

# Development and application of a coupled ECLIPSE – GeoSys modelling system for simulation of CO<sub>2</sub> storage in saline aquifers

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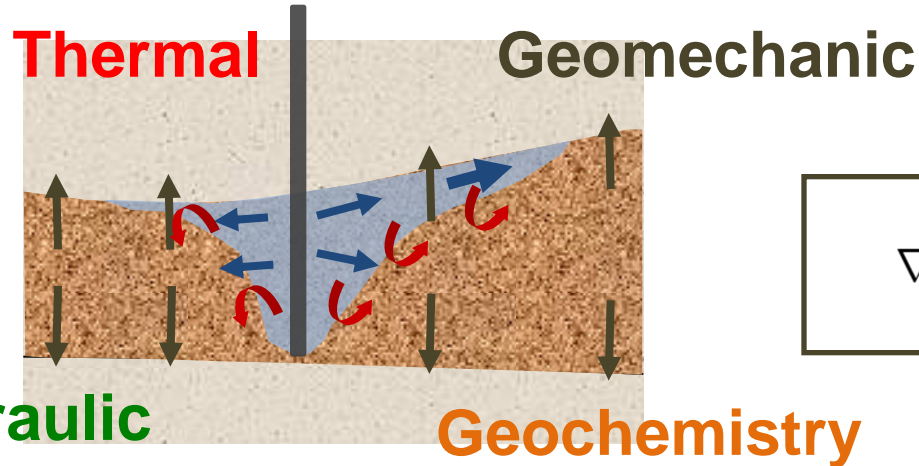
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Simulation of CO<sub>2</sub> injection → THMC modelling system

## Heat transport

$$\frac{\partial(uq)}{\partial t} + \nabla \cdot (uq\mathbf{v}) + \nabla \cdot (p\mathbf{v}) - \nabla \cdot (\lambda \nabla T)$$



Deformation

$$\nabla \vec{\sigma} - \rho \vec{g} = 0$$

## Fluid dynamics

$$\frac{\partial(\rho_\alpha n S_\alpha)}{\partial t} - \nabla \cdot \left( \rho_\alpha \frac{k_{r\alpha} k}{\mu_\alpha} (\nabla p_\alpha - \rho_\alpha g) \right) - q_\alpha \rho_\alpha = 0$$

$$\sum_{\alpha=1}^{\alpha=n_{phase}} S_\alpha = 1$$

$$p_{c\alpha\beta} = p_\alpha - p_\beta = f(S_1, \dots, S_{n_{phase}})$$

## Transport und Reactions

$$\frac{\partial(n S_\alpha \rho_\alpha X_\alpha^\kappa)}{\partial t} + \nabla \cdot (v_\alpha \rho_\alpha X_\alpha^\kappa - n S_\alpha D_{DD, \alpha}^\kappa \nabla(\rho_\alpha X_\alpha^\kappa))$$

$$x_\alpha^\kappa = H_\alpha^\kappa p_g^\kappa \quad - q_\alpha^\kappa = 0$$

$$K_m = \prod_i a_i^{\nu_i} = \exp\left(\frac{-\Delta G_0}{RT}\right)$$

$$r_m = k_{m, T_0} \left( \frac{-E_a}{R} \left( \frac{1}{T} - \frac{1}{T_0} \right) \right) A_m \left[ 1 - \frac{Q_m}{K_m} \right]$$

# OpenGeoSys (OGS)

= Scientific open-source code development for coupled **Thermo-Hydro-Mechanical-Chemical (THMC)** systems in the environment

- Object-oriented programming
- Using Finite-Element-Method

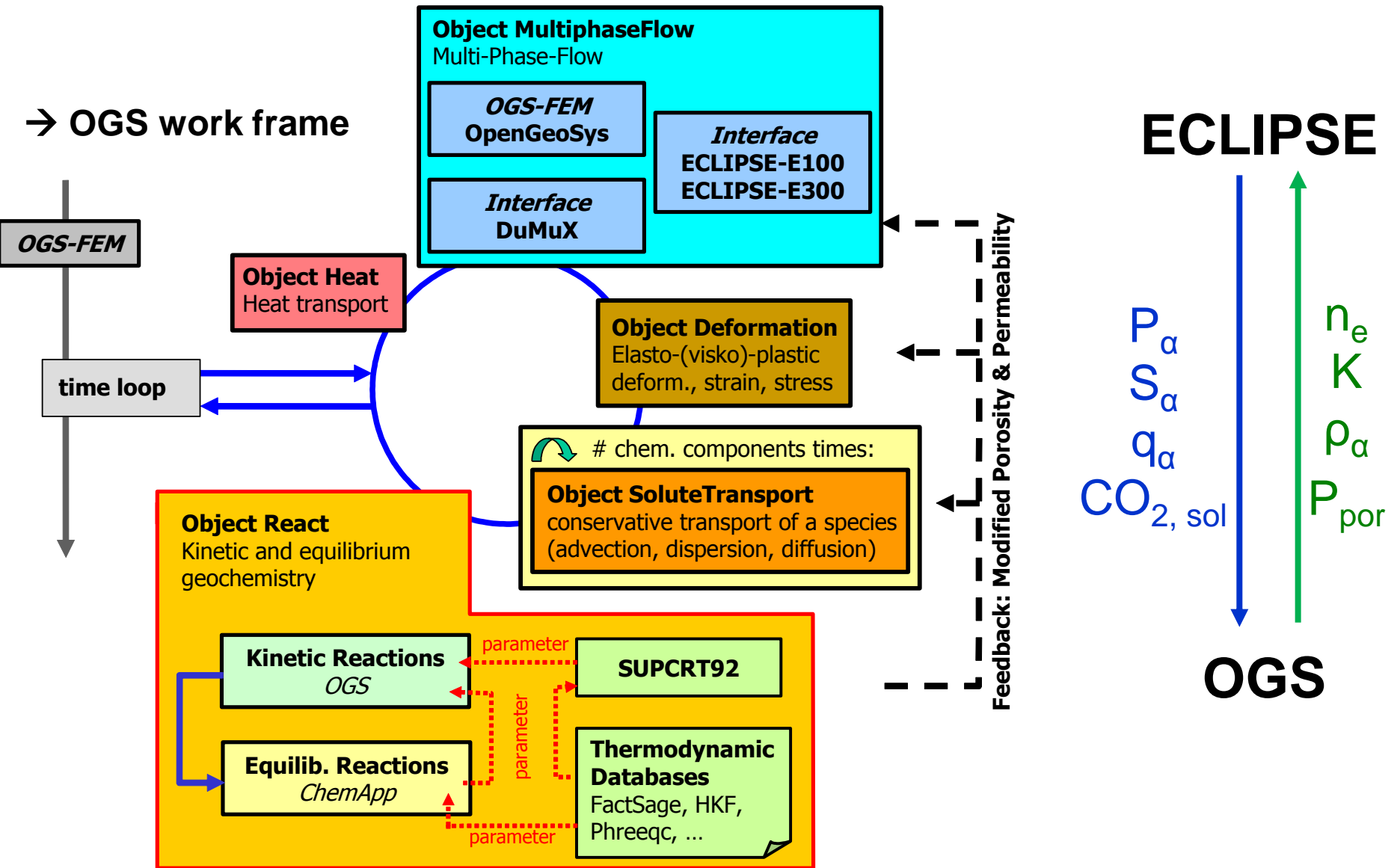
## Schlumberger © **ECLIPSE**

- widely used multiphase simulator
- Finite-Difference-Method
- CO<sub>2</sub>Store option

