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# PetraSim Example Manual

## September 2008

# TOUGHREACT Example Guide

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# Chapter 1. Overview

## Using TOUGHREACT

TOUGHREACT is an extension of the original TOUGH2 simulation that is available as a simulator mode in PetraSim. The TOUGHREACT simulator supports a subset of the TOUGH2 EOS modules. The supported EOS moduels are EOS1, EOS2, EOS3, EOS9, and ECO2. You can perform a TOUGHREACT simulation by selecting the TOUGHREACT simulator mode and one of the available EOS modules in the PetraSim Preferences dialog.

In PetraSim, options relating to TOUGHREACT are presented under the Tough React menu item in the main window. These options allow you to configure the reactive transport solver, simulation output, chemical zones, and other TOUGHREACT-specific parameters.

It is also possible to disable reactive transport during a TOUGHREACT simulation. This will effectively revert the simulator to TOUGH2 mode. This option is available in the Global Properties dialog, on the Analysis tab.

## Input Files

A TOUGHREACT simulation requires four input files. These files are listed below:

- flow.inp -- This is the standard TOUGH2 input file
- solute.inp -- The chemical "geography" of the analysis
- chemical.inp -- The chemical parameters
- thermodb.txt -- The thermodynamic database

PetraSim creates each of these files into your simulation directory. However, the filenames cannot be changed. The naming scheme for the simulation output files follows a similar pattern. To avoid overwriting previous simulation input and output data, you must run each analysis in a separate directory.

## Thermodynamic Database

A thermodynamic database listing the composition of many different species and minerals has been included with PetraSim. PetraSim will automatically load this database. A valid database must be loaded prior to the inclusion of species or the definition of reactive zones. This is because the species used to build up the zones are loaded from the thermodynamic database. If choose to use a custom databse, you must ensure that it is loaded before configuring any species or zones. To load a custom thermodynamic database:

1. On the **Tough React** menu, click **Thermodynamic Database...**

2. Select your custom thermodynamic database
3. Click **OK**

# Chapter 2. Aqueous Transport with Adsorption and Decay (EOS9)

## Description

This problem is the first example in the TOUGHREACT manual. It is a 1-D problem, 12 m in length, with a unit area, divided into 60 blocks of 0.2 m thickness, as shown in Figure 2.1.

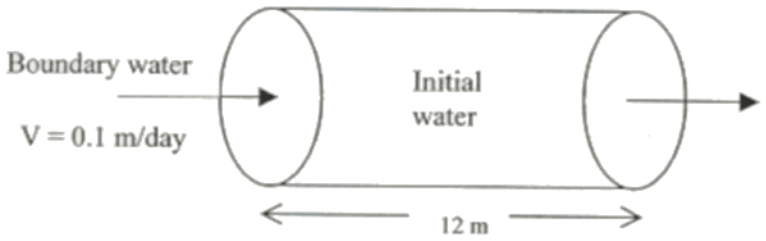


**Figure 2.1. Aqueous Transport with Adsorption and Decay Model (after [Xu, Sonnenthal, Spycher, and Pruess, 2004])**

## Specify the Equation of State (EOS)

To ensure that PetraSim uses TOUGHREACT and EOS9, you may need to edit your PetraSim preferences. To edit preferences, you can use the **Preferences** dialog.

1. On the **File** menu, click **Preferences...**
2. In the **Simulator Mode** list, select **TOUGHREACT**
3. In the **Default Equation of State (EOS)** list, select **EOS9**
4. Click **OK**

Your preferences have now been updated, but you will need to create a new model or restart PetraSim before they become active. If you already have a model open, it will not be affected by the preference change.

On the **File** menu, click **New**.

## Create the Model Boundary

To create the boundary for this model, use the **Define Model Boundary** dialog. The boundary parameters for this model are shown in Table 2.1.

**Table 2.1. Model Boundary Dimensions**

| Axis | Min (m) | Max (m) |
|---|---|---|
| x | 0.0 | 12.0 |
| y | 0.0 | 1.0 |
| z | 0.0 | 1.0 |

To create the model boundary:

1. On the **Model** menu, click **Define Boundary...**
2. Enter the values from Table 2.1
3. Click **OK** to create the model boundary

## Create the Grid

To create the gridas shown in Figure 2.2:

1. On the **Model** menu, click **Create Grid...**
2. In the **Division Method** box, select **Regular**
3. In the **X Cells** box, type `60`
4. In the **Y Cells** box, type `1`
5. In the **Z Cells** box, type `1`
6. Click **OK** to create the grid

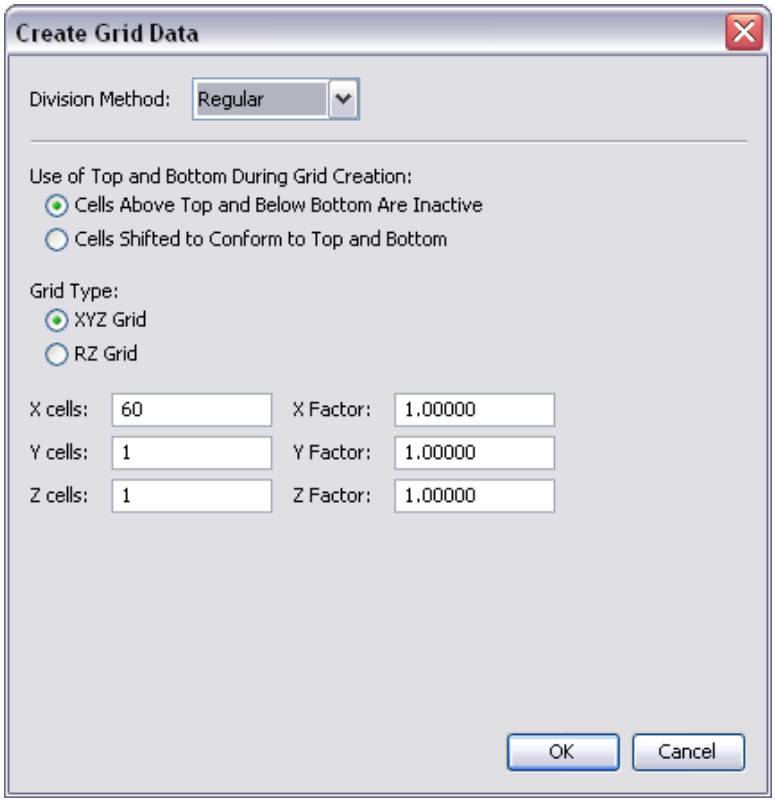


**Figure 2.2. The Create Grid dialog. The values shown will create a regular 60x1x1 grid**

## Global Properties

Global properties are properties that apply to the entire model. In this example, the only thing we will change is the analysis name. To edit global properties, you can use the **Global Properties** dialog.

On the **Properties** menu, click **Global Properties...**

### Simulation Name

1. In the **Global Properties** dialog, click the **Analysis** tab
2. In the **Name** box, type `TOUGHREACT Example 1`

### EOS Data

The EOS (Equation of State) tab displays options for EOS9.

1. In the **Global Properties** dialog, click the **EOS** tab
2. In the **Reference Pressure (Pa)** box, type `1.0E5`

3. In the **Reference Temperature (C)** box, type `4.0`

Click **OK** to close the **Global Properties** dialog.

## Material Properties

To specify the material properties, you use the **Material Data** dialog. This example requires one material.

To open the **Material Data** dialog: on the **Properties** menu, click **Materials...**

### Material Data

1. In the materials list, select **ROCK1**
2. In all three **Permeability** boxes (X, Y, and Z), type `6.51E-12`
3. In the **Wet Heat Conductivity** box, type `0.0`
4. In the **Specific Heat** box, type `952.9`
5. Click **Apply** to save the changes

In addition to the physical rock parameters, we also need to specify the relative permeability and capillary pressure functions for this material. These options can be found in the **Additional Material Data** dialog. To open this dialog, click the **Additional Material Data...** button.

### Relative Permeability

To specify the relative permeability function:

1. Click the **Relative Perm** tab
2. In the **Relative Permeability** list, select **Linear Functions**
3. In the **Slmin** box, type `0.333`
4. In the **Slmax** box, type `1.0`
5. In the **Sgmin** box, type `-0.1`
6. In the **Sgmax** box, type `0.0`

### Capillary Pressure

To specify the capillary pressure function:

1. Click the **Capillary Press** tab

2. In the **Capillary Pressure** list, select **Linear Function**

3. In the **CPmax** box, type `9.7902E3`

4. In the **A** box, type `0.333`

5. In the **B** box, type `1.0`

Click **OK** to exit the **Additional Material Data** dialog.

Click **OK** again to save your settings and exit the **Material Data** dialog.

## Initial Conditions

The initial state of each cell in the model must be defined. The **Default Initial Conditions** dialog is used to define initial conditions that will be applied to the entire model. You can also specify initial conditions by cell, by region, or by importing the results of a previous analysis. For any analysis, the specific initial conditions will depend on several factors including EOS selection, simulator mode, and the initial state of the simulation.

Correct specification of initial conditions is essential for proper convergence and obtaining a correct result. In general, the initial conditions need to physically meaningful. Often this requires an initial state analysis in which a model is run to obtain initial equilibrium conditions before the analysis of interest (geothermal production, VOC spill, etc.) is run.

