

Unconventional Reservoir Engineering Project

Colorado School of Mines

May 3, 2013

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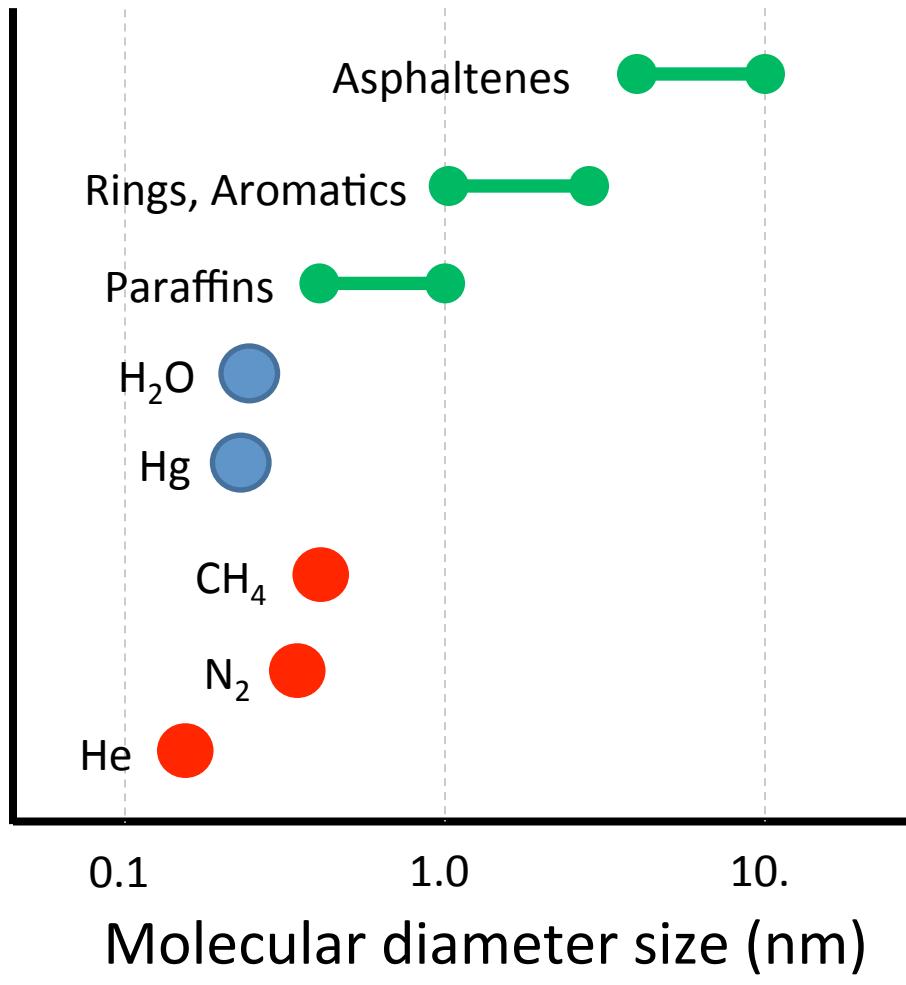
NITEC Research Topics

- Pore Size Related Issues
- Hydrodynamic Equilibrium In Black Oil Simulators
- Development of n-Porosity Simulator

In this presentation,

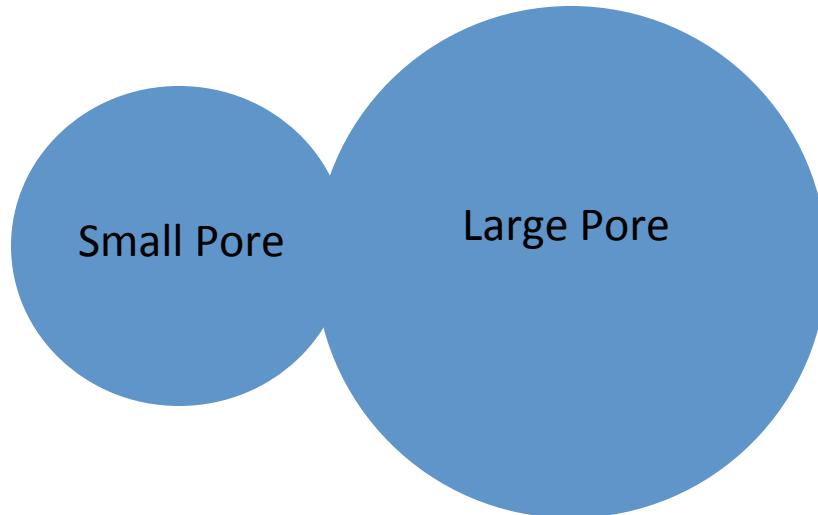
- items that are colored in green show completed work as of May 1, 2013
- Items that are colored in black reflect future work

Pore Size Related Issues



Modified from Nelson (2009)

Pore Size Related Issues



Smaller Molecules

Lower IFT

Higher GOR

Lower Pressure

Higher Surface Forces

Higher Capillary Pressure

Excessive P_c Suppression?