## Benefits of the OpenGeoSys-ECLIPSE coupling:

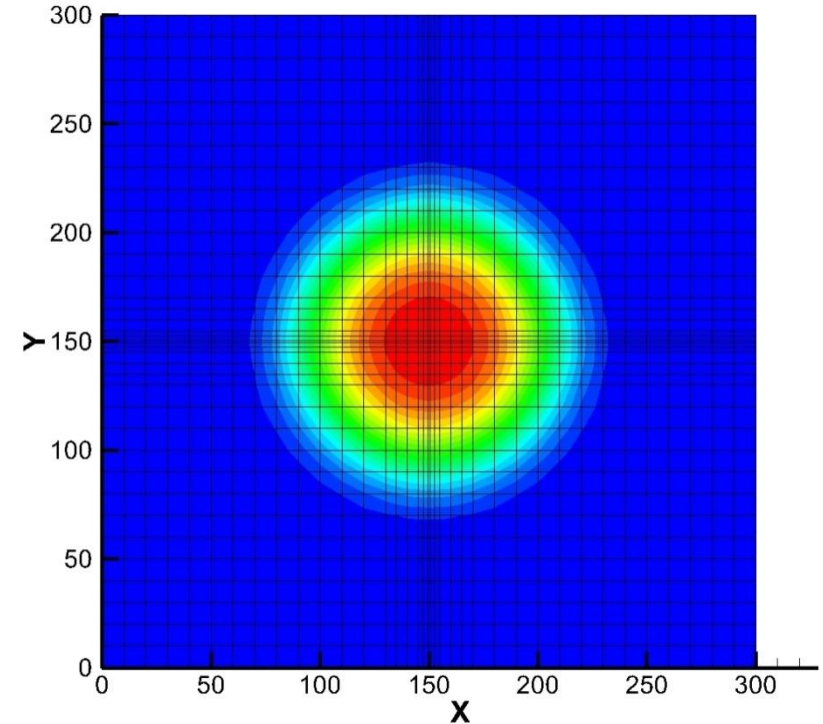
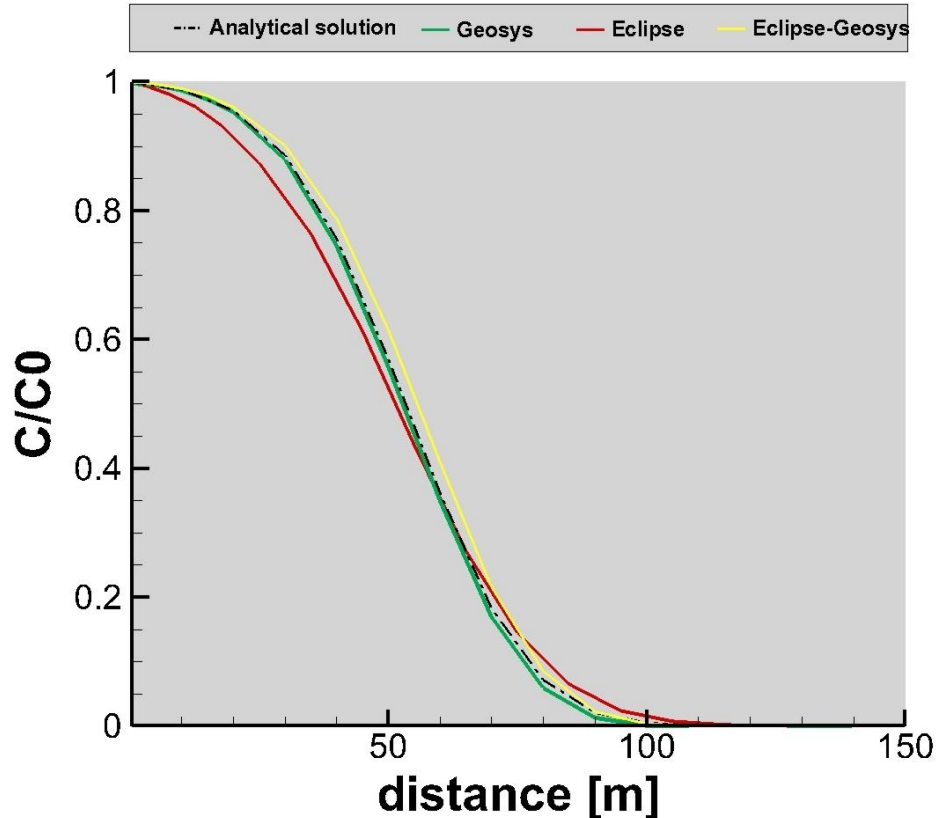
- benefit from the fast computation time of ECLIPSE
- Extend the capabilities of the ECLIPSE simulator for modelling CO<sub>2</sub> injection regarding Geochemistry and Geomechanics
- Using already existing fluid flow models build in ECLIPSE

# Implementation of the coupled code



## Benchmark 1:

### 2D Tracer Transport in a 1Phase radial flow field (after Moench and Ogata, 1981)

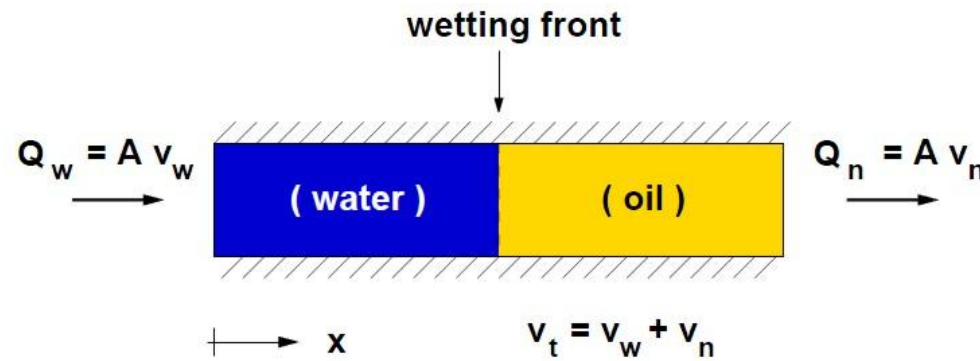


Good agreement of the coupled ECLIPSE-GeoSys code with the analytical solution.