To edit global initial conditions: on the **Properties** menu, click **Initial Conditions...**

To set the initial conditions:

1. In the list, select **Pressure**

2. In the **Pressure** box, type `1.001E5`

Click **OK** to exit the **Default Initial Conditions** dialog.

## Define Boundary Condtions

To edit cells, you can use the **Grid Editor**. In this case, we will use the Grid Editor to define boundary conditions in the model.

To open the **Grid Editor**: on the **Model** menu, click **Edit Grid**.

### Water Source

We will inject into the cell on the left and produce from the cell on the right. Click on the leftmost cell in the grid (cell #1). To edit the properties of this cell, on the **Edit** menu, click **Properties**.

Click the **Properties** tab. Then, in the **Cell Name** box, type `Input`.

Click the **Sources/Sinks** tab. To define the source:

1. Under **Injection**, select **Water**
2. In the **Rate** box, type `1.16E-4`
3. Because EOS9 is an isothermal analysis, the enthalpy value need not be set

Next, click the **Print Options** tab. Select **Print Cell Time Dependent Flow and Generation (BC) Data**. This will output data for this cell at every time step, which can then be used to make detailed time history plots.

Click **OK** to close the **Edit Cell Data** dialog.

Similar steps are followed to define production in the model.

### Production

We will produce from the cell on the right. Click on the rightmost cell in the grid (cell #60). To edit the properties of this cell, on the **Edit** menu, click **Properties**.

Click the **Properties** tab. Then, in the **Cell Name** box, type `Output`.

Select the **Sources/Sinks** tab. To define the production:

1. Under **Production**, select **Mass Out**
2. In the **Rate** box, type `1.16E-4`

Next, click the **Print Options** tab. Select **Print Cell Time Dependent Flow and Generation (BC) Data**.

Click **OK** to close the **Edit Cell Data** dialog.

## Print Center Cell Data

In the grid editor, we will also choose a cell in the center of the model for which time history data will be printed. Right-click on a cell near the center (for example, cell #30) to edit the cell properties.

Click the **Properties** tab. Then, in the **Cell Name** box, type `Center`

Unlike the Input and Output cells, do not set any boundary condition data for the Center cell

Click the **Print Options** tab. Select **Print Cell Time Dependent Flow and Generation (BC) Data**.

Click **OK** to close the **Edit Cell Data** dialog.

On the **File** menu, click **Close** to exit the grid editor.

## Solution Controls

We will now define the solution options. Options relating the time step and other solution controls can be found in the **Solution Controls** dialog.

To open the **Solution Controls** dialog: on the **Analysis** menu, click **Solution Controls...**

### Times

1. In the **Solution Controls** dialog, click the **Times** tab
2. In the **End Time** box, type `8.64E6`
3. In the **Time Step** box, type `10.0`
4. In the **Max Num Time Steps** list, type `1000`
5. In the **Max Time Step** list, select `User Defined`
6. In the **Max Time Step** box, type `8.64E3`

For consistency with the example problem, we will now change some of the default values. In most cases, these values would not be changed.

### Weighting

1. In the **Solution Controls** dialog, click the **Weighting** tab
2. As the **Density at Interface** option, select **Average of Adjacent Elements**

### Convergence

1. In the **Solution Controls** dialog, click the **Convergence** tab
2. In the **Relative Error Criterion** box, type `1.0E-6`

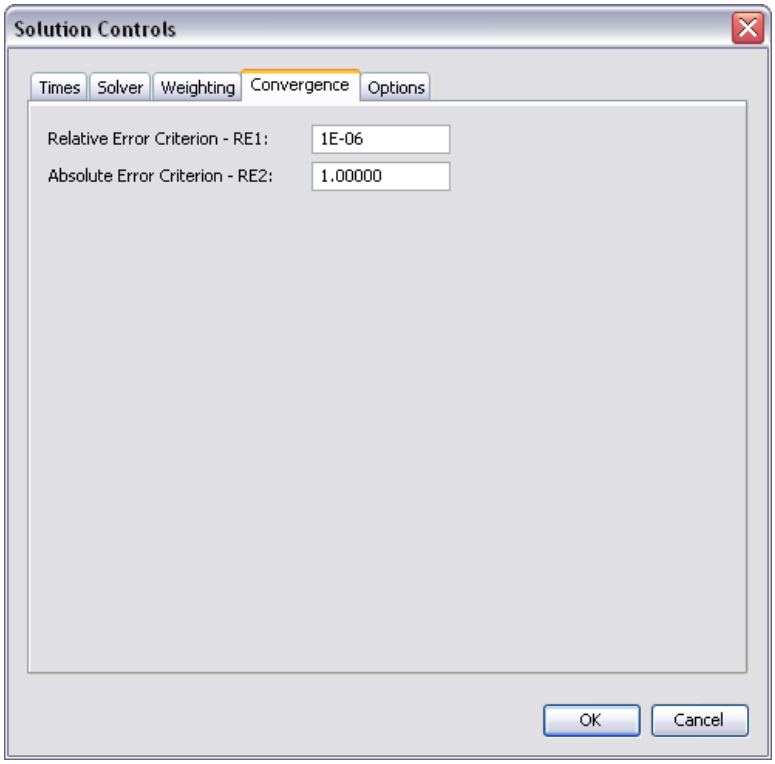


**Figure 2.3. Solution Controls - Convergence**

Click **OK** to exit the **Solution Controls** dialog.

## Output Controls

By default, the simulation will print output every 100 time steps. We can change the resolution of the output in the **Output Controls** dialog.

1. On the **Analysis** menu, click **Output Controls...**

2. In the **Print and Plot Every # Steps** box, type `500`

3. In the **Additional Output Data**group, select **Fluxes and Velocities**, **Primary Variables**, and **Additional T2React Variables**

In addition to printing output every 500 steps, we can also specify times for which we want to view data in the **Additional Print Times** dialog.

To specify additional times for output:

1. On the **Output Controls** dialog, click the **Edit** button to open the **Additional Print Times** dialog

2. In the **Times** table, type `4.32E6` and `8.64E6`

3. Click **OK** to exit the **Additional Print Times** dialog.

Click **OK** to exit the **Output Controls** dialog.

## TOUGHREACT Solution Parameters

We will now set the TOUGHREACT parameters. In this example, we are doing this last, since the entire model will lie in the same zone. However, if we wanted to define different zones in the model, we would specify the TOUGHREACT parameters first.

TOUGHREACT solution parameters can be entered on the **Solution Parameters** dialog. To open the **Solution Parameters** dialog: on the **Tough React** menu, click **Solution Parameters...**.

Select **Advanced** from the list on the left, to display the **Advanced Options** pane. Under the Advanced Options, select **Print Porosity, Permeability, Capillary Pressure Changes**.

Next, select **Times and Convergence** from the list on the left, to open the **Time Stepping and Convergence Options** pane.

In the **Max Iterations to Solve Geochemical System** box, type `300`.

In the **Relative Sorption Concentration Tolerance** box, type `1.0E-6`.

Click **OK** to exit the **Solver Parameters** dialog.

## TOUGHREACT Output Options

TOUGHREACT output options can be changed on the **Output Options** dialog.

1. On the **Tough React** menu, click **Output Options...**.
2. In the **Grid Block Output Frequency(s)** box, type `40`.
3. For **Aqueous Concentration Output**, select **Write Total Aqueous Component Concentrations**.
4. For **Aqueous Concentration Units**, select **mol/L Liquid**.
5. For **Mineral Abundance Units**, select **Change in Volume Fraction**.

Click **OK** to exit the **Output Options** dialog.

## TOUGHREACT Chemical Components

TOUGHREACT chemical components can be specified in the **Chemical Components** dialog. To open the **Chemical Components** dialog: on the **Tough React** menu, select **Chemical Components...**.

To define the primary species:

1. In the list on the left of the **Chemical Components** dialog, select **Primary Species**

2. In the **Thermodynamic Database** list in the middle of the dialog, select **h+**, **h2o**, **na+**, **skdd1**, **skdd2**, and **skdd3**

3. Click the **-->** button to move the selected species into the **Current Simulation** list on the right, as shown in Figure 2.4.

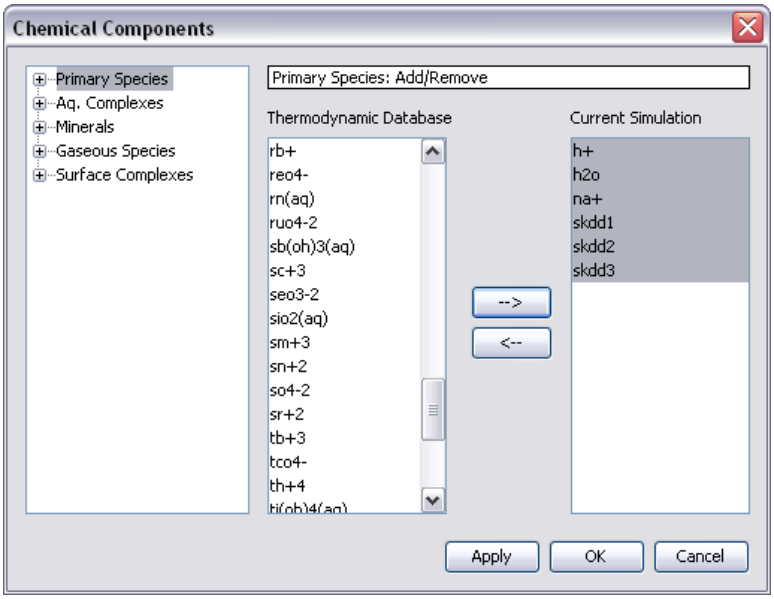


**Figure 2.4. Primary Species**

4. Click **Apply** to add the selected species to the analysis

The parameters specific to each type can be viewed and changed by clicking on that type in the subtree under **Primary Species** in the list on the left.

