modeling the miscibility of the CO<sub>2</sub> and reservoir fluids (oil and brine). Hence, in the current paper, we have improved the model by using a compositional simulator (CMG-GEM), which allows us to account for the CO<sub>2</sub> miscibility in oil and brine. Compositional models are better suited to deal with interaction between the various chemical components in the system, such as gas exsolution and dissolution. Accurate modeling of the distribution of gas is of a particular interest in this study and will be discussed in greater detail later in this paper.

In addition, in this study a new set of seismic interpretations were performed and compared to the interpretations by Carter (2014) and Ditkof et al. (2011) . Such a comparison helps us to understand how various processing and workflow choices impact seismic interpretations.

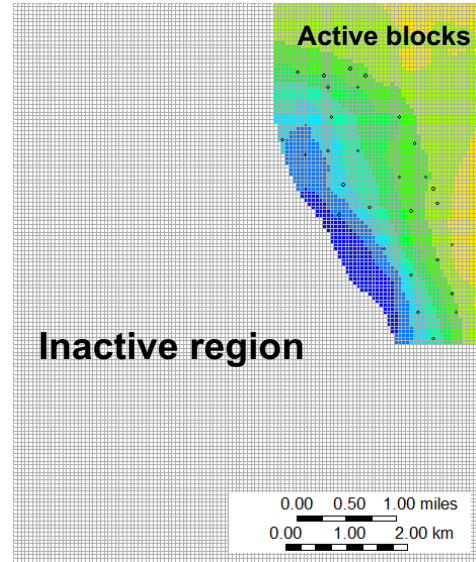
This paper begins with a descriptive explanation of the reservoir static and dynamic model. Oil, gas, and water production data are history matched and the model is tuned accordingly. Moreover, we explain how seismic data were prepared and interpreted. Seismic interpretations (from this paper and previous works) are then compared with dynamic simulation results, following a critical discussion on the role of different reservoir related factors on seismic data interpretation.

## **1. Numerical model description**

80 In this study, we use data collected from the Cranfield site to history match fluid flow simulations.  
81 The Cranfield field is located in Adams and Franklin Counties, Mississippi, east of the town of  
82 Natchez. The original productive area of the reservoir was estimated to be 31.3 km<sup>2</sup> with a  
83 producing depth range of 3060 to 3193 m (Weaver and Anderson, 1966). An active aquifer  
84 downdip of the reservoir provides a constant pressure boundary. The initial reservoir  
85 temperature was reported to be 125°C with an initial reservoir pressure of 32.4 MPa at 3040 m.  
86 A central graben-bounding sealing fault, down dropped to the west, divides the productive zone  
87 into two reservoirs, and creates a trap on the east side (the dashed NW-SE line in Figure 1a).  
88 More details about reservoir specifications, production history, simulation projects, and  
89 monitoring efforts can be found in other publications (Alfi and Hosseini, 2016; Alfi et al., 2015;  
90 Choi et al., 2011; Hosseini et al., 2013; Hovorka et al., 2013; Weaver and Anderson, 1966). The  
91 current work focuses on the northeastern part of the reservoir, assuming a negligible fluid  
92 movement between this part and the rest of the reservoir, as the small contact area and balanced  
93 field development plan minimizes such flow. This approach has substantially decreased the  
94 number of active grid blocks in the system, in order to reduce the computational cost because  
95 compositional models increase the computational burden due the higher number of components  
96 in the system. A Cartesian model with 124 ×149×20 (X×Y×Z) grid cells is used to model a total  
97 reservoir area of 7.5 km by 9.1 km with a total reservoir thickness of 24.4 m. Total number of grid  
98 blocks is 369,520, out of which 82,559 grid blocks located at the northeastern part of the  
99 reservoir are considered to be active (Figure 1b). All grid blocks have a uniform dimension of  
100 61×61×1.2 m. The simulation model is isothermal and does not account for the possible reactions  
101 between minerals, injected and in situ fluids.



a) Cranfield site



b) Simulation gridding

Figure 1-a) Structural contour map of Cranfield. The black dashed line represents the sealing fault that separates the northeastern section of the reservoir from the rest. b) Reservoir simulation model to simulate the CO<sub>2</sub> injection process. Simulation effort focuses on the northeastern side of the reservoir so the rest of model is inactivated to reduce the computational load. Color indicates depth, showing the structural trap formed against the fault.

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103 Black oil models, although very efficient in less complicated cases, fail to address  
 104 thermodynamics of CO<sub>2</sub>/hydrocarbon interactions that are necessary to correctly simulate  
 105 miscible processes at work at Cranfield. In this study, we employ the computer modeling group's  
 106 (CMG) compositional simulator (GEM), an advanced general equation of state compositional  
 107 simulator, designed for various complicated fluid flow problems including CO<sub>2</sub> miscible flood. The  
 108 simulator was set up to model three fluid phases including water, oil, and gas.

109 The Peng Robinson equation of state (Peng and Robinson, 1976) is used in this study to model  
 110 reservoir fluid properties. Our fluid model is composed of seven different components including  
 111 CO<sub>2</sub>. The thermodynamics model and component properties were tuned based on the fluid data  
 112 published by Weaver and Anderson (1966). The data used for this purpose included bubble point  
 113 pressure, solution gas-oil-ratio, formation volume factor, oil and gas viscosities. The CO<sub>2</sub>/brine  
 114 solubility in the current model is accounted for by using the Henry's constant. The experimental

data from Duan and Sun (2003) are used to find the appropriate solubility parameters. Relative permeability curves are generated from the data published by Weaver and Anderson (1966). Keeping the residual values intact from the original relative permeability models, the relative permeability values at the endpoints were slightly modified to match the field production data. The slight modification of the relative permeability data (Figure 2) shows a agreement between the relative permeability set used in this study and the originally published data by Weaver and Anderson (1966).

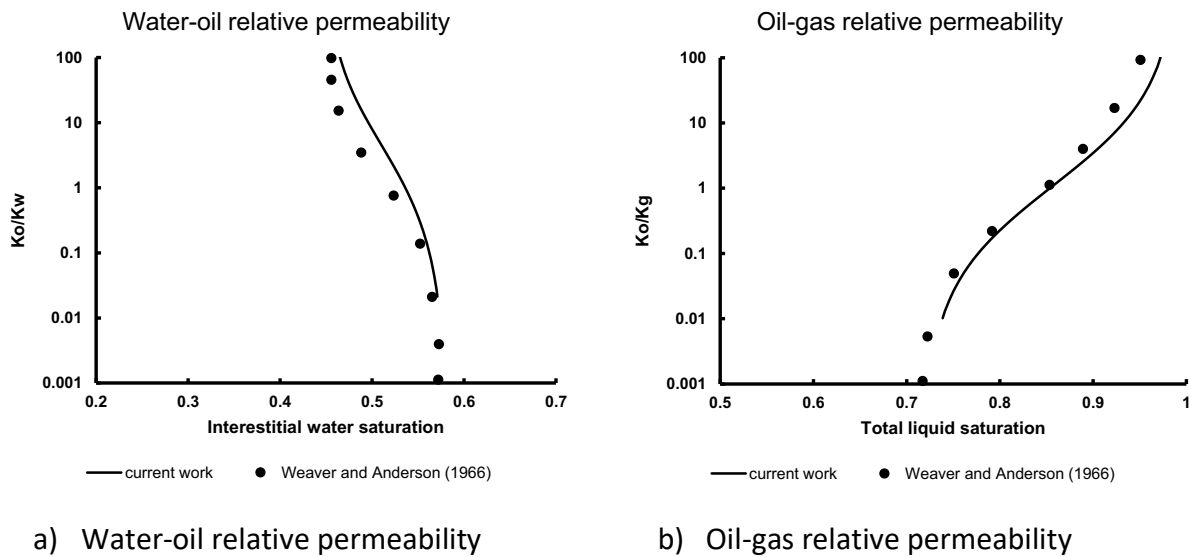


Figure 2-Relative permeability curves used in the current study are slightly modified from the original Weaver and Anderson (1966) values to match the oil, gas, and water production data during the CO<sub>2</sub> injection period. The new permeability sets are still in a good agreement with the original experiments.

The original permeability and porosity model is from Hosseini et al. (2013). They used a purely mechanistic approach to recognize eight different facies (flow zones) in the reservoir based upon combined permeability and porosity data sets. Each flow zone is given a porosity and permeability, which is constant for the entire zone. These flow zones are categorized under two main groups of low permeability (shale) and high permeability (sand) zones. High permeability channels in the lower Tuscaloosa Formation at Cranfield have been mapped and confirmed by various fluid-flow responses (Lu et al., 2013; Mukhopadhyay et al., 2015). However, the geometry of high permeability channels throughout the field presented by Hosseini et al. (2013) is based on probabilistic extrapolation of the available data. In this study, the permeability

contrast between the sand and shale facies and the presence of high permeability flow paths was examined as a history matching parameter to fit the observed field data. Table 1 provides some of the more important reservoir properties used in this work.