Larger Molecules

Higher IFT

Lower GOR

Higher Pressure

Lower Surface Forces

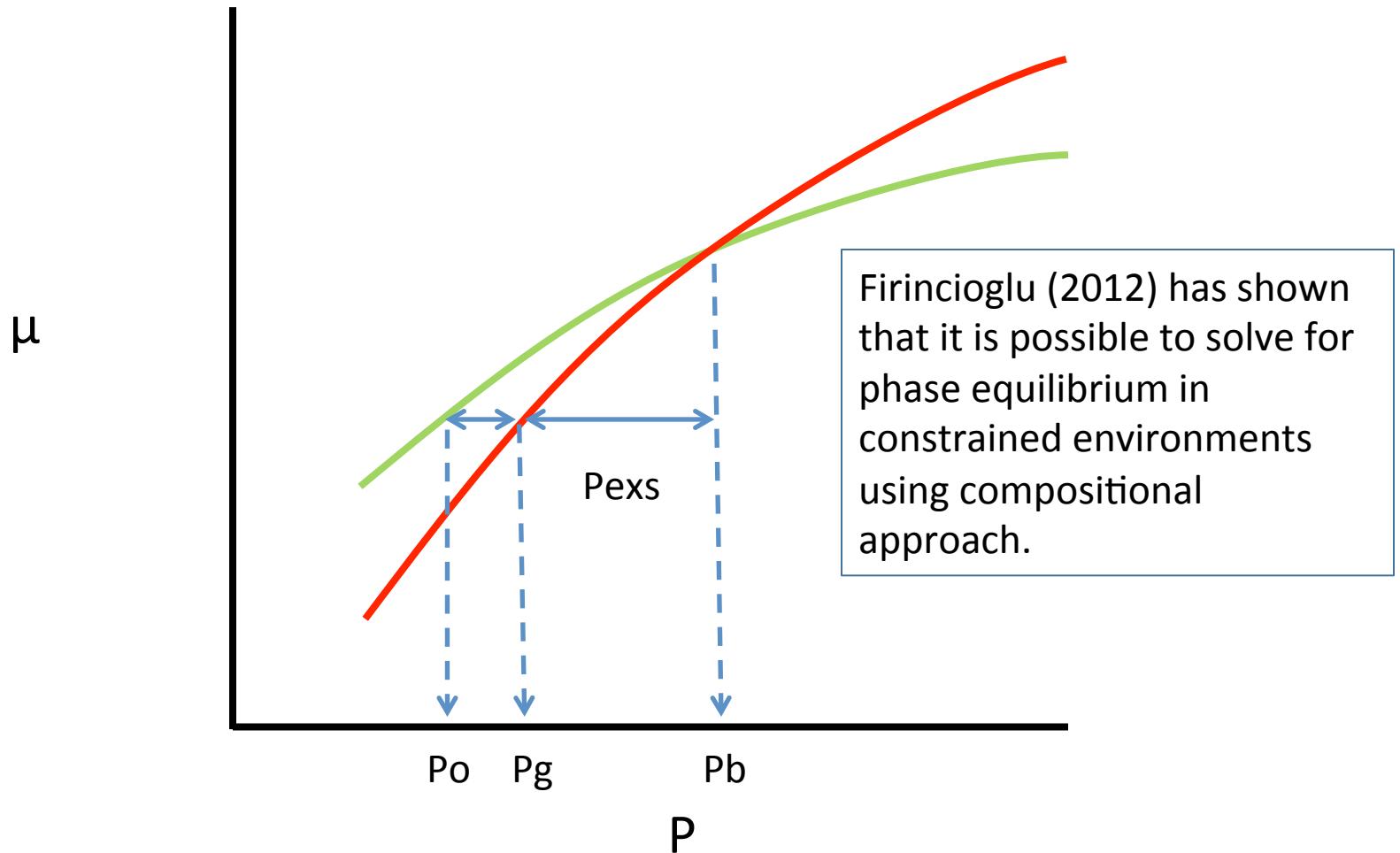
Lower Capillary Pressure

Excessive P_c Suppression?