To edit parameters for na+ :

1. Select **na+** in the list under **Primary Species**.

2. In the pane on the right, select **Output Concentration History at Selected Cells**. This will output additional data for cells that have previously been identified for printing time history data.

To edit parameters for skdd1 :

1. Select **skdd1** in the list

2. In the pane on the right, select **Output Concentration History at Selected Cells**
3. Select **Enable Kd and Decay**.
4. In the **Decay Constant** box, type `0.0`

To edit parameters for skdd2 :

1. Select **skdd2** in the list
2. In the pane on the right, select **Output Concentration History at Selected Cells**
3. Select **Enable Kd and Decay**.
4. In the **Decay Constant** box, type `4.0113E-7`

To edit parameters for skdd3 :

1. Select **skdd3** in the list
2. In the pane on the right, select **Output Concentration History at Selected Cells**
3. Select **Enable Kd and Decay**.
4. In the **Decay Constant** box, type `4.0113E-7`

Click **OK** to exit the **Chemical Components** dialog.

## TOUGHREACT Zone Data

The next task is to create the zone data, which can be done in the **Geochemical Zones** dialog. To open the **Geochemical Zones** dialog: on the **Tough React** menu, click **Zone Data....** Initially, there will not be any zones in the mdoel.

To create the initial water zone:

1. Select **Water (Initial)** in the list on the left
2. Click **Add**
3. In the **Create a New Zone** dialog, type `Water Zone` and click **OK**
4. Click **Apply**
5. In the list on the left, click + beside **Water (Initial)** to expand the subtree.
6. In the subtree under **Water (Initial)**, click **Water Zone** to display the zone parameters to the right of the list.

7. Enter the data that is shown in Table 2.2

**Table 2.2. Water Zone Data**

| Species | Constraint | CGUESS | CTOT |
|---|---|---|---|
| h+ | Amount (mol) | 1.0e-7 | 1.0e-7 |
| h20 | Amount (mol) | 1.0 | 1.0 |
| na+ | Amount (mol) | 1.0e-10 | 1.0e-10 |
| skdd1 | Amount (mol) | 1.0e-10 | 1.0e-10 |
| skdd2 | Amount (mol) | 1.0e-10 | 1.0e-10 |
| skdd3 | Amount (mol) | 1.0e-10 | 1.0e-10 |

To create the boundary water zone:

1. Select **Water (Boundary)** in the list on the left
2. Click **Add**
3. In the **Create a New Zone** dialog, type `Water Boundary Zone` and click **OK**
4. Click **Apply**
5. In the list on the left, click **+** beside **Water (Boundary)** to expand the subtree.
6. In the subtree under **Water (Boundary)**, click **Water Boundary Zone** to display the zone parameters to the right of the list
7. Enter the data that is shown in Table 2.3

**Table 2.3. Water Zone Data**

| Species | Constraint | CGUESS | CTOT |
|---|---|---|---|
| h+ | Amount (mol) | 1.0e-7 | 1.0e-7 |
| h20 | Amount (mol) | 1.0 | 1.0 |
| na+ | Amount (mol) | 1.0e-4 | 1.0e-4 |
| skdd1 | Amount (mol) | 1.0e-4 | 1.0e-4 |
| skdd2 | Amount (mol) | 1.0e-4 | 1.0e-4 |
| skdd3 | Amount (mol) | 1.0e-4 | 1.0e-4 |

To create the permeability porosity zone:

1. Select **Permeability Porosity** in the list on the left
2. Click **Add**
3. In the **Create a New Zone** dialog, type `Perm/Por Zone` and click **OK**
4. Click **Apply**
5. In the list on the left, click + beside **Permeability Porosity** to expand the subtree.
6. In the subtree under **Permeability Porosity**, click **Perm/Por Zone** to display the zone parameters to the right of the list
7. Select **Modified Cubic Law**

To create the linear Kd zone:

1. Select **Linear Kd** in the list on the left
2. Click **Add**
3. In the **Create a New Zone** dialog, type `Linear Kd Zone` and click **OK**
4. Click **Apply**
5. In the list on the left, click + beside **Linear Kd** to expand the subtree.
6. In the subtree under **Linear Kd**, click **Linear Kd Zone** to display the zone parameters to the right of the list
7. For **skdd1** and **skdd3**, in the **Density** box, type `2.6` and in the **Kd** box, type `4.2735E-2`

Click **OK** to exit the **Geochemical Zones** dialog.

## Associate Zones with Grid

The last task is to associate zones with the grid. This can be done in two ways, either by region or for individual cells in the Grid Editor. Initially the model is one region, but it can be subdivided into more regions by using internal boundaries. This can be useful, since it is also possible to assign material data and initial conditions by region. In this case, the model will remain a single region.

Click on the model to select it in entirety. You can better see the model selection by unselecting **Grid** in the **View** menu.

After the model has been selected:

1. Under the **Model** menu, click **Edit Selection** to open the **Edit Region Data** dia-

log

2. Click the **Chemical Zones** tab

3. Select **Specify Zones by Region**

4. From the **Initial Water Zone**, **Boundary Water Zone**, **Permeability/Porosity Zone**, and **Linear Kd Zone** lists, select the respective zones that have previously been created.

Click **OK** to exit the **Edit Region Data** dialog.

## Save and Run

The input is complete and you can run the simulation. If you haven't already, you may want to save your model in a directory specifically intended for the simulation results. For example:

1. On the **File** menu, click **Save As...**

2. In the **File Name** box, type `C:\t2react_example1\t2react_example1.sim`

3. Click **Save**

To run the simulation: on the **Analysis** menu, click **Run T2REACT**

## View 3D Results

To view the 3D results for a simulation: on the **Results** menu, click **3D Results**. The data for the current simulation will be automatically loaded and displayed.

Because this is a 1-D model, the isosurfaces do not display the data well. To turn off the isosurfaces: on the **View** menu, click **Show Isosurfaces**

We will instead define a slice plane through the model to better display the output data. To create a slice plane:

1. On the **Results** menu, click **Slice Planes**

2. In the topmost **Axis** list, select **Z**

3. In the topmost **Coord** box, type `0.5`

4. Click **Close** to close the **Slice Planes** dialog

To view 3D results, as shown in Figure 2.5:

1. In the **Time(s)** list, select **0.13689**

2. In the **Scalar** list, select **na+**

3. Use the mouse to rotate the model to a good viewing perspective

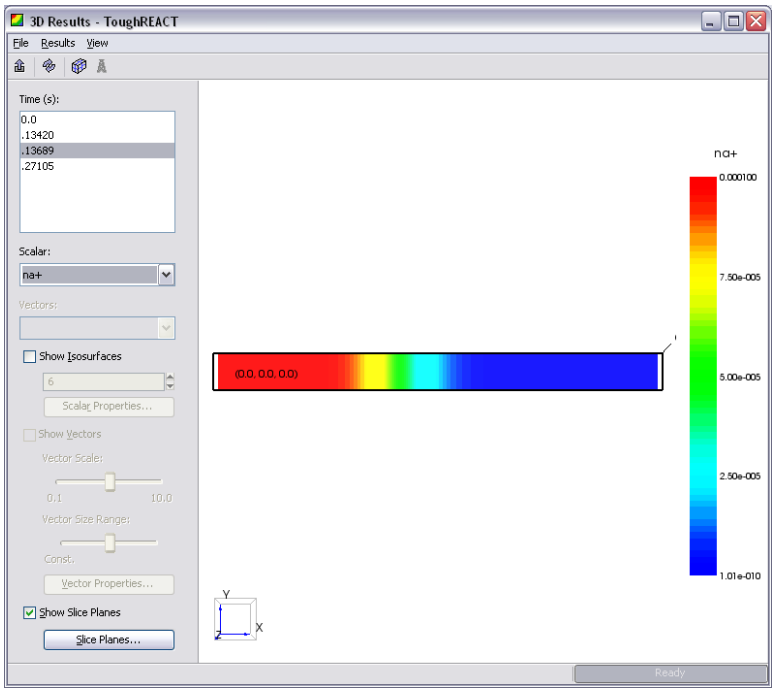


**Figure 2.5. 3D Results**

On the **File** menu, click **Close** to close the **3D Results** dialog.

## View Cell History Plots

You can view time history plots with the Cell History dialog. On the **Results** menu, click **Cell History Plots....**.