Parameter	Value	Parameter	Value
Kh/Kv (permeability anisotropy)	0.05	Permeability range	0.016 – 4422 mD
Brine salinity	150,000 ppm	Porosity range	0.0002 – 0.45
Initial bubble point pressure	31.7 MPa	Irreducible water saturation	0.45
Original water oil contact depth	3064 m	Critical gas saturation	0.01
Original gas oil contact depth	3008 m	Aquifer model	Carter and Tracy (1960)

Table 1—Reservoir related parameters used in numerical modeling of Cranfield

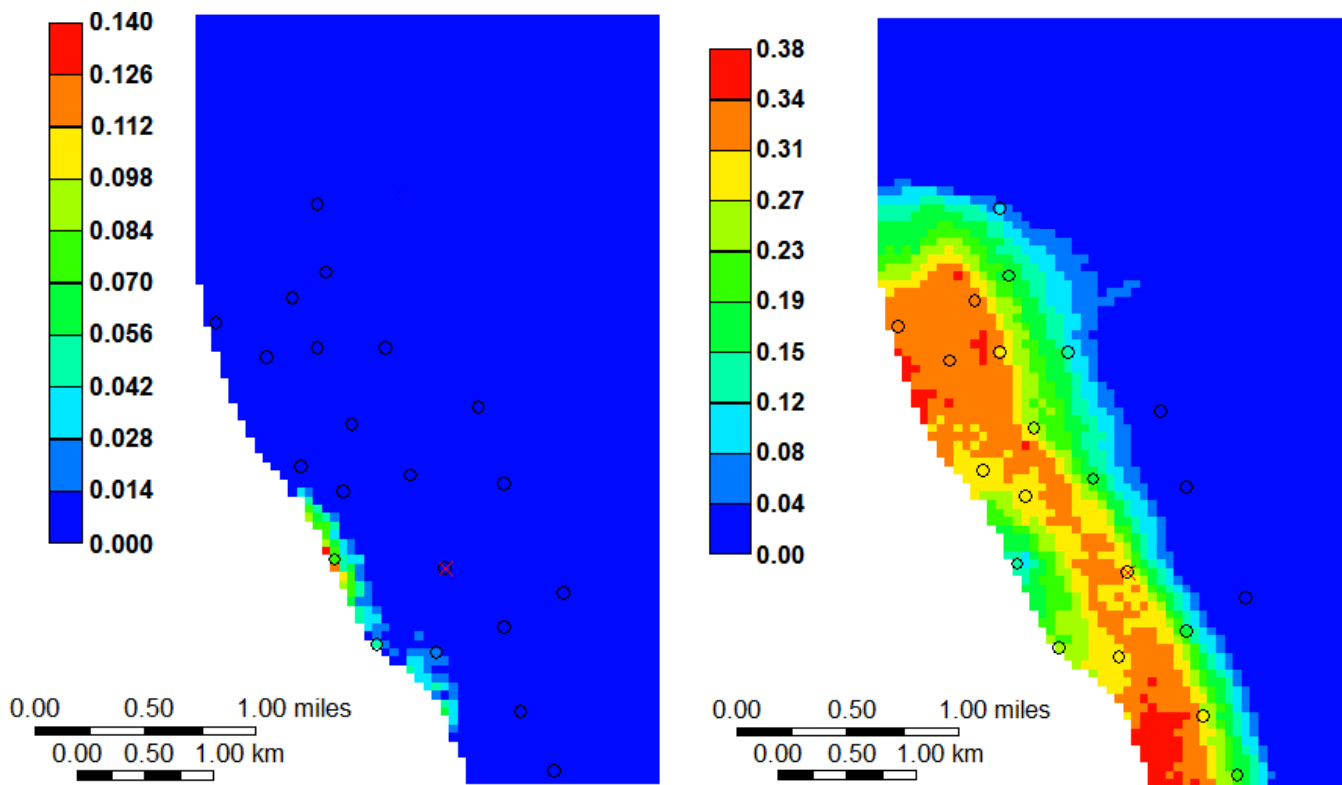
## 2. History matching and reservoir performance

The first oil producing well from Cranfield was drilled in 1944. Since then, a productive area of about 7,750 acres was defined by 93 producing wells. The oil wells were drilled based on a 40-acres spacing while the spacing for the gas wells were 320 acres. The reservoir consists of an oil ring overlain by a large central gas cap (Weaver and Anderson, 1966). A cycling and condensate extraction plant was used to reinject the produced methane from Tuscaloosa and Paluxy reservoirs back into the gas cap. By 1951, the injected gas had broken through to many of the oil zone wells. The gas cycling continued until 1960 with dry gas sweeping the gas cap and the oil zone. Although the gas injection was intended to avoid, or slow down, the pressure depletion in the reservoir, reservoir pressure gradually fell below 27.6 MPa (4000 psi) causing water to encroach into the oil zone with increasing production. By the beginning of 1960, most of the wells had either a 100% water cut or a GOR greater than 100,000 scf/STB with an average field water cut equal to 88% and GOR equal to 85,000 scf/STB. The sale of gas cap methane and condensate started at that point. At the same time, water was produced in high volumes to prevent the aquifer from pushing the remaining oil into the gas cap. Gas injection stopped in 1964 when the project was near its economic limit. Production from the field was halted in 1966 and the



reservoir was abandoned. Over the next several decades, a strong water drive restored pressure to near-initial levels.

Figure 3 shows the modeled oil and gas saturation in the abandoned reservoir after four decades of recovery, just before the initial seismic survey was conducted in 2007. According to the model, residual gas migrated toward the top of the structure and accumulated in a structural trap east of the sealing fault. The modeled residual oil saturations is as high as 0.35 in the oil rim. As explained later, the presence of oil and gas with compressibilities larger than brine, can influence the sensitivity of seismic signals as CO<sub>2</sub> is injected. Therefore, it is important to quantify the amounts of residual oil and gas before conducting the seismic survey. No log or core-based measurements of hydrocarbon saturations from this pre-injection interval are available.



a) Gas saturation distribution at the time of baseline seismic survey.

b) Oil saturation distribution at the time of baseline seismic survey.

Figure 3- Modeled oil and gas saturation in the reservoir at the time of the first seismic survey in 2007.

In 2008, CO<sub>2</sub>-based enhanced oil recovery (EOR) was initiated by Denbury Onshore, LLC to sweep the bypassed residual oil. Between 2008 and 2015, more than half of the oil ring (Figure 1a) was

accessed using an irregular five-spot injection pattern with continuous CO<sub>2</sub> injection. Unlike many enhanced oil recovery efforts, there was no accompanying water injection at Cranfield. Initial patterns began in the northern part of the field and continued clockwise around the oil ring as the field was developed. CO<sub>2</sub> injection in the northeastern part of the reservoir commenced on July 2008 and continued until October 2012. Data gathered over 4 years of CO<sub>2</sub>-EOR were used to calibrate the numerical model for a more reliable forecast of reservoir response. A total of 22 wells were considered for simulation of the northeast segment of the reservoir, 12 injection wells and 10 production wells. Injection wells were modeled based upon the injected CO<sub>2</sub> volumes available at each well location. To account for the amount of CO<sub>2</sub> entering the study area, a well located at the edges of the area of interest was modeled with half of the actual injection rate. The other half is assumed to enter that part of the reservoir that is not included in the model. Oil, water and gas productions rates were also available for all the producing wells during the simulation period. Monthly oil production rates were used as the operational constraint for the production wells. Cumulative production was calculated based on the fluid produced after July 2008. The history matching process in this section was based on observing/matching reservoir response by tuning the relative permeability data, adjusting the geological model to account for the high permeability zones, and altering the permeability contrast in high permeability channels in the reservoir. As discussed above, the presence of high permeability channels in the Tuscaloosa injection zone has been noted by other authors. Analyzing the breakthrough times (the time it takes for CO<sub>2</sub> to reach the production wells) confirms the presence of high permeability channels because CO<sub>2</sub> reaches some of the producing wells in a relatively short time after the start of injection, despite the spacing between the injection and production wells.

Figure 4 compares the cumulative oil, gas, and water production from the simulation and field data. The results show a reasonably good agreement between the field and simulation data, which accounts for the reliability of the numerical model to perform further analyses.

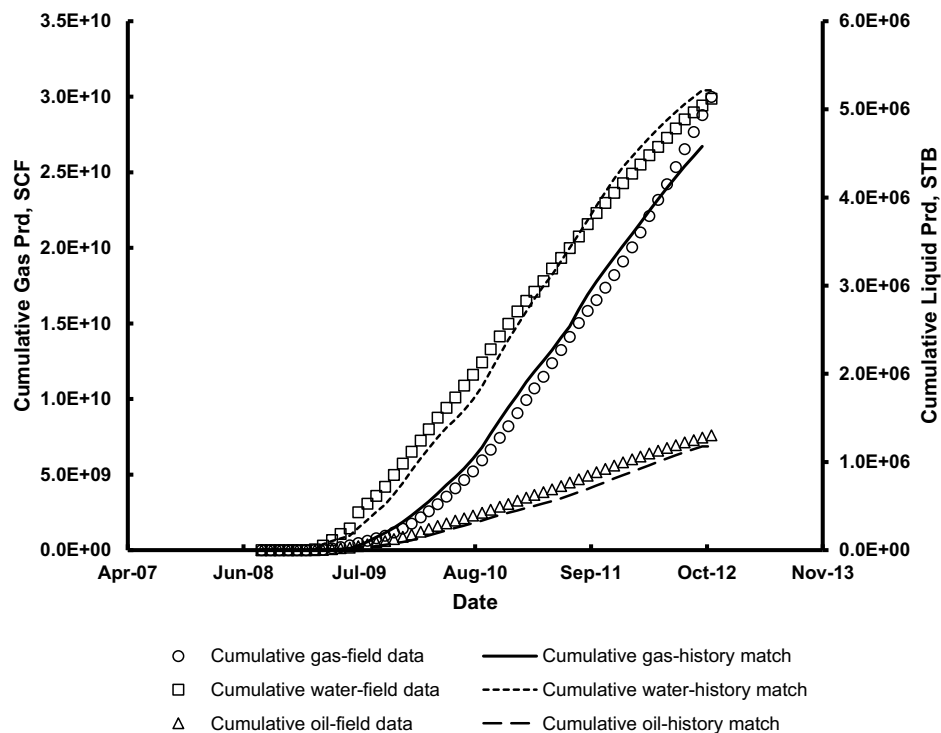


Figure 4-History matching results of 4 years' worth of production data shows an agreement between the field production data and simulation results.