Differences between the tracer fronts result from different mathematical methods which are used within the codes.

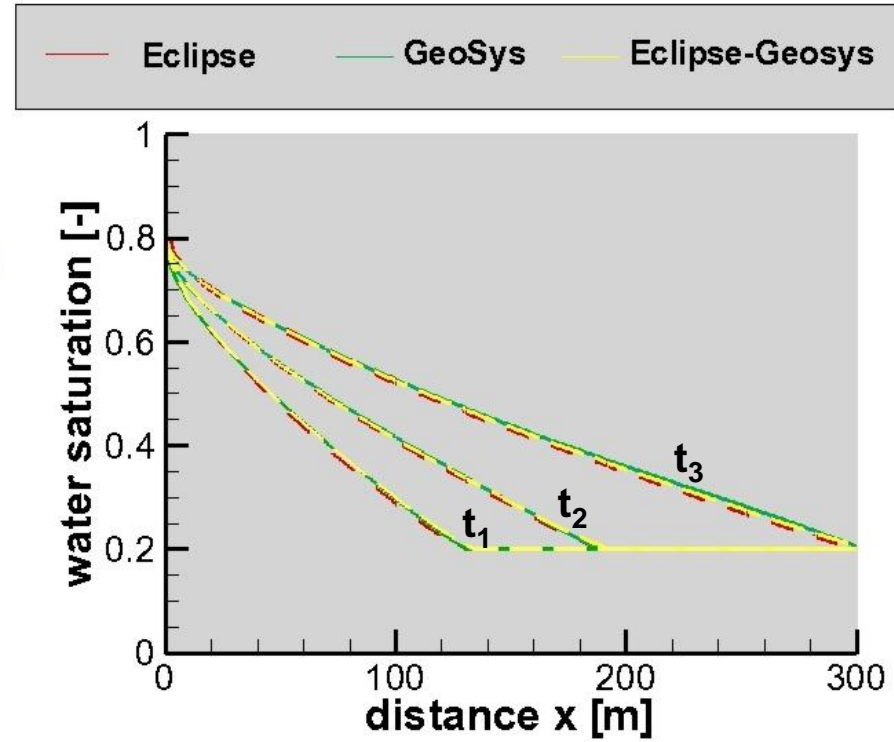
## Benchmark 2:

Advective-dominated two phase flow in a homogeneous system (after Buckley and Leverett, 1942)



(Paul, 2003)

- Linear  $k_r$ - $S_w$ -function and  $p_c$ - $S_w$ -function
- GeoSys: PP formulation
- ECLIPSE: PS formulation

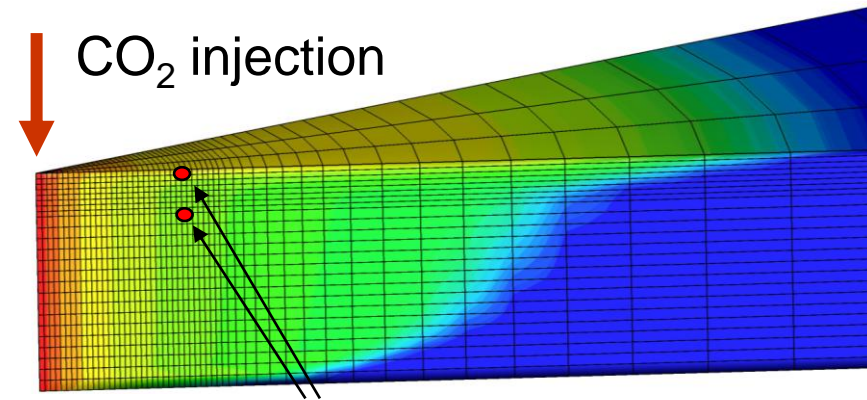




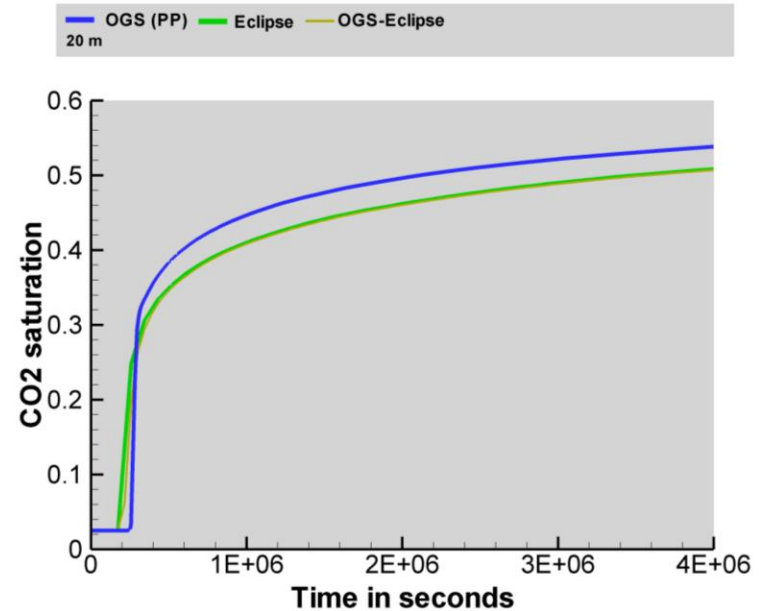
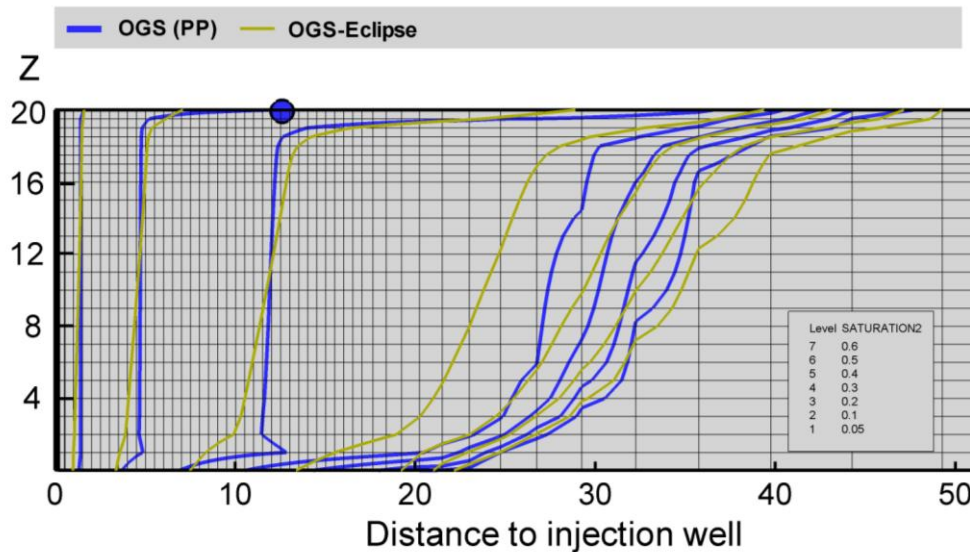
## Benchmark 3:

### 3D multiphase flow field with CO<sub>2</sub> injection

- Radial settings (injection left, boundary condition right)
- Adequate for model development and testing



Breakthrough



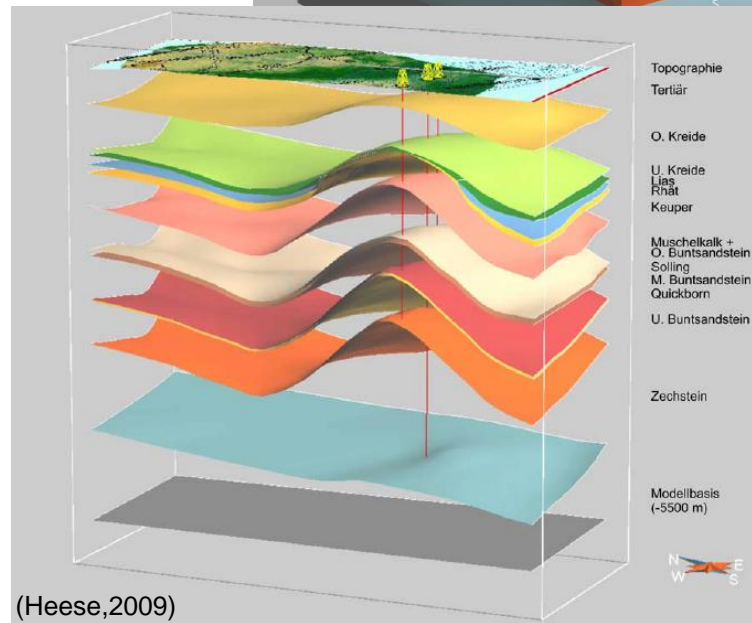
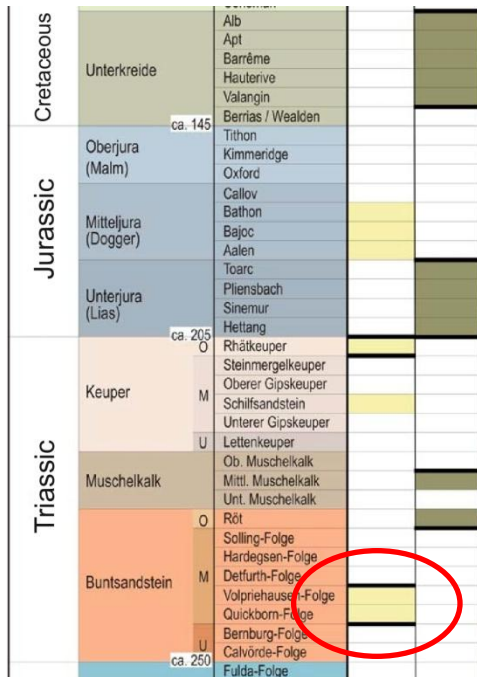
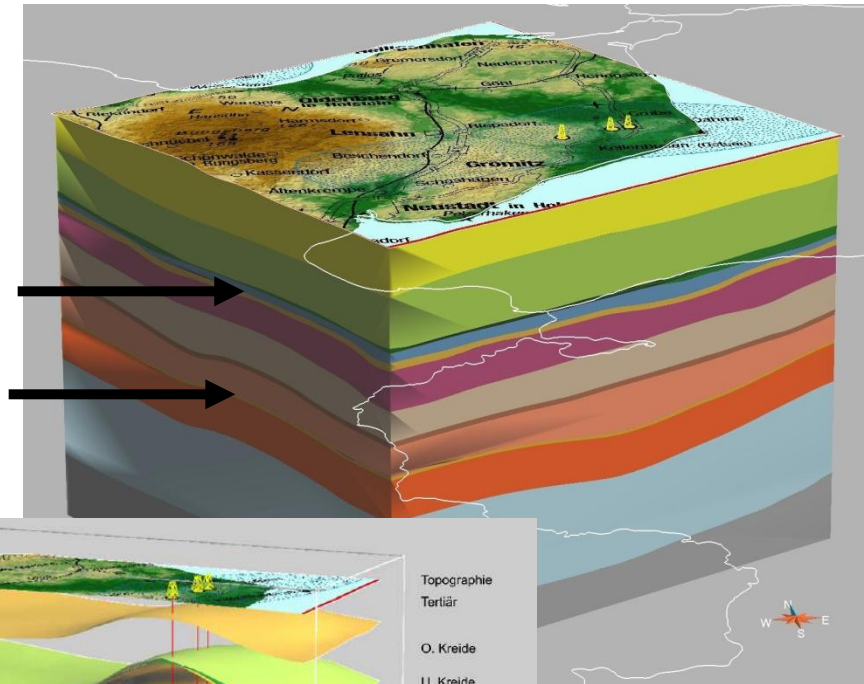
## Geological Model:

Extend: 29 x 28 km

Entire thickness ~ 5000 m

### Possible Storage Formations:

- 1) Rhät Sandstone
- 2) middle Bunter Sandstone



(Heese, 2009)