The **Cell History** dialog will be displayed. In this window, you can display time history data using a plotting parameter and a list of cells. For example, to view the isosurfaces for the Center cell, as shown in Figure 2.6:

1. In the **Variable** list, select **na+**

2. In the **Cell Name (Id#)** list, select **Center**

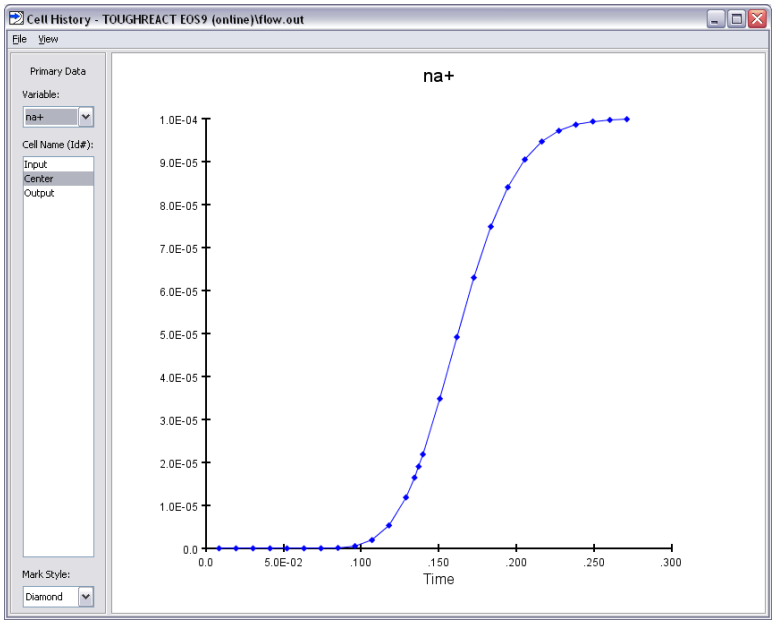


**Figure 2.6. Cell History**

In the **File** menu, click **Close** to close the Cell History window.

# Chapter 3. CO2 Disposal in Deep Saline Aquifers (ECO2)

## Description

This problem is example five in the TOUGHREACT manual. It is a 1-D radial problem in which CO2 is injected into a well field 100 m in depth and extending with a 100,000 m radius. CO2 is injected at a rate of 90 kg/s, which is approximately equivalent to that generated by a 300 MW coal-fired power plant. The CO2 injection continues for 100 years. This problem demonstrates the restart feature to run for an additional 900 years with no additional CO2 injection.

Entering some parameters for the geochemical system in this example problem can be time-intensive. To bypass entering the parameters for chemical components, please open the file `geochem_basic.sim` after updating your EOS preferences (but before editing global properties). This file is available online via the PetraSim documentation page or on your PetraSim install CD.

## Specify the Equation of State (EOS)

To ensure that PetraSim uses TOUGHREACT and ECO2, you may need to edit your PetraSim preferences. To edit preferences, you can use the **Preferences** dialog.

1. On the **File** menu, click **Preferences...**
2. In the **Simulator Mode** list, select **TOUGHREACT**
3. In the **Default Equation of State (EOS)** list, select **ECO2N**
4. Click **OK**

Your preferences have now been updated, but you will need to create a new model or restart PetraSim before they become active. If you already have a model open, it will not be affected by the preference change.

On the **File** menu, click **New**.

## Global Properties

We will first define all material and chemical properties, beginning by defining analysis options related to the equation of state (ECO2).

To edit global properties: on the **Properties** menu, click **Global Properties...**

1. Click the **Analysis** tab
2. In the **Name** box, type `CO2 Injection`
3. Click the **EOS** tab

4. Select **Isothermal**

5. In the **Brine Density in CO2** box, click **Independent**.

Click **OK** to exit the **Global Properties** dialog.

## Material Properties

We next define the material properties. Since this problem uses only one material, we will simply modify the default material.

To edit material properties: on the **Properties** menu, click **Materials...**

### Material Data

1. In the **Name** box, type `SAND`
2. In the **Porosity** box, type `0.3`
3. In the three (x, y, and z) permeability boxes, type `1.0E-13`
4. In the **Wet Heat Conductivity** box, type `2.51`
5. In the **Specific Heat** box, type `920`

### Relative Permeability

1. Click **Relative Perm...** to open the **Additional Material Data** dialog
2. Click the **Relative Perm** tab
3. In the **Relative Permeability** list, select **van Genuchten-Mualem Model**
4. In the **RP(1)** box, type `0.457`
5. In the **Slr** box, type `0.3`
6. In the **Sls** box, type `1.0`
7. In the **Sgr** box, type `0.05`

### Capillary Pressure

1. Click the **Capillary Press** tab
2. In the **Capillary Pressure** list, select **van Genuchten Function**
3. In the **CP(1)** box, type `0.457`

4. In the **Slr** box, type `0.0`
5. In the **1/P0** box, type `5.1E-5`
6. In the **Pmax** box, type `1.0E7`
7. In the **Sls** box, type `0.999`

### Miscellaneous

1. Click the **Misc** tab
2. In the **Pore Compressibility** box, type `4.5E-10`

Click **OK** to exit the **Additional Material Data** dialog.

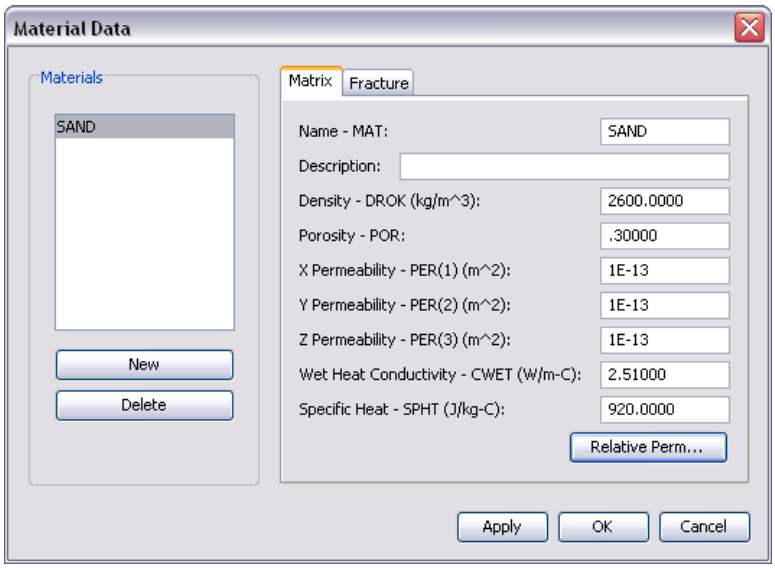


**Figure 3.1. SAND Material Data**

Click **OK** again to exit the **Material Data** dialog.

## Initial Conditions

To open the **Default Initial Conditions** dialog: on the **Properties** menu, click **Initial Conditions...**

1. In the **ECO2** list, select **Two Fluid Phases (P, Xsm, Sg, T)**
2. In the **Pressure** box, type `2.0E7`
3. In the **Temperature** box, type `75.0`

4. In the **Gas Saturation** box, type `0.0`

5. In the **Salt Mass Fraction** box, type `0.06`

Click **OK** to exit the **Default Initial Conditions** dialog.

## TOUGHREACT Solution Parameters

TOUGHREACT solution parameters can be set in the **Solver Parameters** dialog. To open the **Solver Parameters** dialog: on the **Tough React** menu, click **Solution Parameters...**

Click **Standard**

1. Click to select **Enable Gaseous Species Transport**

Click **Advanced**

1. Click to de-select **Ignore Mineral Dissolution/Precipitation Effects on Flow**

2. Under **Effect of CO2 and H2O Reactions on Flow**, click to select **CO2 and H2O**

Click **Validation**

1. In the **Max Stochiometric Ionic Strength** box, type `6.0`

Click **Times and Convergence**

1. In the **Relative Transport Concentration Tolerance**, type `1.0E-6`

2. In the **Relative Sorbtion Concentration Tolerance**, type `1.0E-6`

Click **Diffusion Coefficients**

1. In the **Aqueous Species Diffusion Coefficient** box, type `1.0E-9`

2. In the **Gaseous Species Diffusion Coefficient** box, type `1.1E-5`

Click **OK** to exit the **Solver Parameters** dialog.

## TOUGHREACT Output Parameters

TOUGHREACT output parameters can be set in the **Output Options** dialog. To open the **Output Options** dialog: on the **Tough React** menu, click **Output Options...**

1. Under **Aqueous Concentration Output**, select **Write Total Aqueous Component Concentrations**

2. Under **Mineral Abundance Units**, select **Change in Volume Fraction**

Click **OK** to exit the **Output Options** dialog.

## TOUGHREACT Chemical Components

TOUGHREACT chemical components can be specified in the **Chemical Components** dialog. To open the **Chemical Components** dialog: on the **Tough React** menu, select **Chemical Components...**.

The thermodynamic database contains the master list of chemical components. To use a species (of any type), we first add the species to the simulation, then configure the simulation-specific properties for that species.

### Primary Species

To define the primary species:

1. In the list on the left of the **Chemical Components** dialog, select **Primary Species**

2. In the **Thermodynamic Database** list in the middle of the dialog, select **alo2-**, **ca+2**, **cl-**, **fe+2**, **h+**, **h2o**, **hco3-**, **k+**, **mg+2**, **na+**, **o2(aq)**, **sio2(aq)**, and **so4-2**.