190 As a way to verify the reliability of the simulation results, Figure 5 compares the CO<sub>2</sub>  
 191 breakthrough times of the simulation with that of the field data. As we can see, the model is able  
 192 to predict the CO<sub>2</sub> migration though the reservoir with an acceptable accuracy. This is significant  
 193 because it implies that the modeled thickness and saturation between the wells are reasonable.  
 194

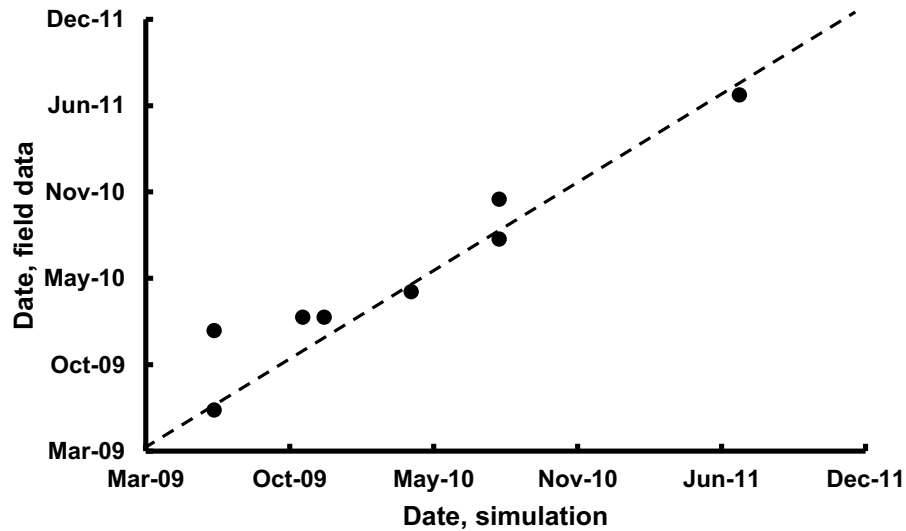


Figure 5-Breakthrough times for the producing wells show a good agreement between the simulation results and field data (x-axis shows the breakthrough times from simulation and y-axis shows that of real field data).

Next, we consider the CO<sub>2</sub> distribution after four years of CO<sub>2</sub> injection into the field. Figure 6 shows the average CO<sub>2</sub> saturation in the northeastern part of the reservoir at October 2012, the saturation of each plotted grid is averaged over the multiple vertical grids at that location. As is evident from the figure, carbon dioxide accumulates around the injection wells. CO<sub>2</sub> is also observed to migrate laterally and vertically below the former oil-water contact from the eastern and northern margins of the reservoir. New CO<sub>2</sub> injectors drilled near or below the oil water contact to assure that the base of the oil column was swept and to inject high volumes to support the SECARB project goals. High permeability in the lower Tuscaloosa channel facies also allowed migration below the oil-water contact as observed by (Lu et al., 2013). Looking at the average reservoir pressure from the simulation results ([Figure 7, Choi et al. (2011)], one notes that reservoir pressures rise as high as 37 MPa due to CO<sub>2</sub> injection. Such a high pressure in the reservoir can direct CO<sub>2</sub> into the lower pressure edges of the reservoir where the aquifer pressure is around 32 MPa, resulting in CO<sub>2</sub> migration off the structure.

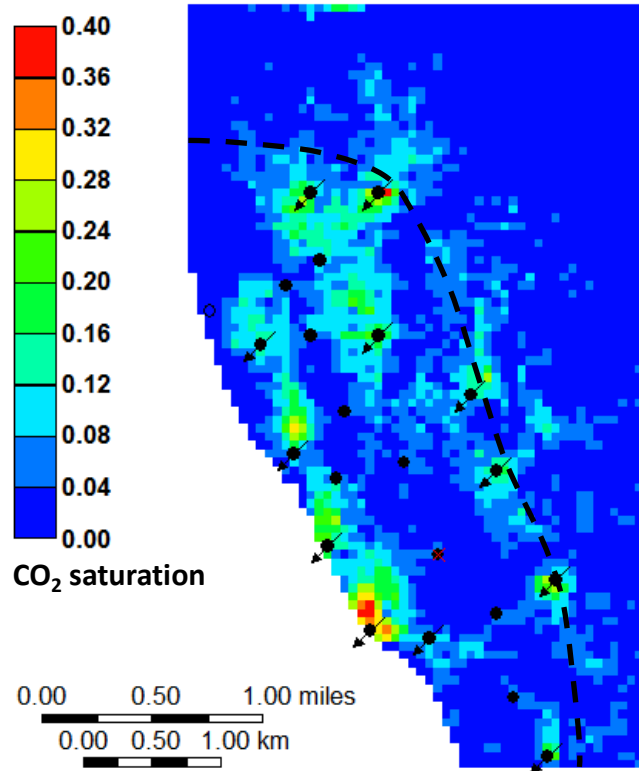
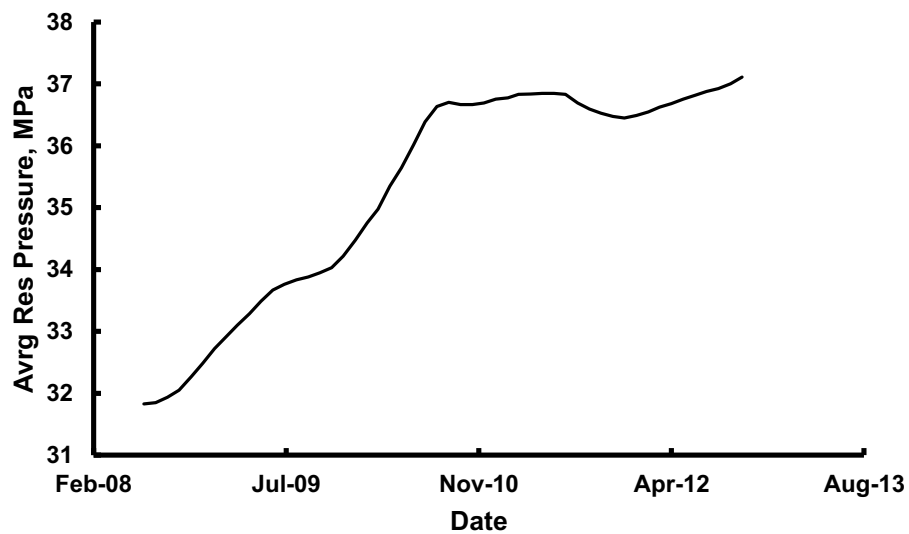


Figure 6-Modeled CO<sub>2</sub> saturation distribution map at October 2012 shows CO<sub>2</sub> accumulated around the injection wells (black dots with arrows). CO<sub>2</sub> saturations of each grid cell is averaged over the local thickness of the zone. Water-oil contact illustrated by dash line.

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[Figure 7-Reservoir pressure increases after CO<sub>2</sub> injection commences at 2008, providing a driving force to move CO<sub>2</sub> below the original oil-water contact.

## 2. Seismic data interpretations

The changes in fluid saturations due to the injection and production operations at the Cranfield site lead to seismic velocity variations within the reservoir. Such changes can induce temporal differences in the amplitude of reflections from the top of the reservoir and within the reservoir itself. In addition, the velocity changes may induce time shifts in reflections within and below the reservoir. Such time shifts will lead to apparent amplitude changes as one set of reflections is offset in time from the other. In an effort to image seismic velocity variations associated with the injection of more than 2 million tons of CO<sub>2</sub>, two seismic surveys were collected, the base survey in 2007 before the start of injection and a follow-up survey in 2010. The two surveys were processed by the commercial contractor Geotrace following identical workflows. The changes in the seismic response between the two surveys have been analyzed previously by Zhang et al. (2013, 2012), Ditkof (2013), and Carter (2014). The pattern of amplitude changes is complicated and there is an imperfect correlation between areas where CO<sub>2</sub> was injected and locations of seismic amplitude changes. This shows a limitation in using the 4-D seismic to assess the distribution of CO<sub>2</sub> as well as using it to validate the fluid flow model.