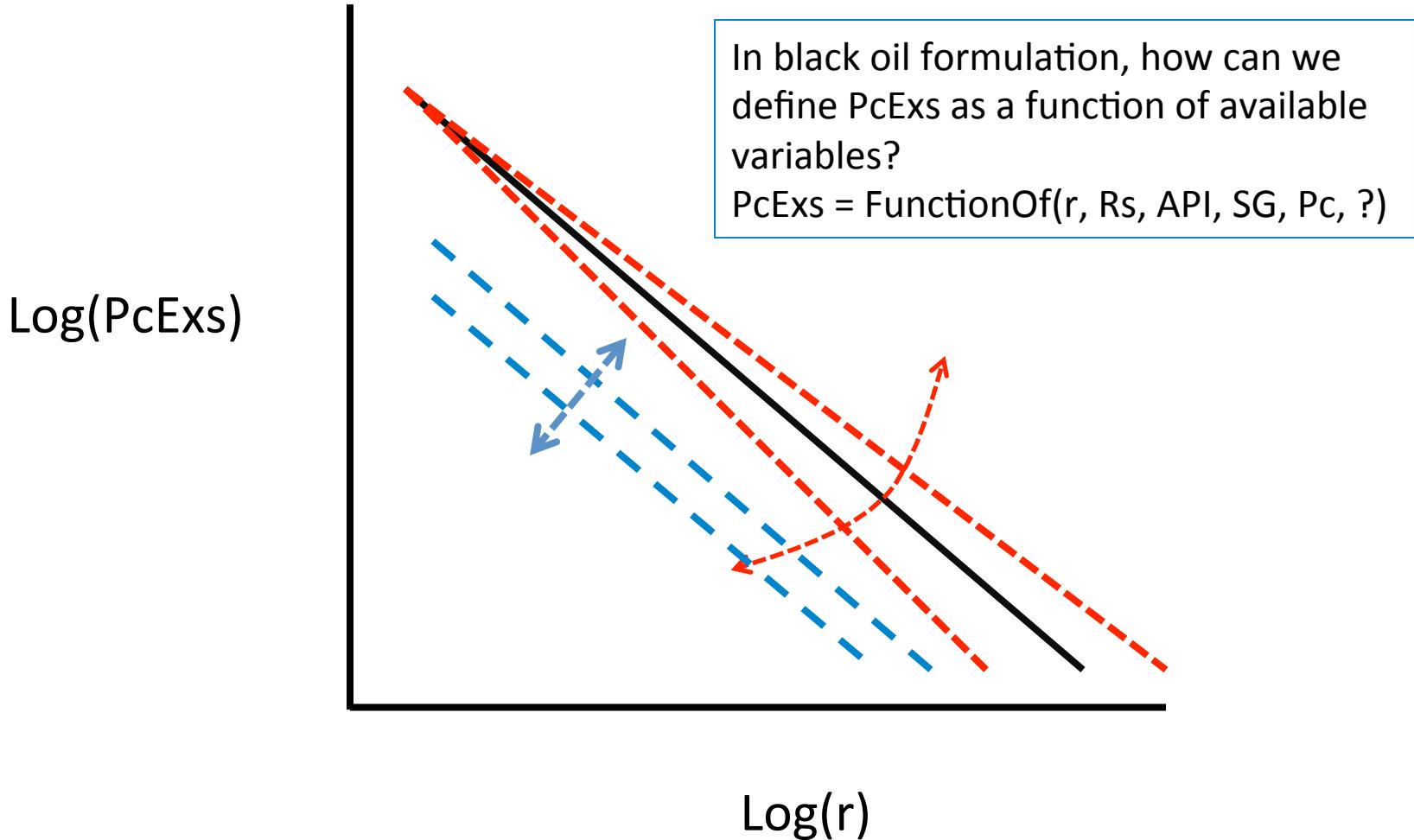
Pore Size Related Issues

- Compositional Variation due to Nano-filtration
 - Laboratory work will commence shortly
- Chemical and Physical Equilibrium
 - Modeling work will commence concurrently
- Diffusion and Osmotic Pressure
 - Diffusion code has been implemented in extended black oil simulator (COZSim) for single-, dual-, and n-porosity formulation
 - Osmotic pressure will be included after the laboratory and modeling work