3. Click the **-->** button to move the selected species into the **Current Simulation** list on the right

4. Click **Apply** to add the selected species to the analysis

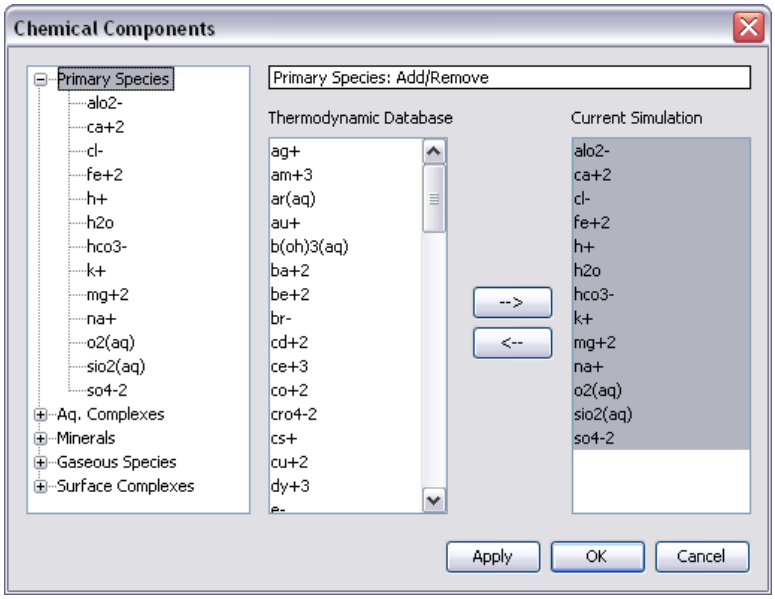


**Figure 3.2. Primary Species**

The parameters specific to each type can be viewed and changed by clicking on that type in the subtree under **Primary Species** in the list on the left. For this simulation, there is no need to edit any additional parameters.

## Aqueous Complexes

A similar process is used to define the aqueous complexes:

1. In the list on the left of the **Chemical Components** dialog, select **Aq. Complexes**

2. For **Specify Secondary Species**, select **Select Species Individually**

3. In the **Thermodynamic Database** list in the middle of the dialog, select **acetic~acid(aq)**, **al(oh)2+**, **al(oh)3(aq)**, **al+3**, **aloh+2**, **cacl+**, **cacl2(aq)**, **caco3(aq)**, **cahco3+**, **caoh+**, **caso4(aq)**, **ch4(aq)**, **co2(aq)**, **co3-2**, **fe+3**, **fecl+**, **fecl4-2**, **feco3(aq)**, **fehco3+**, **h2(aq)**, **h2s(aq)**, **h3sio4-**, **halo2(aq)**, **hs-**, **hso3-**, **kcl(aq)**, **kso4-**, **mgcl+**, **mghco3+**, **mgso4(aq)**, **naalo2(aq)**, **nacl(aq)**, **naco3-**, **nahco3(aq)**,**nahsio3(aq)**, **naoh(aq)**, **naso4-**, **oh-**, and **so2(aq)**

4. Click the **-->** button to move the selected species into the **Current Simulation** list on the right

5. Click **Apply** to add the selected species to the analysis

## Minerals

A similar process is used to define the minerals:

1. In the list on the left of the **Chemical Components** dialog, select **Minerals**

2. In the **Thermodynamic Database** in the middle of the dialog, select **albite~low**, **ankerite-2**, **calcite**, **chlorite**, **dawsonite**, **dolomite-2**, **hematite**, **illite**, **k-feldspar**, **kaolinite**, **magnesite**, **oligoclase**, **pyrite-2**, **quartz**, **siderite-2**, **smectite-ca**, and **smectite-na**

3. Click the **-->** button to move the selected species into the **Current Simulation** list on the right

4. Click **Apply** to add the selected minerals to the analysis

The parameters specific to each mineral can be viewed and changed by clicking on that type in the subtree under **Minerals** in the list on the left.

Each mineral has unique dissolution, precipitation, and additional mechanisms properties. The values for each of these properties for each mineral are shown in Table 3.1, and Table 3.2. Pyrite is a bit different than the other minerals and its properties are shown in Table 3.3 and Table 3.4. This section will demonstrate how to specify the properties for `albite~low`, then you should be able to enter the remaining mineral properties by following the same pattern and referring to the tables.

Please note, if table entries are blank, then those parameters should not be set. The mineral `calcite` is considered to be at equilibrium and no parameters for dissolution, precipitation, or additional mechanisms should be specified. The mineral `quartz` has no additional mechanisms. Several minerals have only 1 additional mechanism.

Select **albite~low** in the list under **Minerals**.

To edit dissolution rate parameters for `albite~low`:

1. Under **Specify Kinetic Constraints**, click to select **Dissolution**.
2. Click **Edit...** This will open the **Edit dissolution Rate Parameters** dialog.
3. In the **Activation Energy - EA** box, type `69.8` (from Table 3.1).
4. In the **Rate Constant $k_{25}$** box, type `2.7542e-13` (from Table 3.1).
5. Click **OK** to save changes and exit the **Edit Dissolution Rate Parameters** dialog.

To edit precipitation rate parameters for `albite~low`:

1. Under **Specify Kinetic Constraints**, click to select **Precipitation**.
2. Click **Edit...** This will open the **Edit Precipitation Rate Parameters** dialog.
3. On the **Options** tab, in the **Initial Volume Fraction** box, type `1.0e-6`.
4. Click the **Energy** tab.
5. In the **Activation Energy - EA** box, type `69.8` (from Table 3.1).
6. In the **Rate Constant $k_{25}$** box, type `2.7542e-13` (from Table 3.1).
7. Click **OK** to save changes and exit the **Edit Dissolution Rate Parameters** dialog.

To add additional mechanisms for `albite~low`:

1. Under **Rate Constant Dependence on pH**, click to select **Specify Additional Mechanisms**.
2. Click **Edit...** This will open the **Edit Additional Mechanisms** dialog.
3. Click **New**. This will create **Mechanism (1)**.
4. In the **Weighting Factor** box, type `6.9183e-11` (from Table 3.2)
5. In the **Activation Energy** box, type `65.0` (from Table 3.2)
6. In the table, in the **Species** list, select **h+** (from Table 3.2).

7. In the table, in the **Exponent** box, type `0.457` (from Table 3.2).
8. Click **New**. This will create **Mechanism (2)**.
9. In the **Weighting Factor** box, type `2.5119e-16` (from Table 3.2)
10. In the **Activation Energy** box, type `71.0` (from Table 3.2)
11. In the table, in the **Species** list, select **h+** (from Table 3.2).
12. In the table, in the **Exponent** box, type `-0.572` (from Table 3.2).
13. Click **OK** to save changes and exit the **Edit Dissolution Rate Parameters** dialog.

All other selected minerals should be edited in this fashion, using the values shown in Table 3.1 and Table 3.2. Data for `pyrite-2` can be found in Table 3.3 and Table 3.4. Empty table cells indicate that the existing data is already correct and that new data need not be entered (i.e. if there is no entry for a particular option, do not enable that option).

**Table 3.1. Dissolution and Precipitation Data for Minerals**

| Mineral | Activation Energy | Rate Constant $k_{25}$ |
|---|---|---|
| albite~low | 69.8 | 2.754E-13 |
| ankerite-2 | 62.76 | 1.260E-09 |
| calcite | | |
| chlorite | 88 | 3.020E-13 |
| dawsonite | 62.76 | 1.260E-09 |
| dolomite-2 | 52.2 | 2.951E-08 |
| hematite | 66.2 | 2.514E-15 |
| illite | 35 | 1.660E-13 |
| k-feldspar | 38 | 3.890E-13 |
| kaolinite | 22.2 | 6.918E-14 |
| magnesite | 23.5 | 4.571E-10 |
| oligoclase | 69.8 | 1.445E-12 |
| quartz | 87.7 | 1.023E-14 |
| siderite-2 | 62.76 | 1.260E-09 |
| smectite-ca | 35 | 1.660E-13 |
| smectite-na | 35 | 1.660E-13 |

**Table 3.2. Additional Mechanism Data for Minerals**

| | Mechanism(1) | | | Mechanism(2) | | |
|---|---|---|---|---|---|---|
| **Mineral** | **Weighting Factor $k_{25}$** | **Activation Energy** | **n(H+) Exponent** | **Weighting Factor $k_{25}$** | **Activation Energy** | **n(H+) Exponent** |
| albite~low | 6.918E-11 | 65.0 | 0.457 | 2.512E-16 | 71 | -0.572 |
| ankerite-2 | 6.457E-04 | 36.1 | 0.500 | | | |
| calcite | | | | | | |
| chlorite | 7.762E-12 | 88.0 | 0.500 | | | |
| dawsonite | 6.457E-04 | 36.1 | 0.500 | | | |
| dolomite-2 | 6.457E-04 | 36.1 | 0.5 | | | |
| hematite | 4.074E-10 | 66.2 | 1.0 | | | |
| illite | 1.047E-11 | 23.6 | 0.340 | 3.020E-17 | 58.9 | -0.400 |
| k-feldspar | 8.710E-11 | 51.7 | 0.500 | 6.310E-22 | 94.1 | -0.823 |
| kaolinite | 4.898E-12 | 65.9 | 0.777 | 8.913E-18 | 17.9 | -0.472 |
| magnesite | 4.169E-7 | 14.4 | 1.0 | | | |
| oligoclase | 2.138E-10 | 65.0 | 0.457 | | | |
| quartz | | | | | | |
| siderite-2 | 6.457E-04 | 36.1 | 0.500 | | | |
| smectite-ca | 1.047E-11 | 23.6 | 0.34 | 3.020E-17 | 58.9 | -0.400 |
| smectite-na | 1.047E-11 | 23.6 | 0.340 | 3.020E-17 | 58.9 | -0.400 |