Because the seismic amplitude response depends upon both the changes in the reflection coefficients and velocity-induced time-shifts of the peaks and troughs of the seismic traces, it may not have a simple relationship to velocity changes in the reservoir. As an alternative, we use the simpler method of interpreting the time shifts in the seismic traces. Ignoring any stress and strain changes in the reservoir amplitude changes should be directly related to velocity changes. To this end, we examined the time shifts between traces of the two seismic cubes produced by Geotrace. This approach is somewhat similar to the work of Ditkof (2013) but the implementation is significantly different. In particular, Ditkof (2013) used two large time windows from 0.35 to 1.80 s above the reservoir and 2.50 to 3.20 s below the reservoir. The lower time window is substantially below the reservoir and is likely to be influenced by a number of factors related to the seismic processing. Any time shift due to velocity changes in the reservoir should begin just below the reservoir and extend to a large distance below it. In an effort to isolate the time shift due to velocity changes within the reservoir, we line up traces using a time window of 0.20 s, extending upward from the top of the reservoir (Figure 8). The segments of the two traces are

aligned by cross-correlating them to produce a cross-spectrum (Figure 9). To obtain sub-sample precision we fit a quadratic polynomial function to the five points surrounding the peak of the cross-spectrum. Such local approximations are often used in determining minima and maxima, as quadratics are the lowest degree polynomials that can represent a single extremum. A linear form cannot represent a peak or a trough and a higher-order polynomial would introduce additional extrema. The time delay associated with the peak of the quadratic is used as the time shift. Once the traces are aligned in the window above the reservoir, the shift is applied to the segments of the traces in the window below the reservoir. Then, the relative shift is computed for these corrected traces using the cross-spectrum and the quadratic fitting technique.

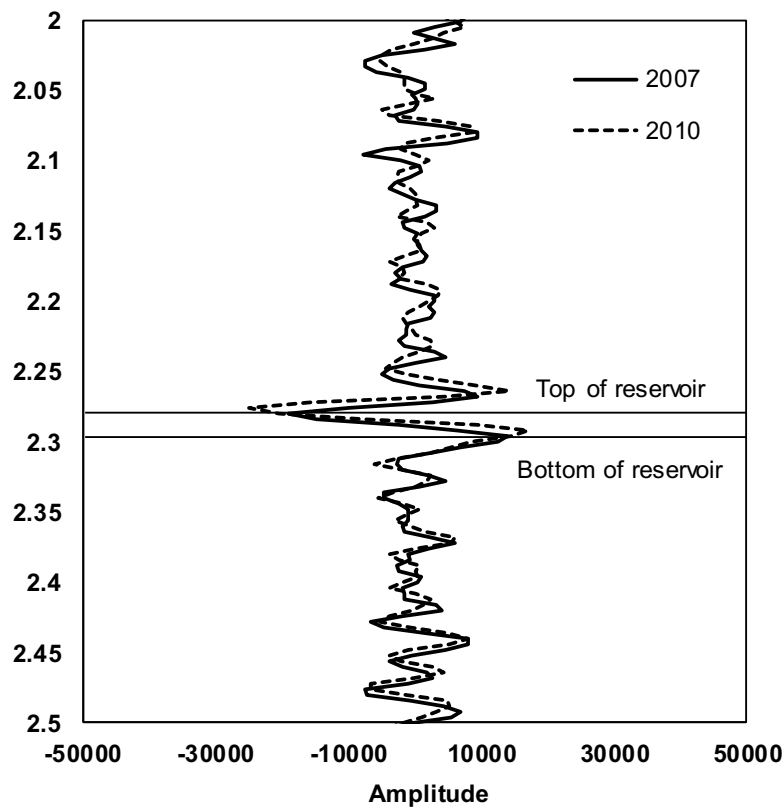


Figure 8- Two traces from the baseline (2007) and follow-up (2010) seismic surveys. The top and bottom of the reservoir are indicated by the horizontal lines.

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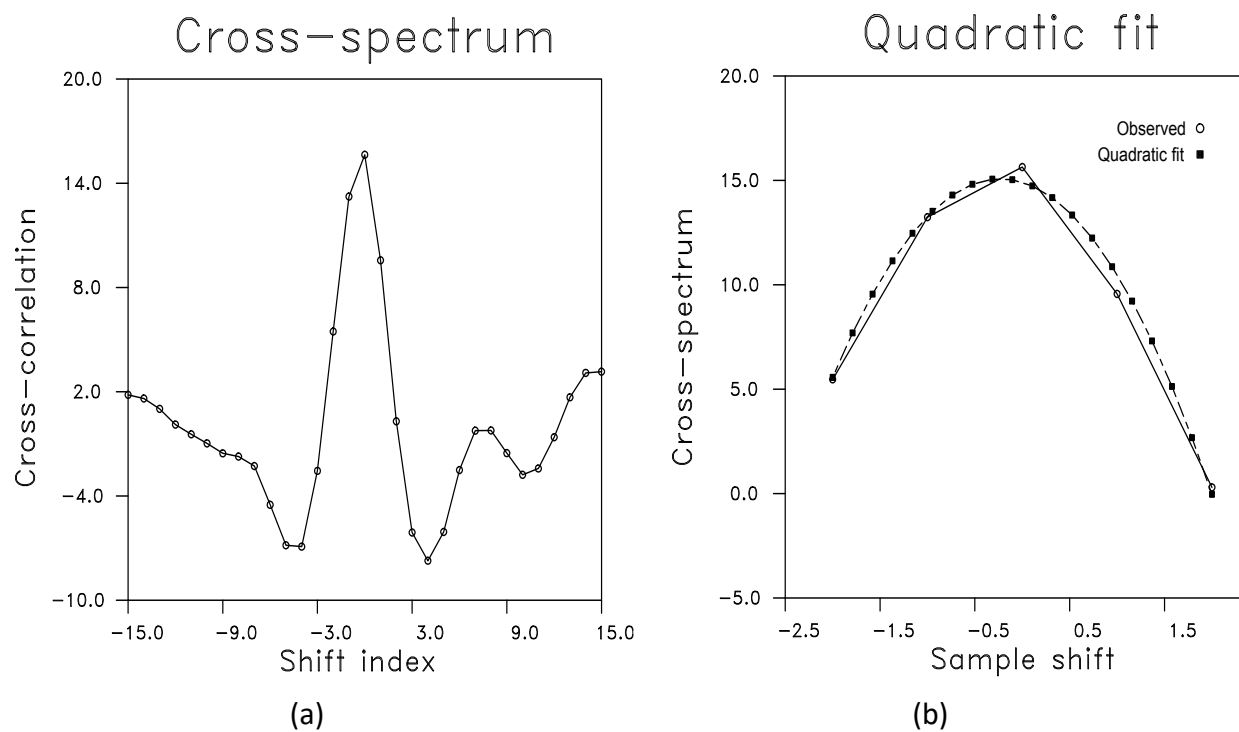


Figure 9-a) Cross-spectrum of the segment of the two traces for the window above the reservoir. The magnitude of the cross-correlation is plotted as a function of the shift. b) Quadratic fit (filled squares) to the five points surrounding the peak.

251

252 The approach was applied to every location for which there were traces for both the 2007 and  
 253 2010 surveys. Figure 10a shows the time shifts for the northeastern side of the reservoir. The  
 254 regions with greater time shifts indicate areas where there has been larger changes in acoustic



properties, primarily because CO<sub>2</sub> has replaced the reservoir fluid (either oil or brine). On the right side of Figure 10b we show the simulated CO<sub>2</sub> saturation at the time of the second seismic shoot (2010). The black circles in Figure 10b provide the locations of injection wells. Larger radii indicate more CO<sub>2</sub> injection. Field development and CO<sub>2</sub> injection commenced from the northern part of the study area, so more CO<sub>2</sub> has accumulated in those regions (larger circles). Comparing the history-matched simulation results with that of the seismic interpretations, one can claim that the seismic data provides a reasonable estimate of the CO<sub>2</sub> plume and its areal extent. In fact, the agreement suggests that time-lapse seismic surveys may be used to monitor subsurface CO<sub>2</sub> injection, provided that the actual physical characteristics of the site does not limit the applicability of seismic methods, and data acquisition and interpretation methods are optimized with respect to the problem specifications.

The time-lapse seismic map (Figure 10a) shows larger increases in time shifts in regions where higher amounts of CO<sub>2</sub> have been injected (Region A in the map). In fact, the greater the volume on injected CO<sub>2</sub> injection, the more pronounced the variation of the acoustic properties should be. According to Figure 10b, the amount of CO<sub>2</sub> injected in this region is less than the average amount in Region A. Hence, fluid properties in this region do not change as much as in Region A, resulting in a lighter alteration in acoustic properties in this region. As a result, although the time-lapse maps show a slight alteration of the acoustic properties in this particular region, the variations are not as intensified as Region A. It is important to add that there have been different interpretations of the same seismic data from Cranfield (Alfi and Hosseini, 2016). The main difference between the current interpretation and the ones done by other authors is that the previous interpretations appear to underestimate the changes in the region around the sealing fault (Region B). In other words, although the volume of injected CO<sub>2</sub> in Region B is less than the volume injected into Region A, the current seismic interpretations reveal a detectable change in seismic responses from Region B. It worth to mention that the previous efforts failed to identify any significant changes occurring in Region B. The next section discusses some of the possible reasons for a weak signal in the structural trap east of the fault (Region B).