Constrained Thermodynamics



Excess Suppression



Hydrodynamic Equilibrium

- Excess Capillary Pressure as a function of ?
 - Described by Firincioglu for Compositional Formulation
 - Developed by Firincioglu for Black Oil Formulation as a function of R_s and P_c
 - $(P_{c,exc} + P_c) / P_c = \text{functionOf}(R_s)$

Hydrodynamic Equilibrium

- COZSim Enhancements
 - Updated Initialization Code (static)
 - Calculate Phase Densities and Phase Pressures
 - Calculate Capillary Pressures
 - Estimate R_s (for multi-phase conditions)
 - $P_{cExc} = f(P_c, R_s)$
 - Bulk Pressure = $P_{gas} + P_{cExc}$
 - Iterate until convergence for multi-phase conditions

Hydrodynamic Equilibrium

- COZSim Enhancements
 - Updated Simulation Code (dynamic)
 - Assume P_c and R_s (or obtain from previous solution)
 - Estimate $P_{c,exc}$ and calculate p_{bulk}
 - Perform VLE at p_{bulk}
 - Calculate phase properties at phase pressures
 - Calculate saturations and P_{cs} s from saturations
 - Correct P_{cs} s for IFT and integrate (dynamic pseudo)
 - Iterate until convergence in P_c and R_s