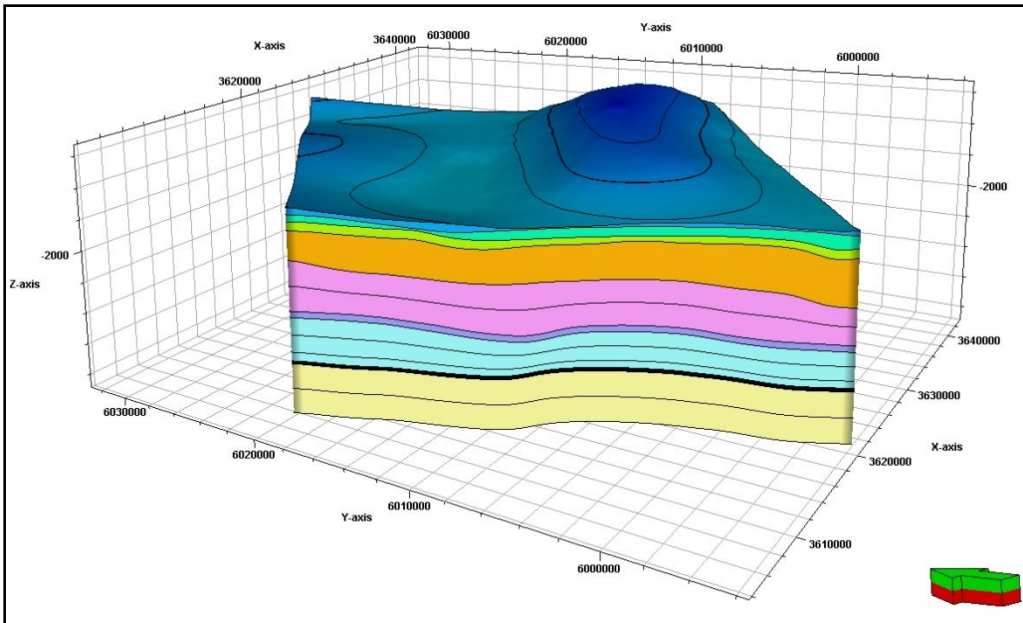
# Simulation model

Entire thickness ~ 1500 m

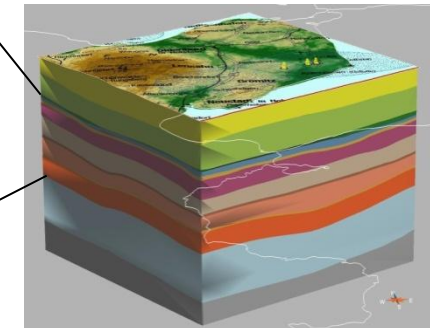
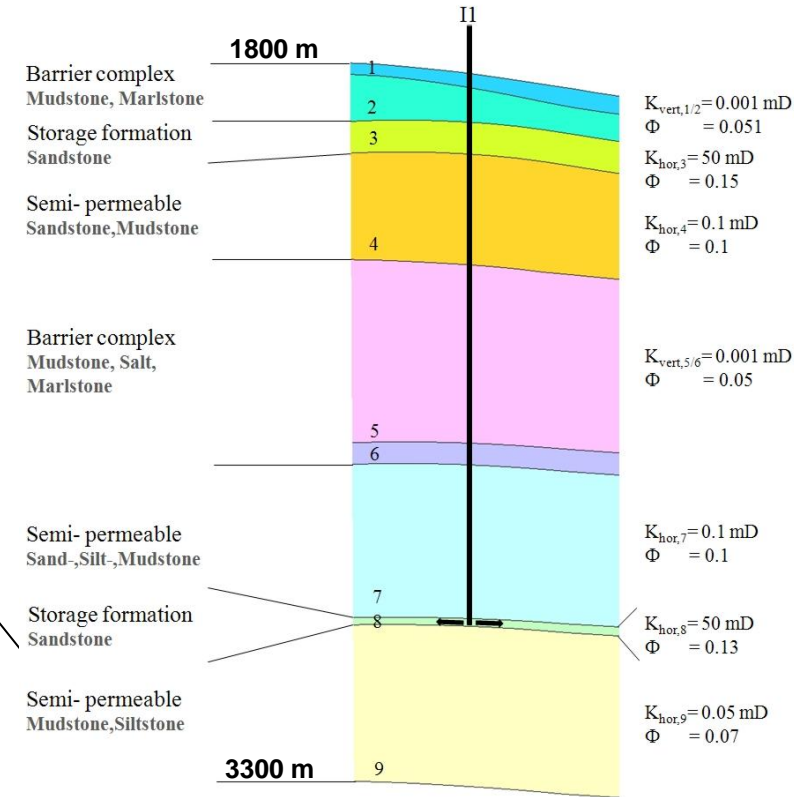
Discretization: 65 x 111 x 23 cells

## Szenario:

- CO<sub>2</sub> injection into the lower storage formation over a period of 20 years (1 Mio tons per year)
- 70 years post-injective simulation