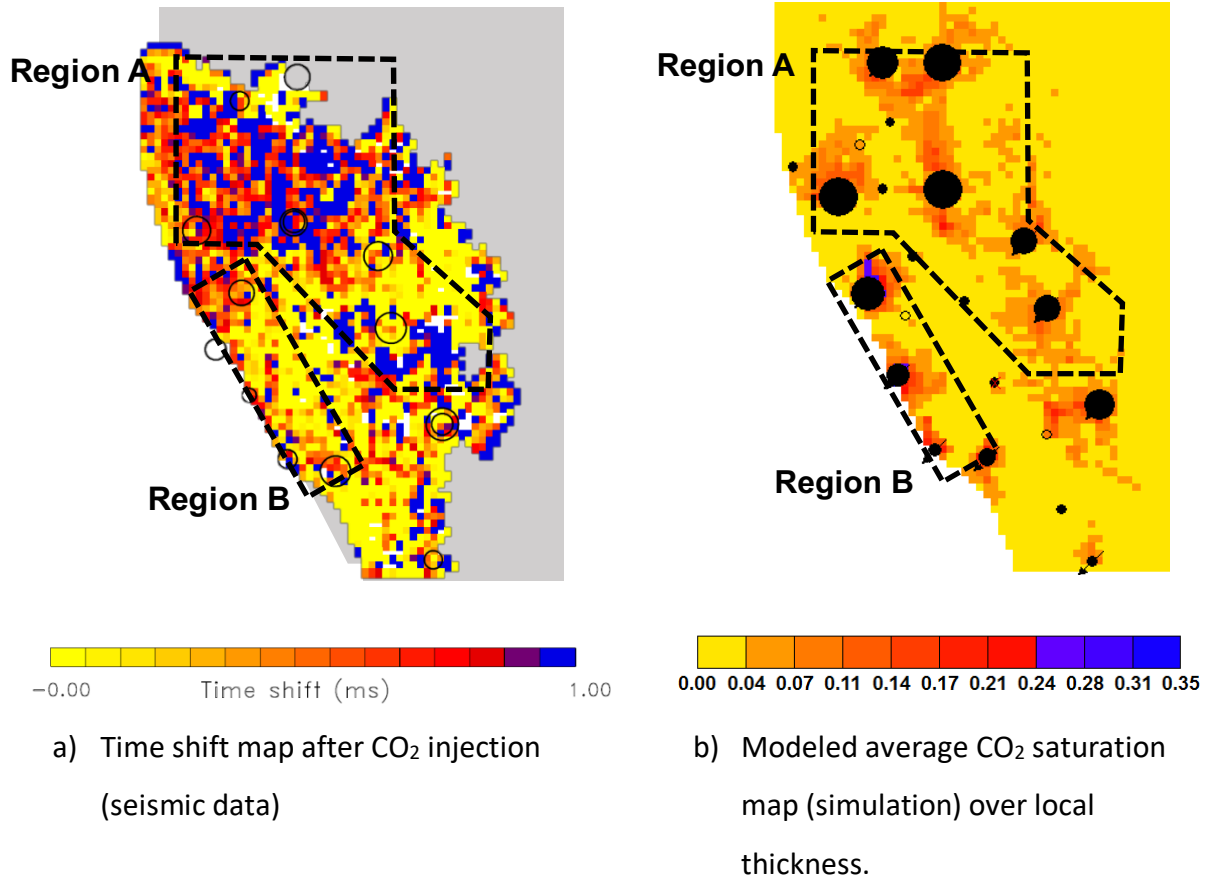


Figure 10-Seismic time shifts compared to the simulated CO<sub>2</sub> saturation distribution (both maps show the study area on September 2010). Time shifts are more pronounced in regions with a higher CO<sub>2</sub> saturation. The simulated saturations are averaged over the local thickness of the reservoir. The circles show the location of wells where the larger circles indicate higher amount of injected CO<sub>2</sub>.

### 3. The curious case of Cranfield

In this section, we provide a heuristic discussion of the Cranfield case, beyond the simple observation that areas with higher volumes of injected CO<sub>2</sub> should be detected in the seismic surveys. A weaker seismic time-lapse signal in Region B is not only attributed to less total injected CO<sub>2</sub> in that region, but also to the actual physics of the problem. Factors that we consider here include zone thickness and pre-injection fluid composition. In cases where the injection zone is thin, lower resolution seismic surveys may not be able to detect the changes properly and any possible signal change will be diluted by intact overlying/underlying layers with no injected CO<sub>2</sub>. In other words, a signal from the thin CO<sub>2</sub> invaded zone will be mixed with the weaker response of the nearby layers and the composite signal does not represent the actual change in elastic

properties of the porous media induced by CO<sub>2</sub> injection. For the case of Cranfield, the thickness of the CO<sub>2</sub> plume is somewhat uncertain because of non –ideal sweep efficiency at this early stage of injection in the heterogeneous reservoir. However, the operator placed several injection wells within the structural trap produced by the sealing fault, in expectation of a thick oil column target below a small gas cap (Figure 3). Figure 11 shows the productive thickness of the oil zone in Basal Tuscaloosa Sand, Cranfield unit. As we can see, the reservoir thickness in Region B, where seismic signals were relatively weaker is comparable to that of the rest of the oil zone.

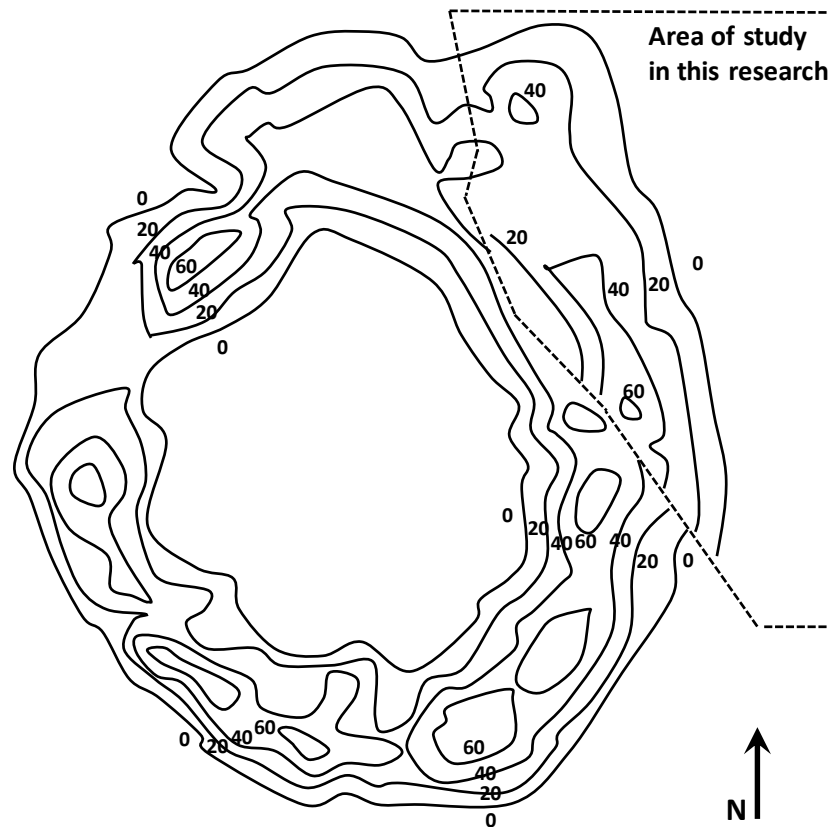


Figure 11-The productive zone thickness map of the study area (denoted by dashed polygon) shows the pay zone in the area of interest, located close to the sealing fault, which has a relatively high thickness compared to the other regions of the oil zone (contour map regenerated after Weaver and Anderson (1966)).

Seismic signals may also be influenced by the presence of gas, specifically methane, in the formation. It is informative to consider the gas saturation distribution prior to the first seismic survey at 2007 (Figure 3a). Given the prior production history, the most effective way to evaluate this is through a compositional reservoir simulation. According to the map of modeled gas

303 saturations, the gas that has evolved in the reservoir during the primary production period (1944-  
 304 1964) will move up-dip towards the center of the reservoir and will accumulate near the trap to  
 305 the east of the sealing fault. The presence of even small gas concentrations during the first  
 306 seismic survey can cause masking effects in which the alteration in acoustic properties of the  
 307 reservoir are not large enough to be detectable during the repeat seismic survey, particularly for  
 308 a case with a relatively high noise ratio.

309 To understand the reasoning behind such behavior, we turn to rock physics modeling and fluid  
 310 substitution theory (Mavko et al. 2009). Fluid substitution uses seismic velocities in rocks  
 311 saturated with one fluid to predict cases where the rock is saturated with a second fluid (here  
 312 CO<sub>2</sub>), or equivalently predicting saturated rock acoustic velocities from that of dry rocks. Fluid  
 313 substitution provides the interpreter with a tool for modeling and quantifying the different fluid  
 314 replacement scenarios, which might give rise to the observed amplitude and travel time variation  
 315 in the time-lapse responses. The most commonly used approach for fluid substitution  
 316 calculations is the low-frequency approach known as Gassmann's fluid substitution (Gassmann  
 317 1951). To model the changes from substituting one fluid with another we should first remove the  
 318 effects of the initial fluid and replace it with the new fluid. Gassmann's primary equation  
 319 correlates the saturated bulk modulus of the rock to its porosity and the bulk moduli of the  
 320 porous rock frame, mineral matrix, and the pore filling fluids (Smith et al., 2003):

$$K_{\text{sat}} = K^* + \frac{\left(1 - \frac{K^*}{K_0}\right)^2}{\frac{\phi}{K_{\text{fl}}} + \frac{1 - \phi}{K_0} - \frac{K^*}{K_0^2}} \quad (1)$$

321  
 322 where  $K_{\text{sat}}$  represents the bulk modulus of the porous media saturated with the fluid of interest,  
 323  $K^*$  is the bulk modulus of the porous rock frame (with no fluid inside the pores),  $K_0$  is the bulk  
 324 modulus of the mineral matrix (depends on the minerology of the rock components),  $K_{\text{fl}}$  is the  
 325 pore-filling fluid's bulk modulus, and  $\phi$  is porosity. In this formula, the value of  $K^*$  does not  
 326 depend on the fluid properties. To use this equation, the rock is drained of the initial fluid to  
 327 determine the bulk and shear moduli of the porous frame ( $K^*$ ). The porous frame properties are

then used to calculate the new effective bulk modulus of the media saturated with the new fluid ( $K_{\text{sat}}$ ). Gassmann equation assumes a connected pore space with equilibrated pressure along the pore network.