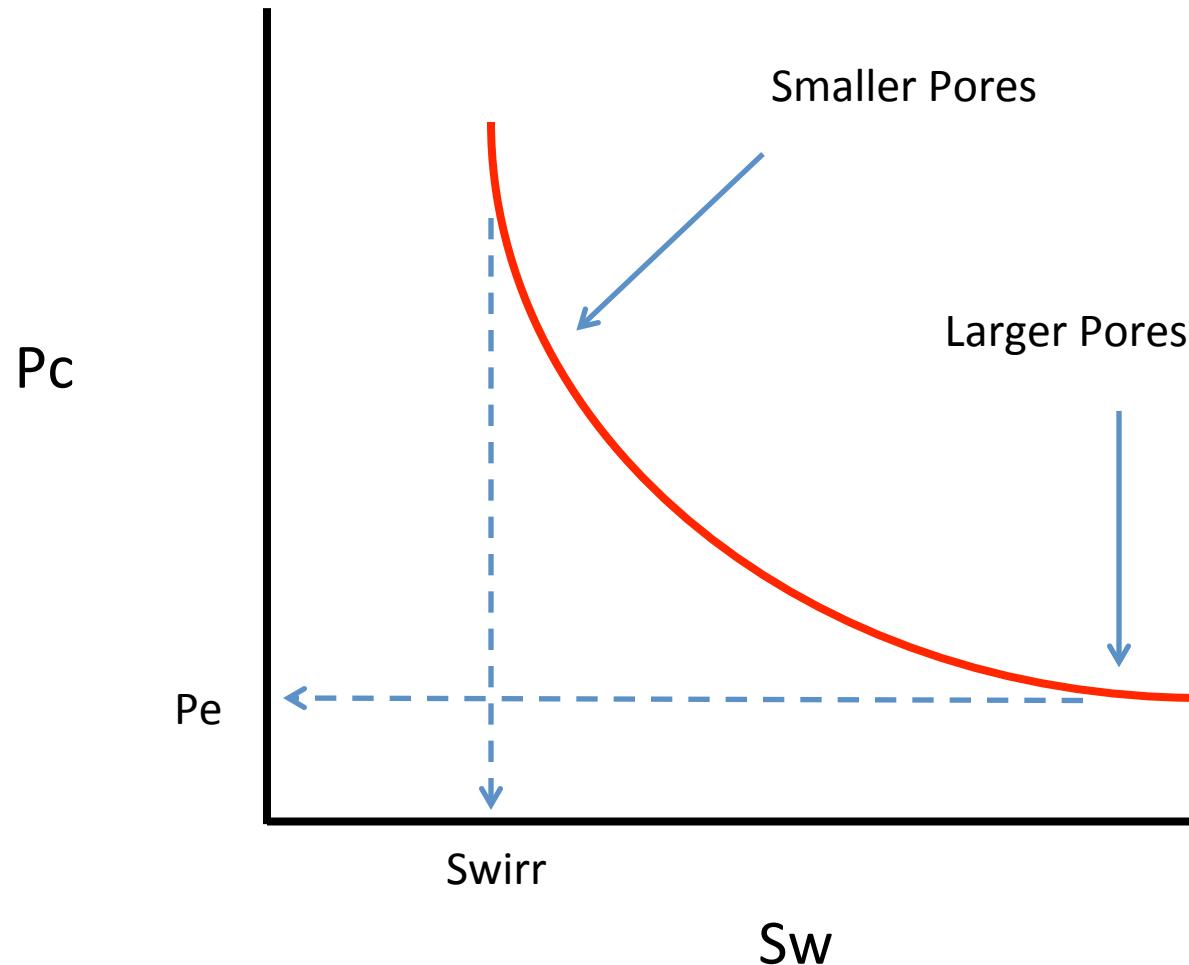
Hydrodynamic Equilibrium

- COZSim Enhancements
 - Other items
 - Dynamic and Static pseudo-Pc (Quandelle)
 - Parachor correlations from Black Oil data
 - Dynamic Interfacial Tension correction for Pc
 - Molecular Diffusion (not tested)

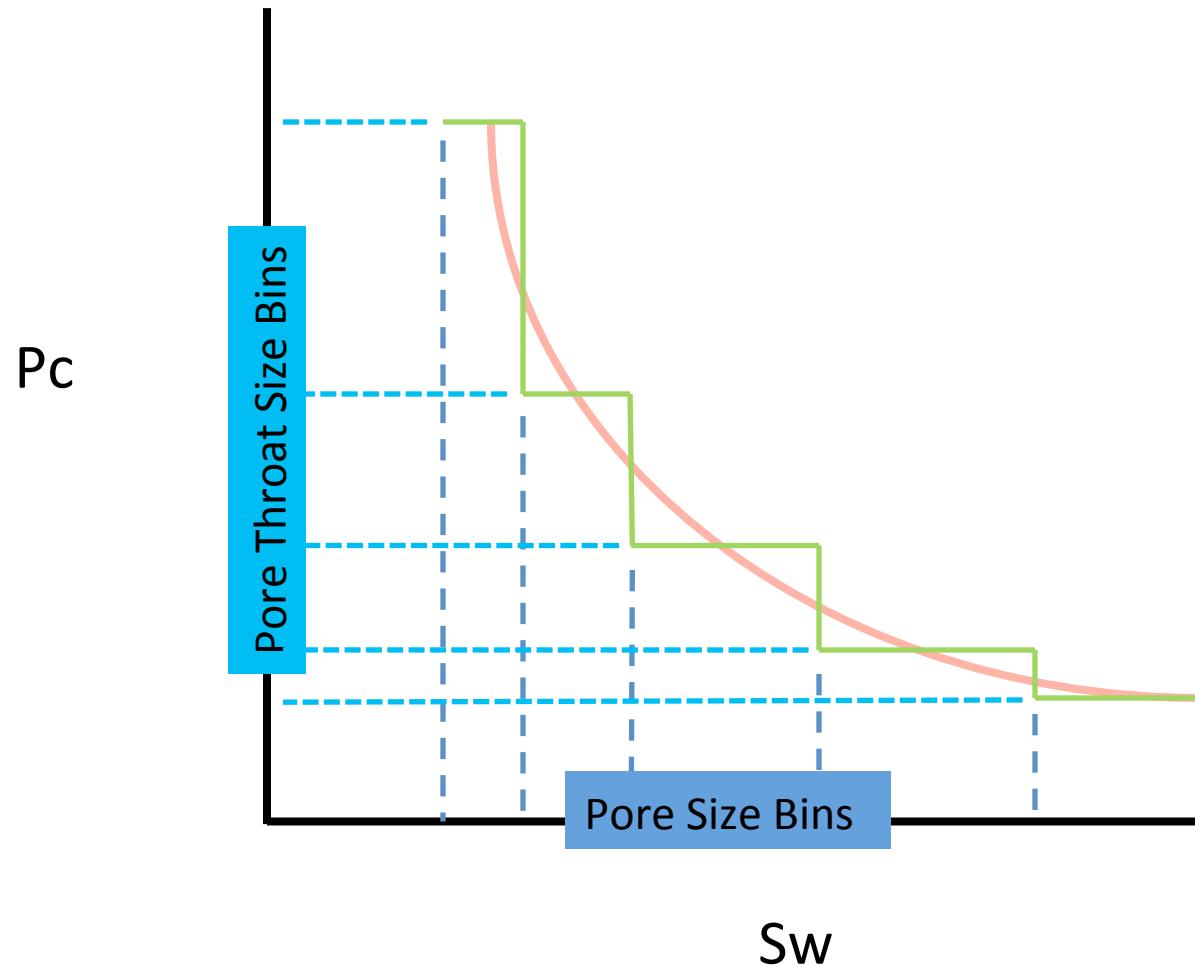
Hydrodynamic Equilibrium

- COZView (Pre-Post Processor) Enhancements
 - Simplified simulation model building for single-porosity cases
 - Post-processing of new arrays