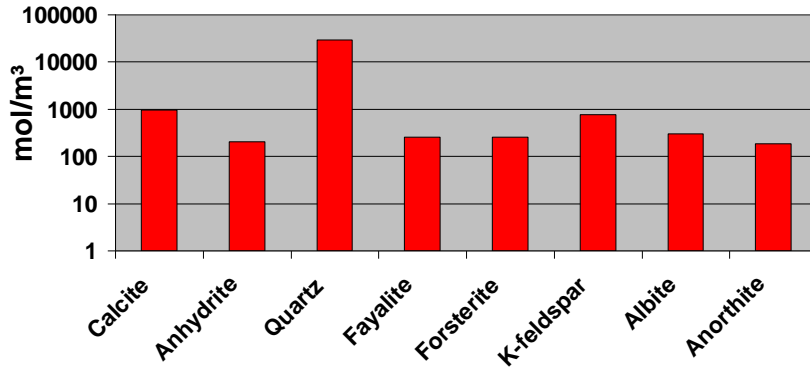
# Field site application



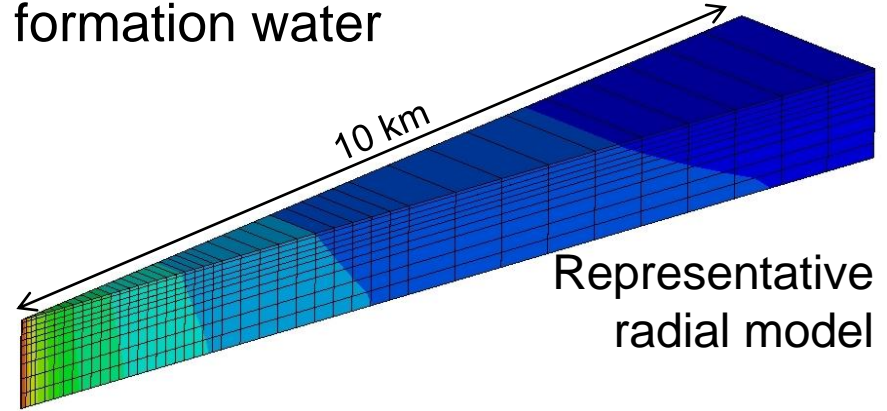
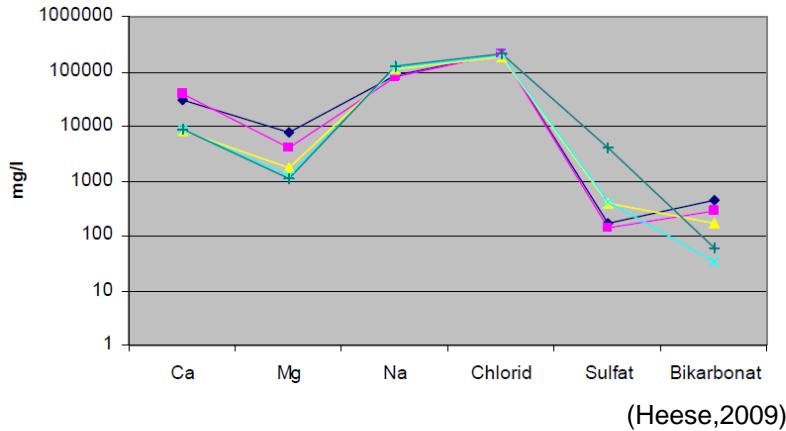
## Geochemical modelling:

Characterisation of the storage rock and formation water

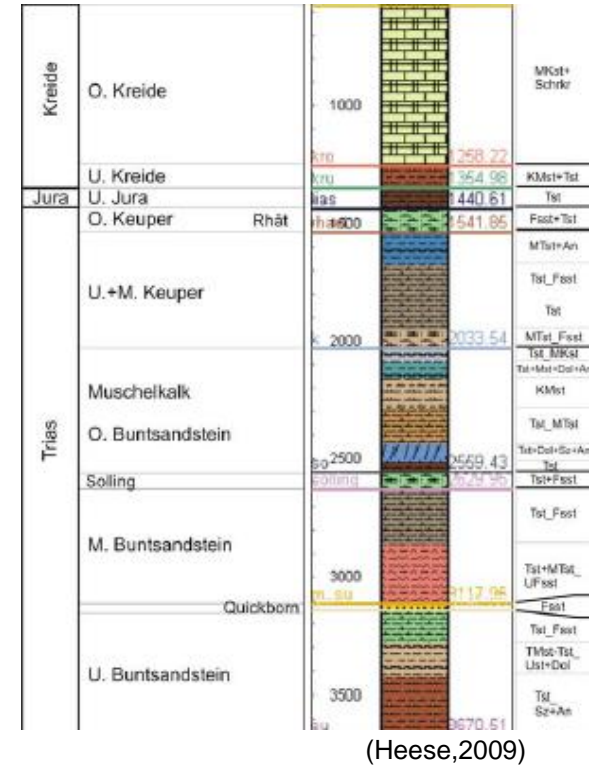
### Mineral composition



### Brine composition



### Lithology information

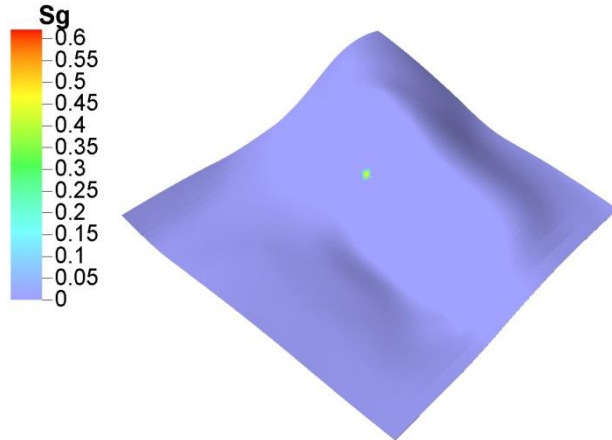


## Simulation results of the coupled code:

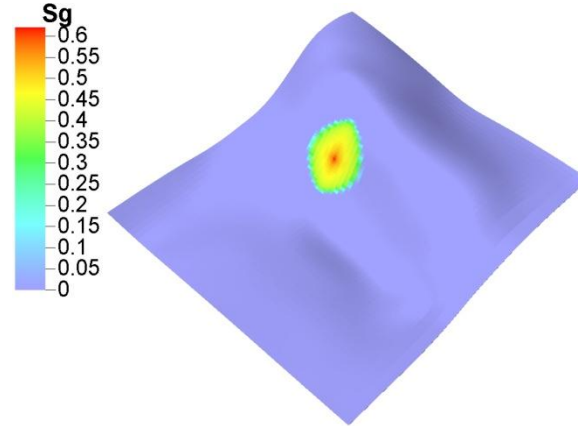
### 1) Pressure build up and phase propagation

Phase propagation after:

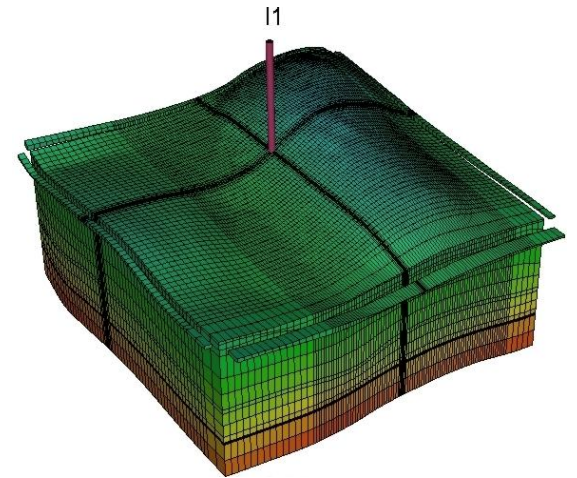
a) Injection start



b) Injection stop

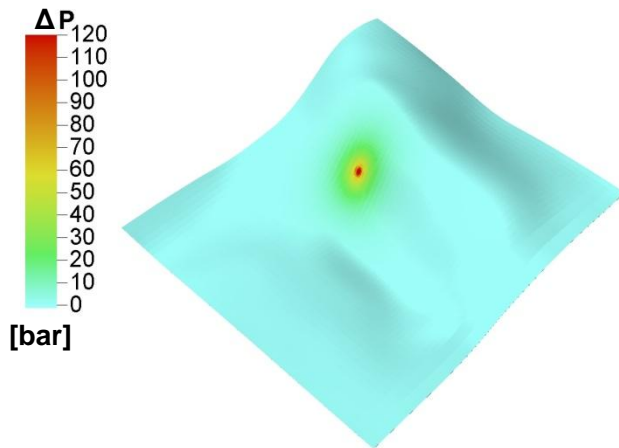


Simulation grid  
(4 times exaggerated)

