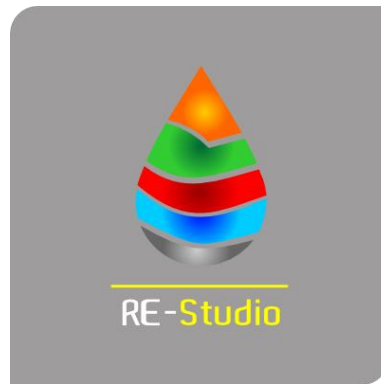
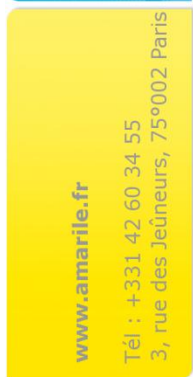


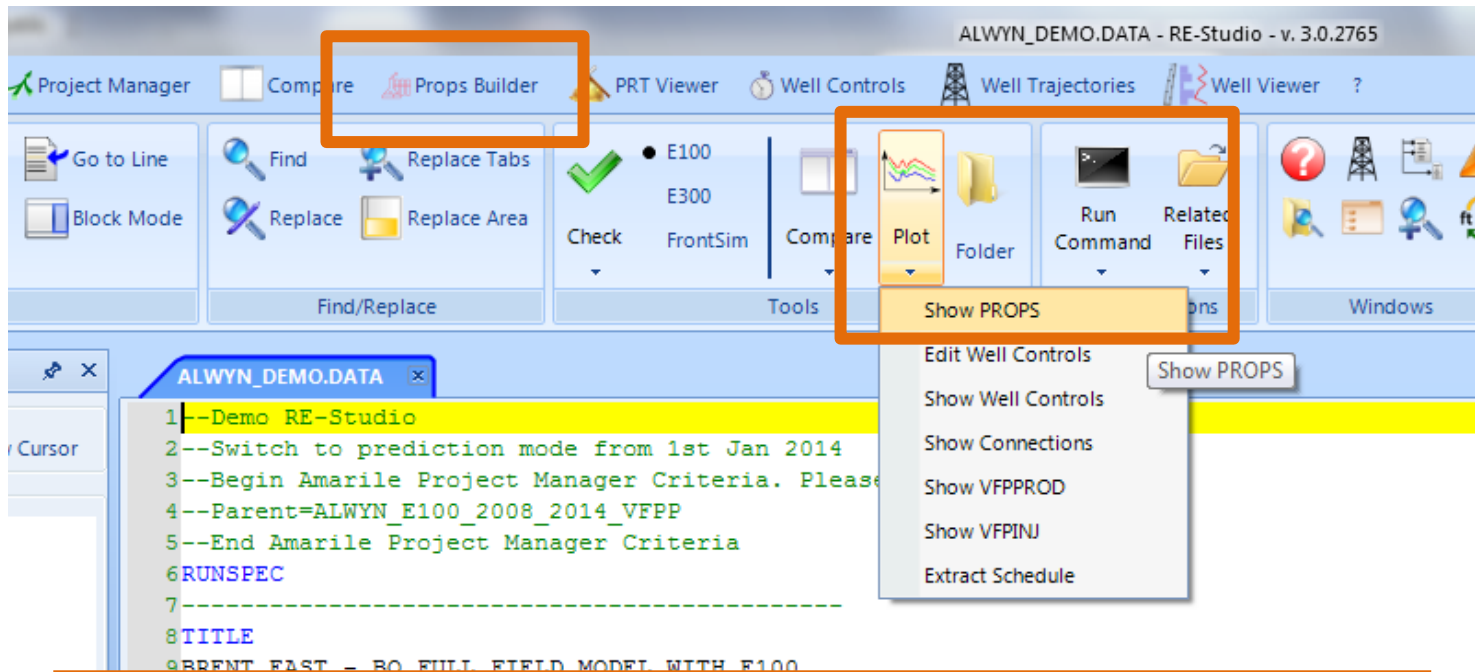
RE-Studio – Props Builder



Relative Permeability Tables Builder

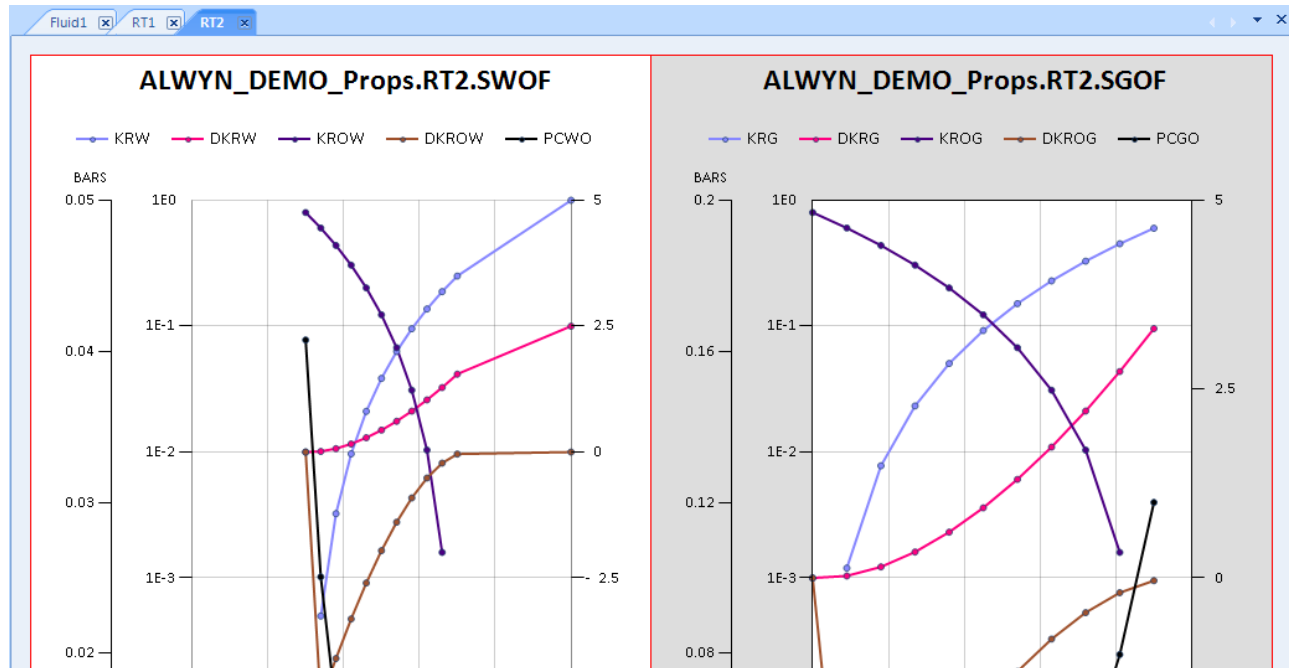


Purpose



📦 This module allows the user to create easily the saturation functions (SWOF and SGOF keywords table) for ECLIPSE datadeck based on Corey correlatoin

Plot – Show Props



- ❖ In For ECLIPSE, click “Show Props” in Plot button, RE-Studio will display Fluid and Rock Types (RT) information
 - ❖ KR: relative permeability
 - ❖ DKR: derived relative permeability
 - ❖ PC: capillary pressures
- ❖ NB: rel perm series are displayed in logarithmic scale
 - ❖ You can change this by double-clicking on the axis

Create your rock type

- First of all, you need to give a name to your rock type then it “Add”

- Select the equation you want
- Select the unit you want
- Tick the buttons to select phases

Rock types

Add rock type

Corey METRIC

Phases

OIL GAS

- You also can Load and Save your rock type
 - The rock type data will be saved in .akr (Amarile Rock Type)
 - Also notice that it's an XML format so you can edit it with a text editor

Create your rock type

- For this example, I named my rock type “RockType1” and “Add” it
- I tick both oil and gas
- I am working with the Corey equation in a Metric system
- I defined parameters in all boxes
- I can click Plot to see the functions, and Apply only if the plots are already displayed