**Table 3.3. Dissolution and Precipitation Data for pyrite-2**

| | Dissolution | | Precipitation | |
|---|---|---|---|---|
| **Mineral** | **Activation Energy** | **Rate Constant $k_{25}$** | **Activation Energy** | **Rate Constant $k_{25}$** |
| pyrite-2 | 0.0 | 0.0 | 56.9 | 4.0e-11 |

**Table 3.4. Additional Mechanism Data for pyrite-2**

| | Mechanism(1) | | | Mechanism(2) | | |
|---|---|---|---|---|---|---|
| **Mineral** | **Weighting Factor** | **Activation Energy (Ea)** | **Exponents** | **Weighting Factor $k_{25}$** | **Activation Energy** | **Exponent** |
| pyrite-2 | 3.02e-8 | 56.9 | [h+, -0.5] [fe+3, 0.5] | 2.8184e-5 | 56.9 | [o2(aq), 0.5] |

To save these changes: click **Apply**

### Gaseous Species

1. In the list on the left of the **Chemical Components** dialog, select **Gaseous Species**
2. In the **Thermodynamic Database** list in the middle of the dialog, select **co2(g)**,
3. Click the **-->** button to move co2(g) into the **Current Simulation** list on the right
4. Click **Apply** to add the selected species to the analysis

Click **OK** to exit the **Chemical Components** dialog.

## TOUGHREACT Zone Data

Geochemical zones describe the initial chemical composition of the model. You can define geochemical zones using the **Geochemical Zones** dialog. To open the **Geochemical Zones** dialog: on the **Tough React** menu, click **Zone Data...**. Initially, there will not be any zones in the mdoel.

To create the water zone:

1. Select **Water (Initial)** in the list on the left
2. Click **Add**
3. In the **Create a New Zone** dialog, type `Water Zone` and click **OK**
4. Click **Apply**
5. In the list on the left, click **+** beside **Water (Initial)** to expand the subtree
6. In the subtree under **Water (Initial)**, click **Water Zone** to display the zone parameters to the right of the list
7. Edit each cell in the table, setting the **Constraint**, **CGUESS**, and **CTOT** values for each species as shown in Table 3.5.

**Table 3.5. Water (Initial) Zone**

| Species | Constraint | CGUESS | CTOT |
|---|---|---|---|
| alo2- | Amount (mol) | 1.078e-8 | 1.361e-8 |
| ca+2 | Amount (mol) | 4.479e-3 | 4.737e-3 |
| cl- | Amount (mol) | 0.9109 | 1.001 |

| Species | Constraint | CGUESS | CTOT |
|---|---|---|---|
| fe+2 | Amount (mol) | 2.615e-7 | 3.022e-7 |
| h+ | Amount (mol) | 8.48e-8 | 0.0432 |
| h2o | Amount (mol) | 1.0 | 1.0 |
| hco3- | Amount (mol) | 1.841e-8 | 4.562e-2 |
| k+ | Amount (mol) | 5.805e-3 | 5.980e-3 |
| mg+2 | Amount (mol) | 2.348e-5 | 2.669e-5 |
| na+ | Amount (mol) | 0.9006 | 0.9905 |
| o2(aq) | Amount (mol) | 2.763e-66 | -8.646e-2 |
| sio2(aq) | Amount (mol) | 9.203e-4 | 1.034e-3 |
| so4-2 | Amount (mol) | 1.443e-16 | 1.324e-9 |

8. Click **Apply**

To create the mineral zone:

1. Select **Mineral** in the list on the left
2. Click **Add**
3. In the **Create a New Zone** dialog, type `Mineral Zone` and click **OK**
4. Click **Apply**
5. In the list on the left, click **+** beside **Mineral** to expand the subtree
6. In the subtree under **Mineral**, click **Mineral Zone** to display the zone parameters to the right of the list
7. Edit each cell in the table, setting the **Vol. Fraction**, **Grain Radius**, **Surface Area**, and **Units** values for each mineral as shown in Table 3.6.

**Table 3.6. Mineral Zone**

| Mineral | Vol. Fraction | Grain Radius | Surface Area | Units |
|---|---|---|---|---|
| albite~low | 0 | 0.001 | 9.8 | cm^2/g mineral |
| ankerite-2 | 0 | 0.001 | 9.8 | cm^2/g mineral |
| calcite | 0.01929 | | | |

| Mineral | Vol. Fraction | Grain Radius | Surface Area | Units |
|---|---|---|---|---|
| chlorite | 0.04556 | 0.001 | 9.8 | cm^2/g mineral |
| dawsonite | 0 | 0.001 | 9.8 | cm^2/g mineral |
| dolomite-2 | 0 | 0.001 | 9.8 | cm^2/g mineral |
| hematite | 0.00497 | 0.001 | 12.87 | cm^2/g mineral |
| illite | 0.00954 | 0.001 | 151.63 | cm^2/g mineral |
| k-feldspar | 0.08179 | 0.001 | 9.8 | cm^2/g mineral |
| kaolinite | 0.02015 | 0.001 | 151.6 | cm^2/g mineral |
| magnesite | 0 | 0.001 | 9.8 | cm^2/g mineral |
| oligoclase | 0.19795 | 0.001 | 9.8 | cm^2/g mineral |
| pyrite-2 | 0 | 0.001 | 12.87 | cm^2/g mineral |
| quartz | 0.57888 | 0.001 | 9.8 | cm^2/g mineral |
| siderite-2 | 0 | 0.001 | 9.8 | cm^2/g mineral |
| smectite-ca | 0 | 0.001 | 151.63 | cm^2/g mineral |
| smectite-na | 0.03897 | 0.001 | 151.63 | cm^2/g mineral |

8. Click **Apply**

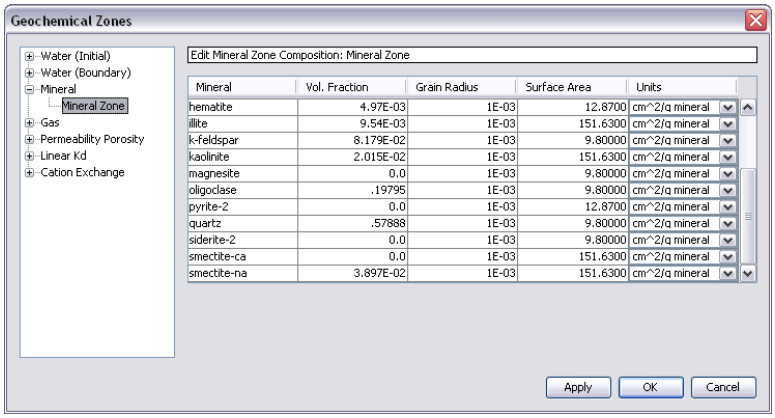


**Figure 3.3. Mineral Zone Data**

To create the permeability porosity zone:

1. Select **Permeability Porosity** in the list on the left

2. Click **Add**

3. In the **Create a New Zone** dialog, type `Perm-Pore Zone` and click **OK**

4. Click **Apply**

5. In the list on the left, click **+** beside **Permeability Porosity** to expand the subtree

6. In the subtree under **Permeability Porosity**, click **Perm-Pore Zone** to display the zone parameters to the right of the list

7. For **Permeability Law**, select **Cubic Law**

8. Click **Apply**

Click **OK** to exit the **Geochemical Zones** dialog.

## Saving the Geochemical Data File as a Starting Point for a New Analysis

All of the problem data that is not specific to a particular geometry and loading condition has now been defined. At this point, save the data by clicking **Save...** on the **File** menu, and typing "`geochemical data`" as the file name. We can now open "geochemical data" and click **Save As...** on the **File** menu, giving the new name as "`eco2_example`". Anytime you want to create a model that uses the same (or similar) geochemical data, you can open the "geochemical data" file and save it as a different file as a starting point for a different analysis. In this manner you can avoid repeating all the data input.

In continuing this example, model specific data will now be entered and simulated.

## Create the Model Boundary

To create the boundary for this model, use the **Define Model Boundary** dialog. The boundary parameters for this model are shown in Table 3.7.