When the fluid saturation in a porous media varies, for example as  $\text{CO}_2$  replaces oil or water in the reservoir, the fluids can distribute within the formation in various configurations and this can lead to different bulk moduli ( $K_{\text{sat}}$ ) for the fluids. For example, in the case of a uniform or homogeneous distribution of fluids, the new value of  $K_{\text{fl}}$  is computed using the Reuss average (Smith et al., 2003):

$$K_{\text{fl}} = \left[ \sum \frac{S_i}{K_i} \right]^{-1} \quad (2)$$

where  $S_i$  represents the saturation of each individual fluid while  $K_i$  is the bulk modulus of that fluid phase. Eq. 2 is the effective modulus for a stack of layers perpendicular to the direction of propagation of the seismic wave. The Reuss average is also a lower bound on the effective fluid bulk moduli because the weakest layer determines the modulus. Note that the assumption of a homogeneous fluid distribution or of fluids arranged in layers perpendicular to the direction of propagation is not always a good approximation. While we cannot consider all of the possible fluid arrangements, we can consider the other extreme distribution that gives an upper bound on the effective fluid bulk modulus. This modulus, known as the Voigt bound, is given by

$$K_{\text{fl}} = \sum S_i K_i \quad (3)$$

and corresponds to the case in which the fluid is arranged in layers oriented parallel to the direction of propagation (Mavko et al., 2009). In that case the strongest layers determine the effective modulus. Given that one typically does not know the distribution of fluid in a heterogeneous reservoir with an extensive history of production and injection, we cannot be certain about the averaging that we should use. However, in making estimates of seismic velocity changes due to fluid saturation one could compare both the Reuss and Voigt approaches in order

to gauge the possible variation. The Hill estimate, which is simply the average of the Reuss and Voigt bounds, represents a reasonable approximation to the fluid bulk modulus in the face of uncertainty regarding the nature of the saturation distribution within a given reservoir (Mavko et al. 2009).

Changing  $K_{fl}$  would ultimately affect the bulk modulus of the saturated porous media ( $K_{sat}$ ). When  $K_{sat}$  changes, the seismic velocities in the reservoir would change and such changes can be detected in the time-lapse surveys and are interpreted as variation in fluid content of the reservoir. Eq. 4 shows how p-wave velocities are related to the saturated bulk modules of the porous media:

$$V_p = \sqrt{\frac{K_{sat} + \frac{4}{3}G}{\rho_B}} \quad (4)$$

In this equation,  $V_p$  is the p-wave velocity of the saturated media,  $G$  is the shear modulus of the saturated media and  $\rho_B$  is the bulk density (fluid and rock). Note that the shear modulus ( $G$ ) does not depend on the fluid content and will not vary with reservoir fluid alteration. Gassmann's fluid substitution will affect the value of  $V_p$  by changing  $K_{sat}$  and  $\rho_B$ . In Figure 12 we plot the seismic velocity variations corresponding to Reuss and Voigt approaches to averaging the fluid moduli. In addition, we also plot the Hill estimate, which is just the arithmetic average of the two bounds.

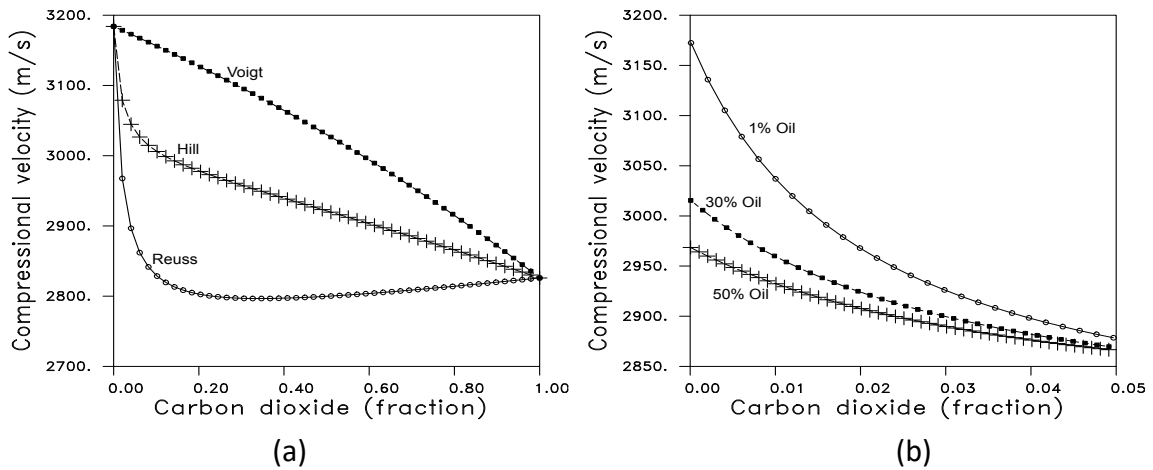


Figure 12-a) Voigt, Reuss, and Hill based estimates of compressional velocity changes in a porous layer saturated with varying fractions of carbon dioxide and water. b) Effect of varying

oil saturation as a third contribution to the fluid mixture for the Reuss average. The horizontal scale indicates the fraction of carbon dioxide trading off with water content.

To understand the importance of the presence of residual gas and its effect on seismic velocity, and to mimic the situation at Cranfield, consider a simple model in which the reservoir has a porosity of 0.24 with the mineral frame primarily composed of quartz. Ditkof (2013) reported an average P- and S-wave velocities of 3328 and 2096 m/s for a 100% brine saturated rock in this deposition environment (S- and P-velocities are used to calculate the bulk modulus of the porous rock frame,  $K^*$ ). According to Mavko et al. (2009), quartz has a bulk modulus of 36 GPa and density of 2.65 g/cc. Now, consider a case in which the reservoir is initially saturated with 40% water and 60% oil and CO<sub>2</sub> is injected into the reservoir to replace oil. **Error! Reference source not found.** uses the Gassmann formula and Reuss average to calculate the p-wave velocities for different injected CO<sub>2</sub> saturations. According to the graph, 30% increase in CO<sub>2</sub> saturation (from 0 to 30) will cause P-velocity to decrease from 3170 to 3010 m/s, which is 160 m/s reduction in the  $V_p$ . Now consider a case in which the reservoir is originally saturated with 40% water, 50% oil, and 10% residual gas. To make the case easier to understand and to be able to still use **Error! Reference source not found.**, we will assume the residual gas in the reservoir (usually methane or low-end light hydrocarbons) have properties comparable to that of CO<sub>2</sub> so we can still use the P-wave velocities in **Error! Reference source not found.**. For this case, the initial P-wave velocity (before any CO<sub>2</sub> is injected) should be read at the gas saturations of 10%, due to the presence of residual gas. According to the graph, the initial  $V_p$  is around 3040 m/s. Now assume we inject CO<sub>2</sub> in the reservoir and add 30% to the gas saturation in the reservoir. The repeat seismic survey in the area of the CO<sub>2</sub> injection will read a P-wave velocity of 3005 m/s (at CO<sub>2</sub> saturation of 40%). This means, 30% increase in CO<sub>2</sub> saturation will decrease the P-wave velocity by only 35 m/s. When we compare this value to the original reduction of 135 m/s for cases with no residual gas in the reservoir, one can conclude that the variation in elastic properties can be masked when residual gas is present in the reservoir. This is important because seismic surveys always contain random and signal-generated noise. If the level of expected changes in seismic responses is comparable to the noise level, such changes could be completely obscured. A similar masking effect is reported by Urosevic et al. (2011) when injecting CO<sub>2</sub> into a depleted gas reservoir in the

394 Otway Basin, Australia. Their results showed that presence of residual gas impacts changes in  
 395 elasticity of the reservoir rock after CO<sub>2</sub> injection in which even carefully implemented time lapse  
 396 seismic methodologies are not capable of capturing the changes.