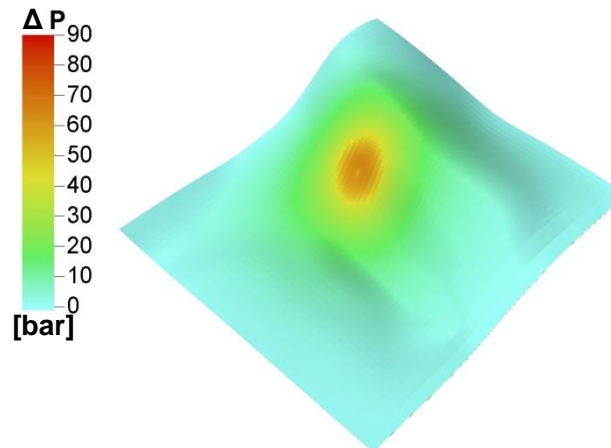


Pressure change after:

a) Injection start

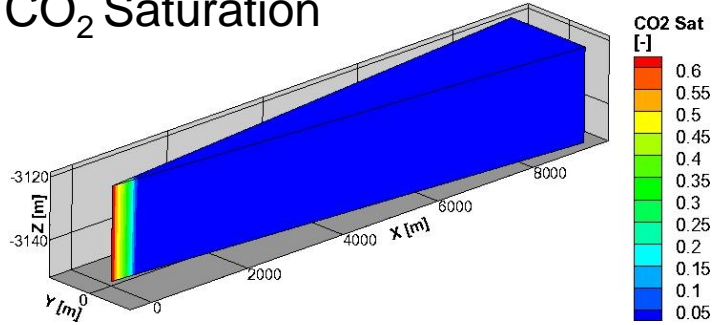


b) Injection stop

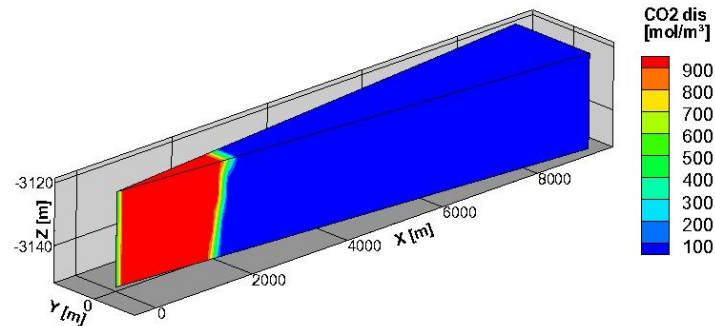
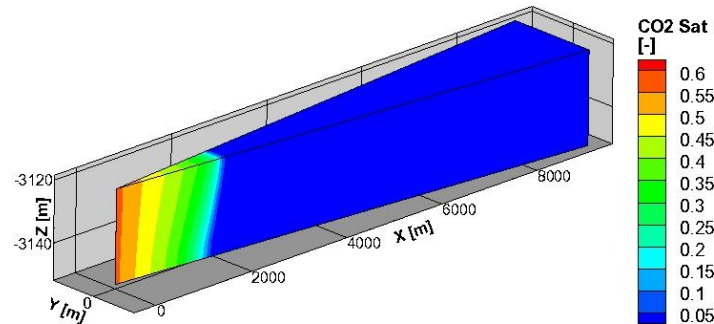
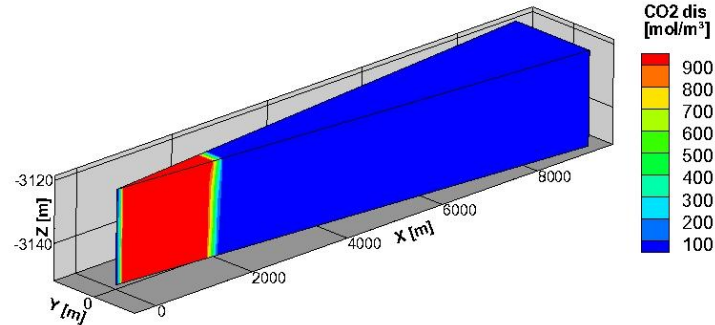
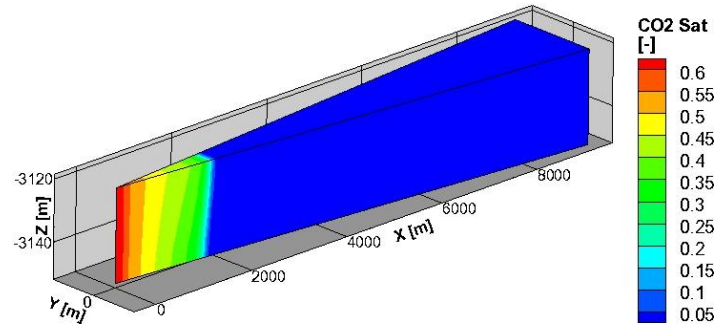
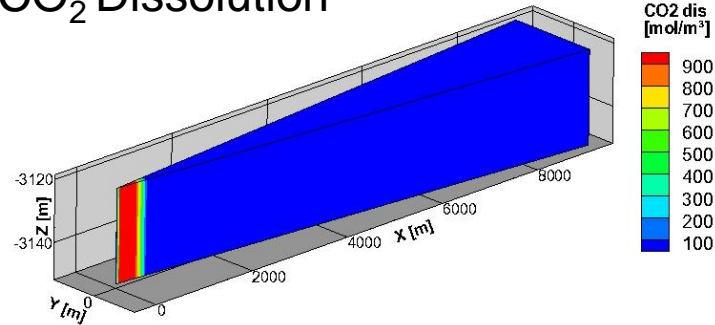