**Table 3.7. Model Boundary Dimensions**

| Axis | Min (m) | Max (m) |
|---|---|---|
| x | 0.0 | 100000.0 |
| y | 0.0 | 1.0 |
| z | -100.0 | 0.0 |

To create the model boundary:

1. On the **Model** menu, click **Define Boundary...**
2. Enter the values from Table 3.7
3. Click **OK** to create the model boundary

## Create the Grid

To create the grid:

1. On the **Model** menu, click **Create Grid...**
2. In the **Division Method** box, select **Regular**
3. For **Grid Type**, select **RZ Grid**
4. In the **X Cells** box, type `100`
5. In the **X Factor** box, type `1.1096251`
6. In the **Z Cells** box, type `1`
7. In the **Z Factor** box, type `1.0`
8. Click **OK** to create the grid

This 100 cell grid used here is different than the 130 cell meshmaker grid used in the example problem writeup from the TOUGHREACT user's manual. The style of the grid used in TOUGHREACT user's manual would require a logarithmic capability in PetraSim's meshmaker grid creater that is currently not supported. The 100 cell logarithmic grid used in this example problem appears to be sufficient based on a sensitivity analysis (Figure 3.4) of the Sg parameter after 100 years of simulation with different grid sizes.

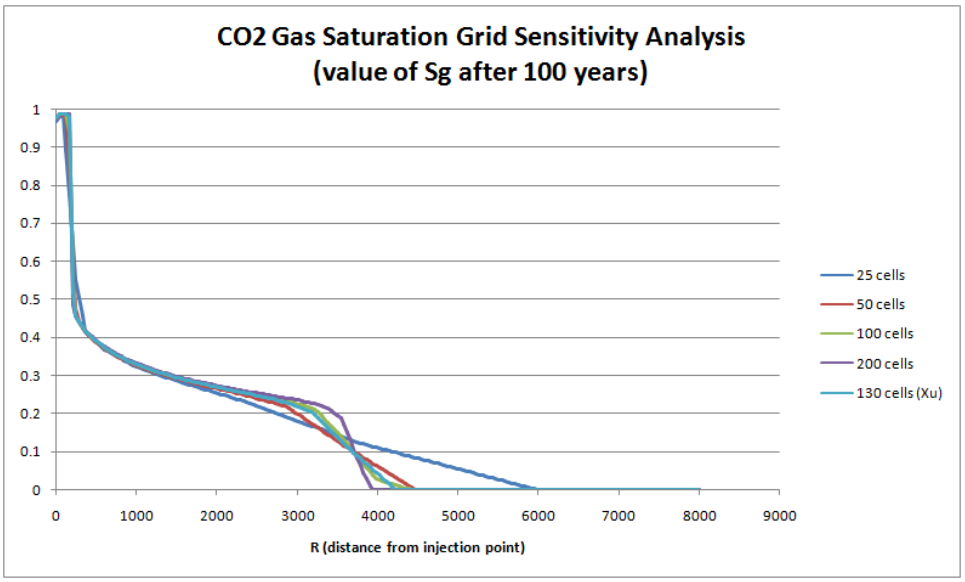


**Figure 3.4. CO2 Injection**

To use a 200 cell grid, change the value of the **X Cells** paramter to `200` and the value of the **X Factor** parameter to `1.04950055`. The running time of this simulation is linear in the number of cells. It takes approximately 10 minutes to run the first 100 years of the simulation with 100 cells, this time would be doubled for 200 cells.

## Define Boundary Condtions

To edit cells, you can use the **Grid Editor**. In this case, we will use the Grid Editor to define boundary conditions in the model.

To open the **Grid Editor**: on the **Model** menu, click **Edit Grid**. In this model, 90 kg/s of CO2 is injected into the center. Since this is an RZ grid, the center cell is the min x cell (far left). Since this is an isothermal simulation, it will not be necessary to specify an enthalpy for the injected CO2 - we will leave this value at 0.0.

To select and edit the injection cell:

1. In the **Find** box, type `1`, then press `Enter`. Cell 1 will be selected and centered in the Grid Editor.

2. On the **Edit** menu, click **Properties**. This will open the Edit Cell Data dialog.

To specify the properties for the injection cell:

1. Click the **Sources/Sinks** tab.

2. Under **Injection**, click to select **CO2**

3. In the **Rate** box, type `90.0`

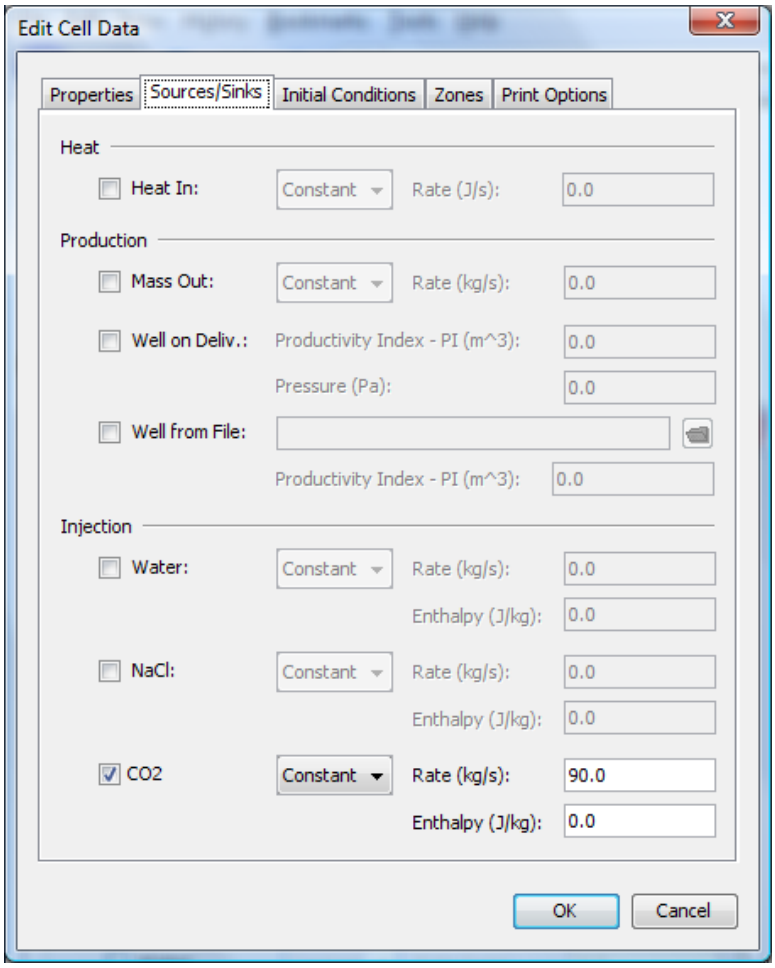


**Figure 3.5. CO2 Injection**

Click **OK** to close the **Edit Cell Data** dialog. On the **File** menu, click **Close** to close the **Grid Editor**.

## Solution Controls

We will now define the solution options. Options relating the time step and other solution controls can be found in the **Solution Controls** dialog.

To open the **Solution Controls** dialog: on the **Analysis** menu, click **Solution Controls...**

### Times

1. In the **Solution Controls** dialog, click the **Times** tab

2. In the **End Time** box, type `3.1557E9` (100 years).
3. In the **Time Step** box, type `1.0`
4. In the **Max Num Time Steps** list, click **Infinite**
5. In the **Max Time Step** list, select `User Defined`
6. In the **Max Time Step** box, type `2.592E6` (30 days).

### Solver

1. Click the **Solver** tab
2. For the **Conjugate Gradient Solvers**, select **Stabilized Bi-Conjugate Gradient**
3. In the **Max CG Iterations (Frac. of Eqns.)** box, type `0.8`
4. In the **CG Convergence Criterion** box, type `1E-7`

Click **OK** to exit the **Solution Parameters** dialog.

## Output Controls

By default, the simulation will print output every 100 time steps. We can change the resolution of the output in the **Output Controls** dialog.

1. On the **Analysis** menu, click **Output Controls...**
2. In the **Print and Plot Every # Steps** box, type `9999` (we will specify specific output times).

To specify specific times for output:

1. On the **Output Controls** dialog, click the **Edit** button to open the **Additional Print Times** dialog
2. In the **Times** table, type `3.15576E7`, `3.15576E8`, `1.57788E9`, `3.15576E9`, `1.57788E10`, `3.15576E10`, `6.31152E10`, and `1.57788E11`.
3. Click **OK** to exit the **Additional Print Times** dialog.

Click **OK** to exit the **Output Controls** dialog.

## Associate Zones with Grid

The last task is to associate zones with the grid. This can be done in two ways, either by region or for individual cells in the Grid Editor. Initially the model is one region,

but it can be subdivided into more regions by using internal boundaries. This can be useful, since it is also possible to assign material data and initial conditions by region. In this case, the model will remain a single region and we will specify zones for the entire domain.

To specify zone data:

1. In the tree view at the left, under **Regions**, click to select **Region 1**.

2. Under the **Model** menu, click **Edit Selection** to open the **Edit Region Data** dialog

3. Click the **Chemical Zones** tab

4. Select **Specify Zones by Region**

5. From the **Initial Water Zone**, **Mineral Zone**, and **Permeability/Porosity Zone** lists, select the respective zones that have previously been created.