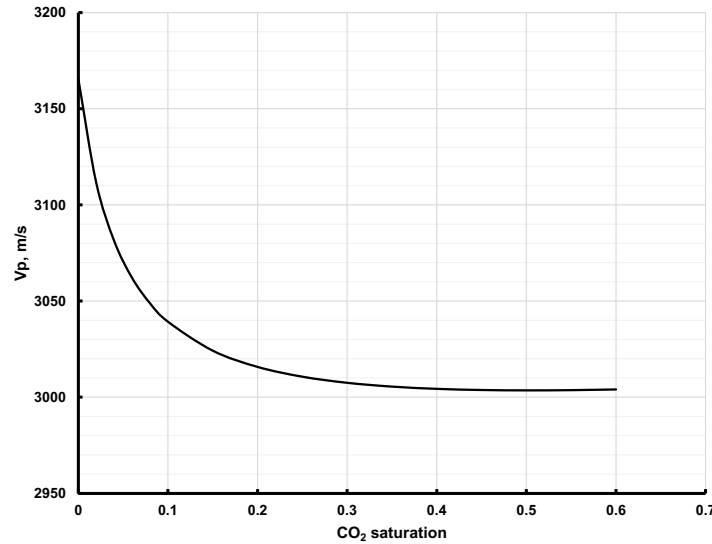


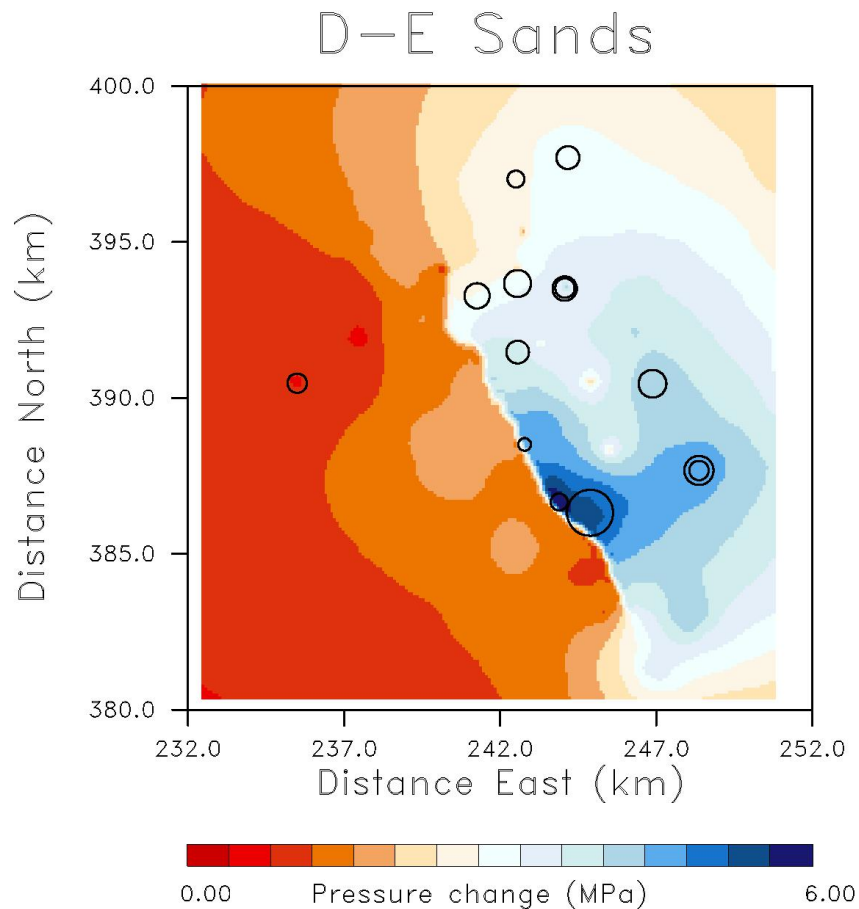
Figure 13-P-wave velocity in the saturated media decreases as CO<sub>2</sub> (fluid with a lower bulk modulus) saturation increases. Seismic wave velocity decline is steeper for lower saturations and slows down as CO<sub>2</sub> saturation increases, according to the Gassmann formulation.

397 Pressure changes in the reservoir, due to injection and production can also lead to changes in  
 398 seismic velocity. This may be due to changes in the bulk moduli of the rock matrix induced by  
 399 variations in effective pressure, or by changes in the nature of the fluid. For example, if the fluid  
 400 pressure is lowered below the bubble pressure gas may ex-solve from the solution. Recently, the  
 401 velocity-stress sensitivity at Cranfield has been characterized using crosswell seismic data  
 402 (Marchesini et al. 2017). This continuous active-source monitoring experiment during fluid  
 403 withdrawal indicated that the fractional change in seismic velocity for a given change in fluid  
 404 pressure is given by

$$\frac{dV}{V} = \frac{dT}{dT} = 1.09 \times 10^{-3} / \text{MPa}$$



406 In order to calculate the influence of pressure changes on the seismic travel time and velocity  
 407 changes we considered the pressure variations calculated by the numerical reservoir simulation.  
 408 The estimated pressure changes between the two seismic surveys in 2007 and 2010 are shown  
 409 in Figure 14.



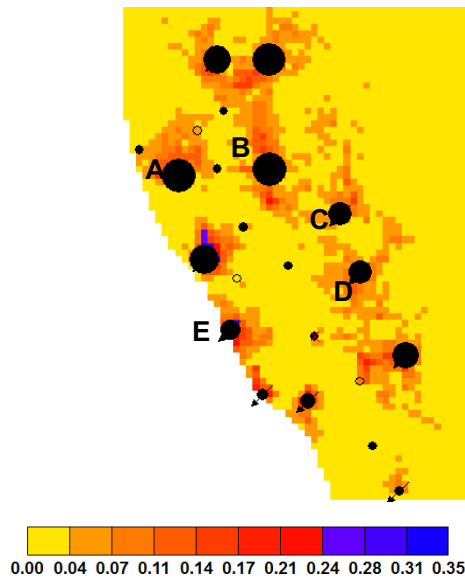
410 Figure 14. Pressure changes between the seismic surveys in 2007 and 2010, calculated using the  
 411 compositional reservoir simulation.

412 Given the pressure changes within the reservoir, we estimate that the change in two-way travel  
 413 time through the reservoir will be less than 0.1 ms and that the change in velocity will be less  
 414 than 1%. As is evident in Figure 14, this change will be largest in an area adjacent to the bounding  
 415 fault and much less outside of this region. Furthermore, it has been hypothesized that the change  
 416 is due to gas coming out of solution due to the pressure decrease during the pumping test. Thus,  
 417 given that the pressure increases within the reservoir during the injection of carbon dioxide, the

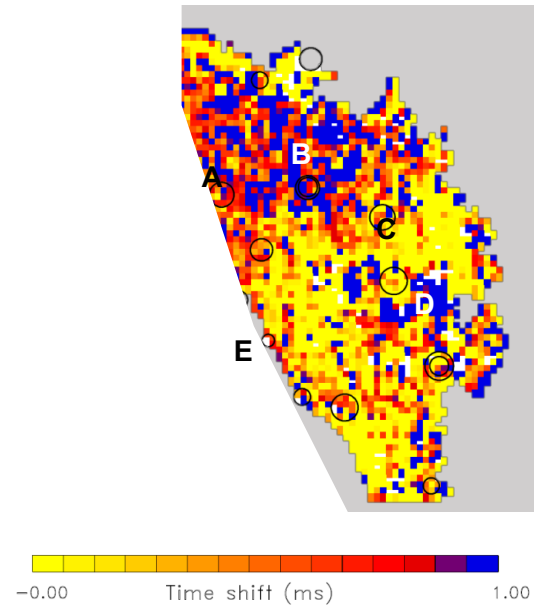
seismic changes due to such pressure variations might be much smaller. Given that the time shifts estimated from the time-lapse data are of the order of 1 ms, the pressure changes are not thought to significantly impact our results.

#### **4. Comparison of different seismic interpretations from Cranfield**

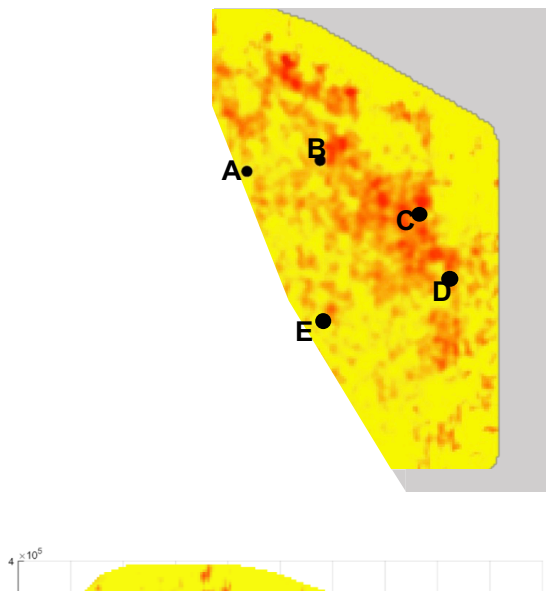
In this section, we compare the results of three different interpretations from the same seismic data acquired from Cranfield (Figure ). This comparison helps us to understand how the actual physics of the problem can influence our interpretation of the same seismic data. In Figure a, five injection wells are marked with letters A to E. The color map shows average CO<sub>2</sub> saturation at the time of the second seismic survey (2010) and the size of the black circles show the relative amount of CO<sub>2</sub> injected into the ground. In Figure 15, the letters are selected in a way that they are geographically consistent in all four graphs. If the seismic interpretation works as predicted, time-lapse seismic signals should show a more pronounced change in those injection locations. Figure b represents the seismic interpretations from the current work. Figure c shows the amplitude change map by flattening the reservoir to 2300 ms to characterize a 4D response (Ditkof et al., 2011; Ditkof, 2013). A relatively high amount of noise was reported in the seismic data, which can influence the repeatability of analysis according to the authors. Carter (2014) used the 3D seismic data along with well logs to probabilistically invert the CO<sub>2</sub> saturation and porosity distribution in Cranfield (Figure d). The well logs were used to calibrate the rock-physics model, which later was used to statistically derive the porosity distribution. A high resolution basis pursuit inversion was then used to invert seismic data for P-impedance and obtain the CO<sub>2</sub> saturation distribution (Carter, 2014).