Rock types

Add rock type

RockType2

Add

Corey

METRIC

Load

Save file

Phases

OIL

GAS

Oil Water parameters

$K_{rw} = K_{rw}(1-S_{orw}) * S_{wn}^{No}$

$K_{row} = K_{row}(S_{wi}) * (1 - S_{wn})^{No}$

Swi: 0.150

Krow(Swi): 0.800

Sorw: 0.220

Krw(1-Sorw): 0.250

No: 2.700

Nw: 2.700

Capillary Pressure

The Brooks-Corey correlation for capillary pressure reads $P_c = c^* S_w^{a(-a)}$

c: 0.000381345

a: 3.88008142

Gas-Oil parameters

$K_{rg} = K_{rg}(1 - S_{wi} - S_{org}) * S_{gn}^{Ngo}$

$K_{rog} = K_{row}(S_{wi}) * (1 - S_{gno})^{Ngo}$

Sgc: 0.100

Sorg: 0.250

Krg(1-Swi-Sorg): 0.600

Nog: 2.700

Ngo: 2.700

Capillary Pressure

The Brooks-Corey correlation for capillary pressure reads $P_c = c^* S_g^{a(-a)}$

Cg: 0.12

Ag: 3.880081

Restore

Auto apply

Apply

Plot

Write

Load your rock type

Saturation Functions

Rock types

Add rock type

RockType2

Define current rock type

RockType1

Corey METRIC

Phases

OIL GAS

Oil-Water parameters

$K_{rw} = K_{rw}(1-S_{orw}) * S_{wn}^{Nw}$ $K_{row} = K_{row}(S_{wi}) * (1 - S_{wn})^{No}$

Swi: 0.150 Krow(Swi): 0.800

Sorw: 0.220 Krw(1-Sorw): 0.250

No: 2.700 Nw: 2.700

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The Brooks-Corey correlation for capillary pressure reads $P_c = c * S_w^{(-a)}$

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Gas-Oil parameters

$K_{rg} = K_{rg}(1 - S_{wi} - S_{org}) * S_{gn}^{Nog}$ $K_{rog} = K_{rog}(S_{wi}) * (1 - S_{gn})^{Nog}$

Sgc: 0.100 Sorg: 0.250

Krg(1-Swi-Sorg): 0.600

Nog: 2.700 Nog: 2.700

Capillary Pressure

The Brooks-Corey correlation for capillary pressure reads $P_c = c * S_g^{(-a)}$

Cg: 0.12 Ag: 3.880081

Auto apply

- I can Save to file the parameters I have input
- I can Load from file such parameters
- Loading a AKR file will overwrite all the parameters for your current rock type
 - Caution: we are loading the parameters values without a name so you need to create and “Add” an empty rock type before

Edit your rock type

Saturation Functions

Rock types

Add rock type

RockType2

Define current rock type

RockType1

Corey

Phases

OIL GAS

Oil-Water parameters

$K_{rw} = K_{rw}(1-S_{orw}) * S_{wn}^{Nw}$ $K_{row} = K_{row}(S_{wi}) * (1 - S_{wn})^{No}$

Swi: 0.150 Krow(Swi): 0.800

Sorw: 0.220 Krw(1-Sorw): 0.250

No: 2.700 Nw: 2.700

Capillary Pressure

The Brooks-Corey correlation for capillary pressure reads $P_c = c * S_w^{[-a]}$

c: 0.000381345 a: 3.88008142

Gas-Oil parameters

$K_{rg} = K_{rg}(1 - S_{wi} - S_{org}) * S_{gn}^{Nog}$ $K_{rog} = K_{rog}(S_{wi}) * (1 - S_{gno})^{Nog}$