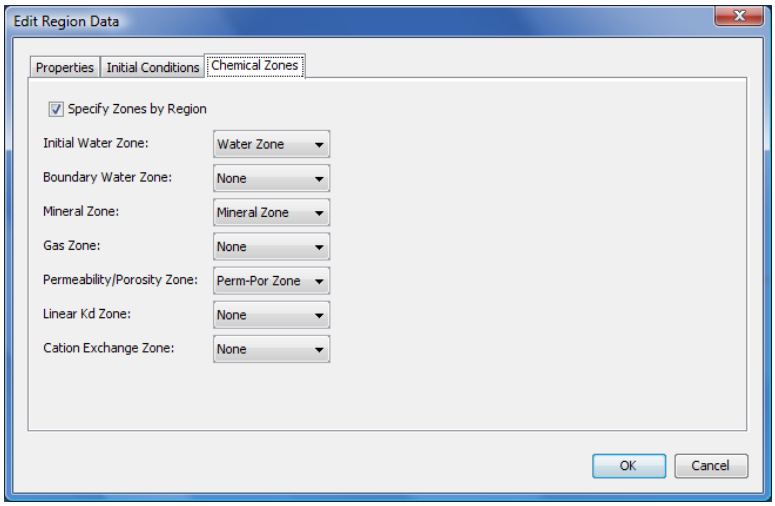


**Figure 3.6. Zones Associated with Grid**

Click **OK** to exit the **Edit Region Data** dialog.

## Save and Run

The input is complete and you can run the simulation. To retain the geochemical data as a reusable starting point, save this model as a different file in a separate directory. For example:

1. On the **File** menu, click **Save As...**

2. In the **File Name** box, type `C:\t2react_co2dsa\eco2_co2sda.sim`

3. Click **Save**

To run the simulation: on the **Analysis** menu, click **Run T2REACT**

## View Results

The best way to view results for this problem is to use the **Line Plot** feature in PetraSim. Line plots allow us to view the data in 2D as a function of distance from a point. In this case, we will view the data as it radiates from the center of the model (as a function of R). Line plots are availble in the **3D Results** view.

To create a line plot:

1. On the **Results** menu, click **3D Results**.

2. In the **3D Results** view, on the **File** menu, click Line Plot...

3. In the **Line Plot** dialog, enter the following two points: Point1=`(0.0, 0.5, -50)`, Point2=`(8000, 0.5, -50)`, then click **OK**.

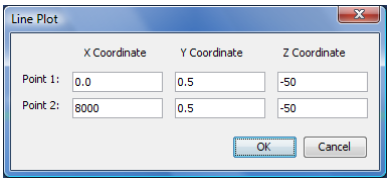


**Figure 3.7. Preparing a Line Plot**

In the **Line Plot** dialog, to show CO2 saturation (Sg):

1. In the **Variable** list, select **SG**

2. In the **Time** list, select **3.156E9**

This result is shown in Figure 3.8

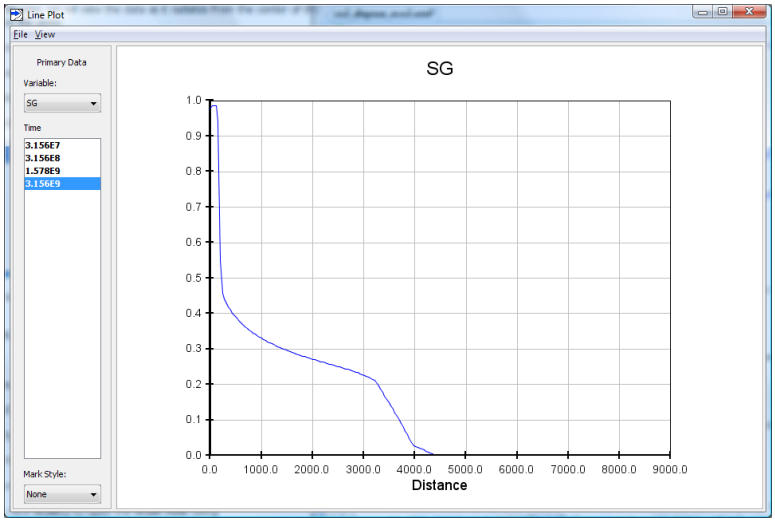


**Figure 3.8. Line Plot of CO2 Saturation (Sg)**

## The Continuation Run (Restart)

During the first part of this simulation, we injected 90 kg/s of CO2 into the aquifer for 100 years. During the second part, we will stop injecting CO2 and monitor the aquifer out to 1,000 years. Before proceeding, create a second folder on your computer for the restart run - otherwise the continuation run will overwrite the results from the first run.

1. On the **File** menu, click **Save As...**

2. In the **File Name** box, type `C:\t2react_co2dsa\restart\restart.sim`

3. Click **Save**

Next, turn off the CO2 injection.

1. In the **Find** box, type `1`, then press `Enter`. Cell 1 will be selected and centered in the Grid Editor.

2. On the **Edit** menu, click **Properties**. This will open the Edit Cell Data dialog.

3. Click the **Sources/Sinks** tab.

4. Under **Injection**, click to de-select **CO2**

5. Under **OK**, click to exit the **Edit Cell Data** dialog.

Click **OK** to close the **Edit Cell Data** dialog. On the **File** menu, click **Close** to close the **Grid Editor**.

### Add Restart Data

We will use the initial conditions and geochemical data from the end of the previous simulation, to start the continuation run. To enable the simulation restart feature:

1. On the **TOUGHREACT** menu, click **Restart Options...**
2. In the **Restart Options** dialog, click to enable **Activate Simulation Restart**.
3. For **Initial Conditions (SAVE)**, click **Browse...** and select the SAVE file from the previous run.
4. For **Geochem Data (savechem)**, click **Browse...** and select the `savechem` file from the previous run.
5. Click **OK**, to save changes and close the **Restart Options** dialog.

### Set a New End Time

To specify the new 1,000 year end time for the simulation:

1. On the **Anaysis** menu, click **Solution Controls...**
2. In the **Solution Controls** dialog, in the **End Time** box, type `3.1557E10`
3. Click **OK**, to save changes and close the **Solution Controls** dialog.

### Start the Continuation Run

On the **Analysis** menu, click **Run T2React...**

## View Results

To look at the total amount of sequestered CO2 after 1,000 years, we can create another line plot.

To create a line plot:

1. On the **Results** menu, click **3D Results**.
2. In the **3D Results** view, on the **File** menu, click Line Plot...
3. In the **Line Plot** dialog, enter the following two points: Point1=`(0.0, 0.5, -50)`, Point2=`(8000, 0.5, -50)`, then click **OK**.

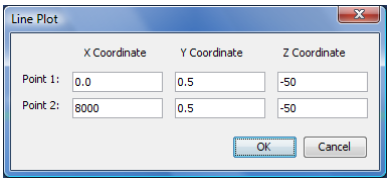

**Figure 3.9. Preparing a Line Plot**

In the **Line Plot** dialog, to show total CO2 sequestered in minerals (SMco2):

1. In the **Variable** list, select **SMco2**
2. In the **Time** list, select **3.156E10**

This result is shown in Figure 3.10

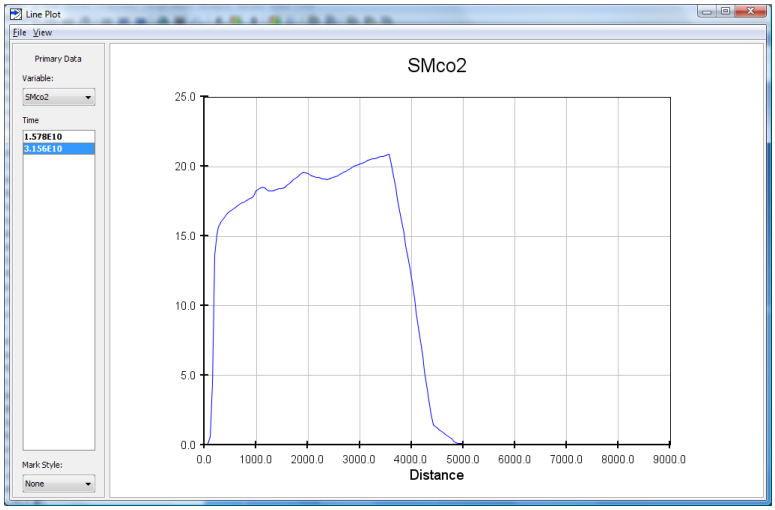


**Figure 3.10. Line Plot of Total CO2 Sequestered in Minerals (SMco2)**

You can also look at the mineral abundance as a function of x (R), using the **Line Plot** dialog.

# References

[Pruess, Oldenburg, and Moridis, 1999] Karsten Pruess, Curt Oldenburg, and George Moridis. *TOUGH2 User's Guide, Version 2.0*. November, 1999. Earth Sciences Division, Lawrence Berkeley National Laboratory. Berkeley CA USA . LBNL-43134.

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[Pruess and Garcia, 2002] Karsten Pruess and J. Garcia. *Solutions of Test Problems for Disposal of CO2 in Saline Aquifers*. December, 2002. Earth Sciences Division, Lawrence Berkeley National Laboratory. Berkeley CA USA . LBNL-51812.

[Pruess et al., 2002] Karsten Pruess, J. Garcia, T. Kovscek, Curt Oldenburg, J. Rutqvist, C. Steefel, and T. Xu. *Intercomparison of Numerical Simulation Codes for Geologic Disposal of CO2*. December, 2002. Earth Sciences Division, Lawrence Berkeley National Laboratory. Berkeley CA USA . LBNL-51813.

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