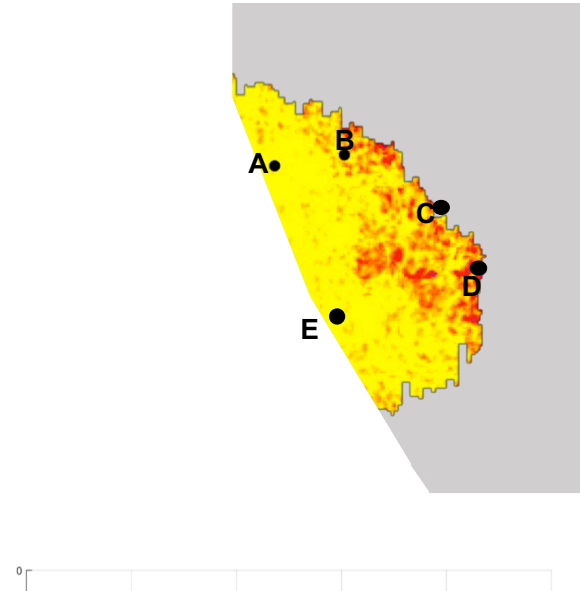
a) CO<sub>2</sub> saturation (simulation)



b) time shift map (current work)



c) amplitude change map (Ditkof et al., 2011)



d) CO<sub>2</sub> saturation from seismic (Carter, 2014)

Figure 15- Comparison of Cranfield seismic interpretations from three different studies with the actual field injection data and reservoir simulation results.

439 Let us consider point A where field data and simulation results show a considerable amount of  
 440 injected CO<sub>2</sub> with high CO<sub>2</sub> saturations around an injection well. The new seismic evaluation from  
 441 this paper (Figure b) shows a noticeable change in the time shift around point A, while the

interpretations by Ditkof et al. (2011) (Figure c) and Carter (2014) (Figure d) do not. This represents an example of an improvement in seismic interpretation that results in an improved match with the fluid flow model. CO<sub>2</sub> injection in points B and D is captured in all three seismic interpretations. These matches increase confidence in the correctness of the fluid flow model and the seismic interpretation. Both Ditkof et al. (2011) and Carter (2014) report no alteration in elastic properties of the media. We believe that correlates to the specific physics of the current problem, in which presence of residual gas migrated to the top of a structural trap adjacent to the sealing fault, which can mask the seismic signals in that area. However, some smaller scale mismatches remain; for example at point D, we can see a mismatch between the new seismic interpretation and the fluid flow model. The reservoir is highly heterogeneous and it is not possible to characterize it on a small enough scale to adequately capture this variability. The history matching of production data can capture the larger-scale variations in reservoir properties but will likely miss variability that is important for the detailed changes in seismic properties. In addition to the variability in static flow properties, there are variations in fluid saturations due to past production and injection. This will lead to spatial variations in the distribution of methane, water, and oil within the reservoir, variations in the thickness and geometry of the zone saturated by CO<sub>2</sub>, in addition to variations in both mechanical and flow properties. Therefore, we might only expect to obtain large-scale agreement between reservoir simulation predictions and observations of seismic time-lapse changes.

## **5. Forward model**

In an effort to capture some of the large-scale variability that is possible in the seismic time-shifts, we used the simulation estimates of total saturation methane, water, oil, and carbon dioxide in 2007 and 2010 and Gassmann's approach to estimate seismic velocity changes within the reservoir. A background velocity model was constructed from three well logs that contained compressional velocity sonic logs to construct an average model. The velocity changes were superimposed on this background model for each location within the seismic velocity cube to construct a series of one-dimensional vertical profiles. Each linear profile was subject to a variable scale change in order to match the location and thickness of the reservoir boundaries. An approximate reflection response for a vertically propagating wave impinging on a stack of elastic

layers was computed using the method of Kennett (Kennett, 1983, 1974). We considered the three methods discussed above for averaging the saturations and fluid bulk moduli in order to compute the effective fluid bulk modulus, namely the Reuss, Voigt, and Hill averaging techniques. The resulting time shifts computed from the synthetic seismograms are shown in Figure . It was found that the most significant effects were due to the locations of interaction of injected  $\text{CO}_2$  with the oil-water contact at the edge of the field. Large changes were found in two main areas: where  $\text{CO}_2$  replaced brine in the water leg and in the area of high  $\text{CO}_2$  accumulation in the structural trap at the sealing fault. The relative size of the time shifts varies, depending on the technique used to average the fluid moduli (Figure ). For the Reuss model the largest changes are in the water leg to the northeast. For the Voigt approach the most significant changes are near the bounding fault. If the Hill method is used then the two major anomalies, near the bounding fault and in the surrounding aquifer, are roughly equal in magnitude. Note that the Voigt model leads to time shifts that are about an order of magnitude lower than the observed values. As mentioned earlier, the Reuss model would appear to be more compatible with property distributions that are largely horizontal or sub-horizontal.

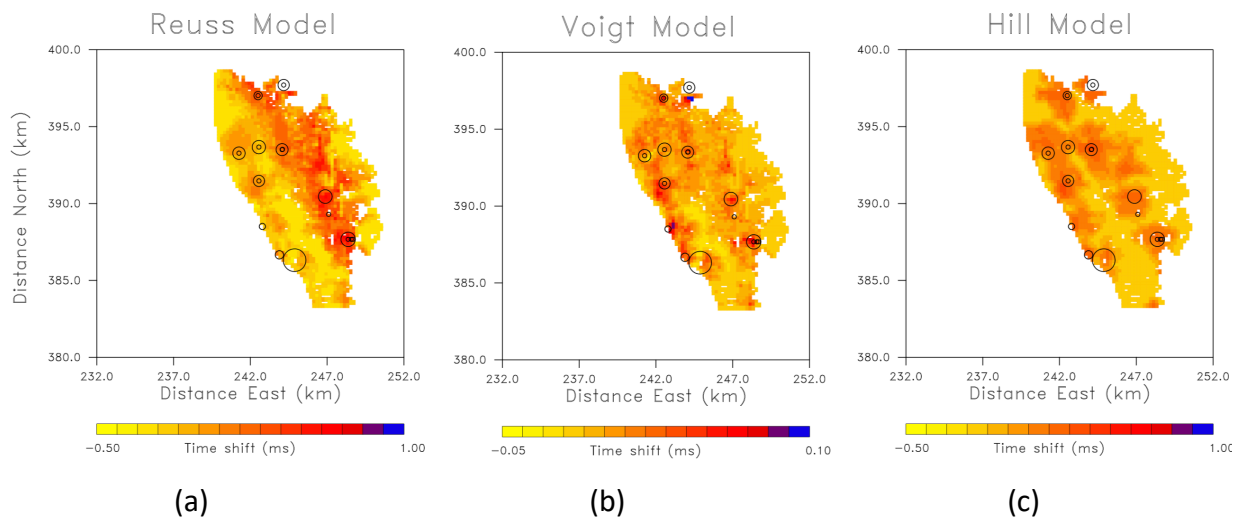


Figure 16-Time shifts calculated using the reservoir simulations and computed changes in methane, oil, water and carbon dioxide saturations. Three different methods, the Reuss bound (a), the Voigt bound (b), and the Hill average of the two (c), were used to compute the effective fluid bulk modulus at each location in the reservoir. Please note that time shift color scale is variable in different plots.

In general, the calculated time shifts (Figure ) are in qualitative agreement with the observed values (Figure 10 and Figure ) with large changes in the water leg of the reservoir and smaller but observable changes in the trap near the bounding fault. The detailed distribution of the anomalies will depend upon smaller scale variations in reservoir properties and will be influenced by heterogeneous static properties as well as variations in background fluid saturations due to earlier production and injection. It may be possible to use the good quality seismic data to infer these properties but that would be difficult with just a single pair of low resolution seismic surveys. In absence of economic limitations, a more robust approach would be to use a number of seismic surveys to image the propagation of the changes and to infer the onset of changes in the seismic properties due to the propagation of the carbon dioxide. Such onset times are more sensitive to flow properties and less sensitive to the rock physics models that govern the mapping from saturations to seismic moduli (Vasco et al., 2015, 2014). Multiple surveys were used at the Sleipner site (Arts et al., 2004) and they clearly documented the movement of the carbon dioxide in three dimensions.

## **Conclusions**

In this paper, we compared the results of reservoir simulation and seismic interpretation to consider a strategy for determining CO<sub>2</sub> movement during EOR operations. A compositional simulation approach is employed to address thermodynamics of CO<sub>2</sub>/hydrocarbon interactions and reliably model the miscibility phenomenon. Of a specific interest is gas exsolution/dissolution, which essentially leads to accurate modeling of gas appearance in the reservoir. The numerical model is history-matched using the available field production data. Seismic interpretations show large changes in the water leg of the reservoir and smaller but observable changes bordering the structural trap against the bounding fault. We obtain large-scale agreement between reservoir simulation predictions and observations of seismic time-lapse changes. In addition to the variability in static flow properties, there are variations in fluid saturations due to past production and injection. This will lead to spatial variations in the distribution of methane, water, and oil within the reservoir, in addition to variations in both mechanical and flow properties. For the specific case of Cranfield, one can conclude that the variation in elastic properties (before and after CO<sub>2</sub> injection) can be masked when residual gas

is present in the reservoir. This is important because seismic surveys often come with uncertainty and noise and if the level of expected changes in seismic responses is comparable to the noise ratios, such changes could be completely ignored or misunderstood.

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