Sgc: 0.100 Sorg: 0.250

Krg(1-Swi-Sorg) 0.600

Nog: 2.700 Ngo: 2.700

Capillary Pressure

The Brooks-Corey correlation for capillary pressure reads $P_c = c * S_g^{[-a]}$

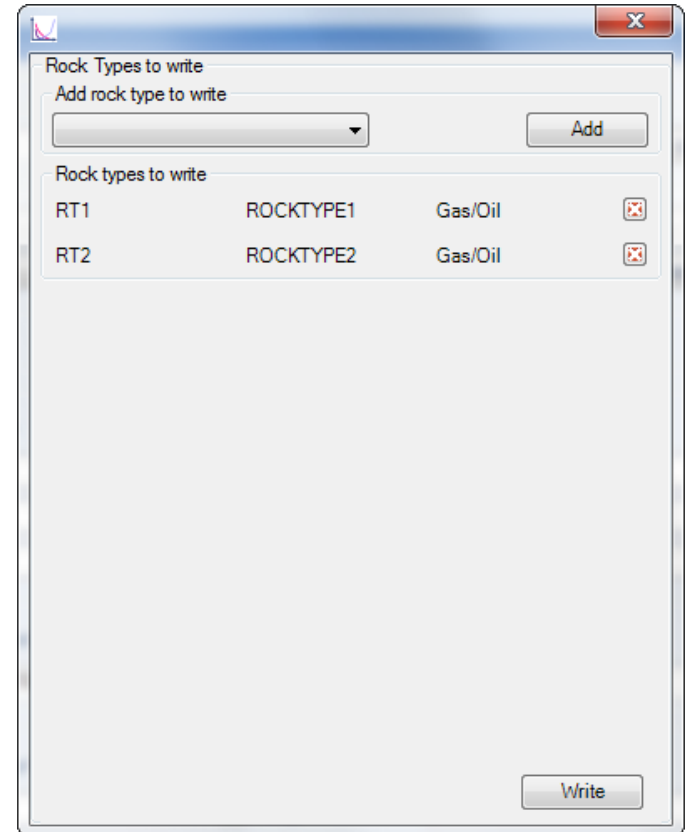
Cg: 0.12 Ag: 3.880081

Auto apply

- Once you set the parameters, you can:
 - “Restore” the last save
 - “Apply” the changes to the series already plot
 - “Plot” the rock type
 - “Write” will generate automatically the keywords for your datadeck

Write the SWOF/SGOF

- When you hit the “Write” button, a window will pop up
- You can use any number of rock types by selecting them in the list then by hitting “Add”
- When you are satisfied with your list of rock types just hit “Write”



Write the SWOF/SGOF

```
PROPS Builder - Gas/Oil
1 SWOF
2 -- Rock type RockType1
3 --Sw      Krw      Krow      Pc
40.15      0        0.8        0.59999973
50.2      0.00026726  0.63992059  0.19650725
60.25     0.00173668  0.50166679  0.08267227
70.3     0.00518999  0.3839033  0.04075024
80.35    0.011285   0.28525636  0.02240635
90.4     0.02061382  0.20430839  0.0133462
100.45   0.03372469  0.13959123  0.00845049
110.5    0.05113335  0.08957714  0.00561486
120.55   0.07333014  0.05266665  0.00387911
130.6    0.10078463  0.02717104  0.00276764
140.65   0.13394902  0.01128542  0.00202875
150.7    0.17326059  0.00304238  0.00152177
160.75   0.21914365  0.00021533  0.00116437
170.78   0.25      0        0.001
181      1        0        0
19 /
20 -- Rock type RockType2
21 --Sw      Krw      Krow      Pc
220.3     0        0.8        0.04075024
230.35   0.00091117  0.55784267  0.02240635
240.4    0.00592077  0.3679217  0.0133462
250.45   0.01769394  0.22488787  0.00845049
260.5    0.03847326  0.12311444  0.00561486
270.55   0.07027746  0.05662059  0.00387911
280.6    0.11497553  0.01894646  0.00276764
290.65   0.17432583  0.00291573  0.00202875
300.7    0.25      0        0.00152177
311      1        0        0
32 /
33
34 SGOF
35 -- Rock type RockType1
36 --Sg      Krg      Krog      Pc
370      0        0.8        0
380.1     0        0.48899072  0
390.15   0.00119716  0.3679217  1.582E-05
400.2    0.00777915   0.26769705  0.00023287
410.25   0.02324763  0.18666604  0.00112297
420.3    0.05054908   0.12311444  0.0034288
430.35   0.09233583   0.07525228  0.00815005
440.4    0.15106353   0.04119672  0.01653446
450.45   0.22904244   0.01894646  0.03007108
460.5    0.32846887   0.00633989  0.05048502
470.55   0.45144623   0.00097566  0.07973307
480.6    0.6        0        0.12
49 /
50 -- Rock type RockType2
51 --Sg      Krg      Krog      Pc
520      0        0.8        0
530.65   0.0015811   0.58207387  2.38E-05
```

RE-Studio automatically generate the SWOF and SGOF keywords with the parameters you fulfilled


Here is an example with my RockType1 and RockType2



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Plot your rock type

 All your rock type creation and importation will be available to plot in the right panel called “Quick Plot” of the Porps Builder Module