## Simulation results of the coupled code: 2) Geochemical modelling

### CO<sub>2</sub> Saturation



### CO<sub>2</sub> Dissolution



10 days

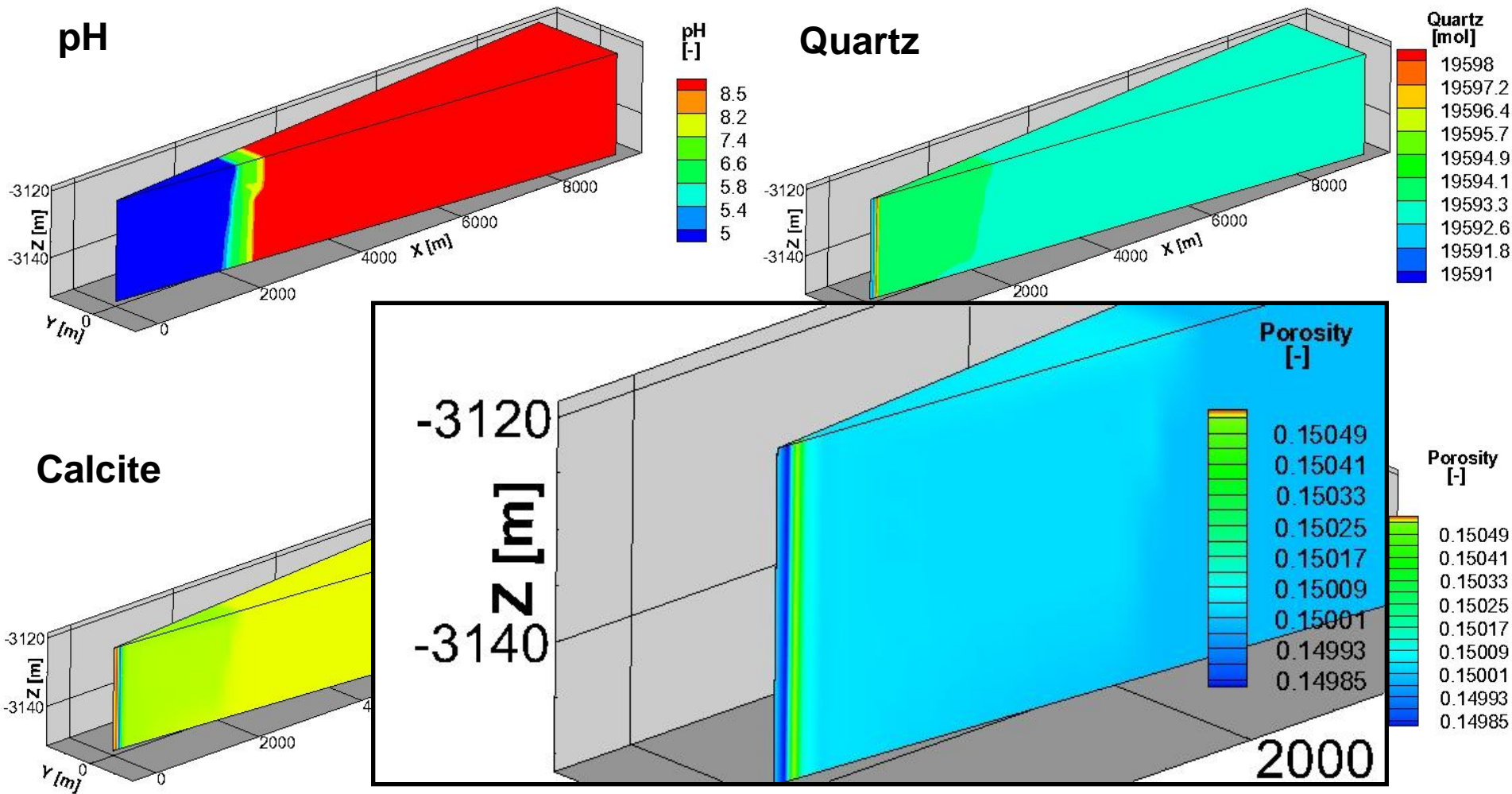
20 years

90 years



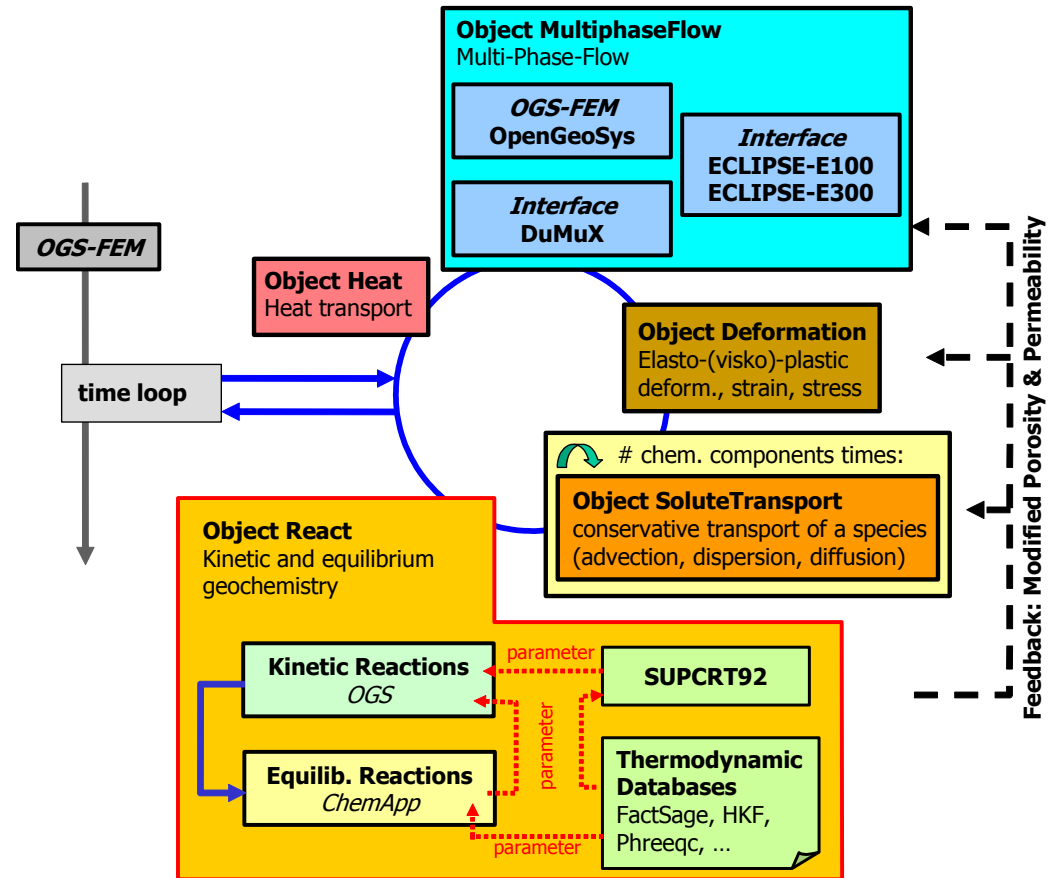


## Simulation results of the coupled code: 2) Geochemical modelling after 90 years



## Next steps

- Increasing computation efficiency
  - Multiphase flow
  - Geochemistry
- Mechanical feedback on multiphase flow (OGS→ECLIPSE)
- Further code verification
- More large scale applications
  - Geochemistry





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**THANKS FOR YOUR  
ATTENTION**

**Greetings from Kiel**