Characteristic P_c Curve



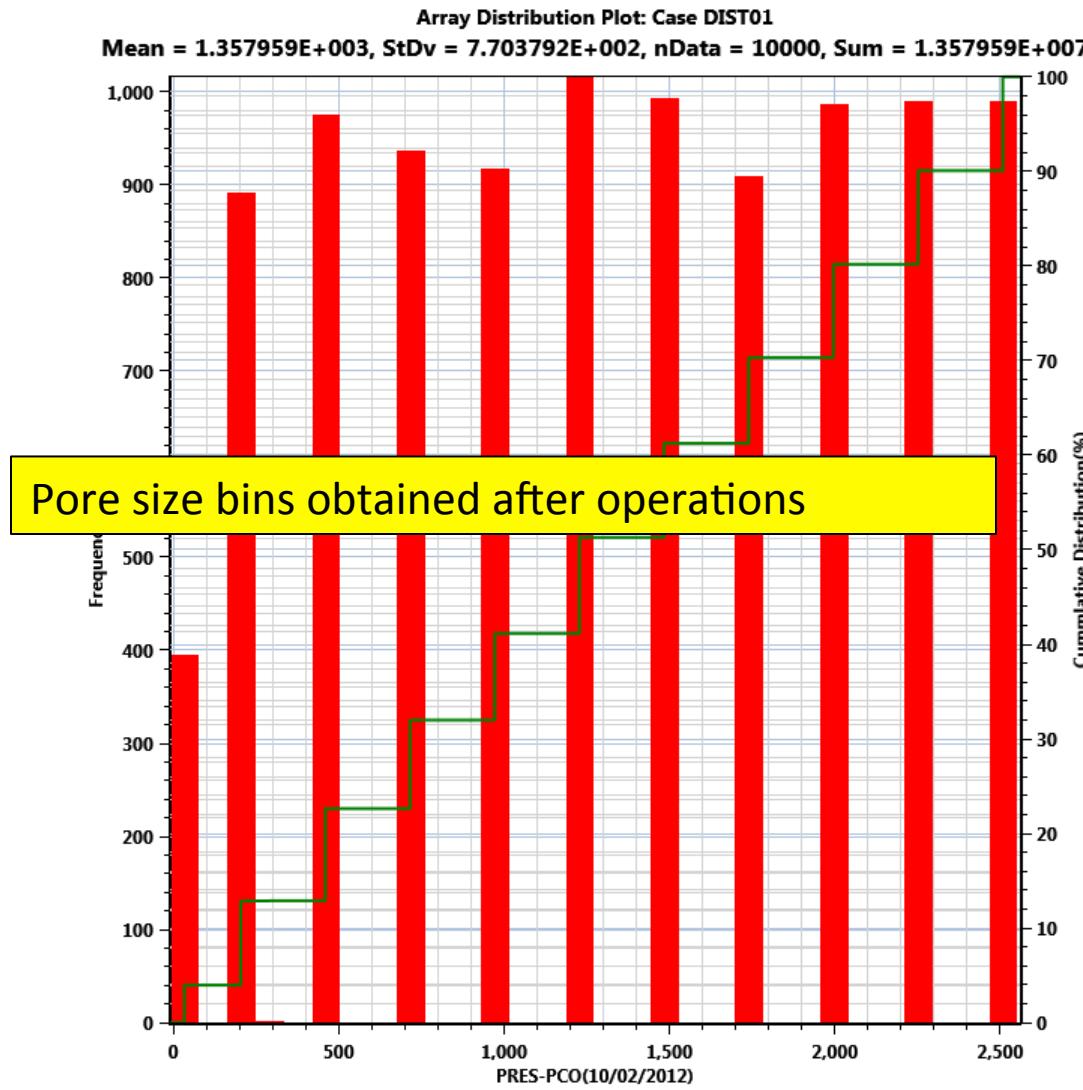
Discretized P_c Curve



Simulation Examples

```
!
POROSITY
10000*0.10          ! set average porosity
!
! for repeatability, set the seed to a value
SETSEED 5000
!
! initialize pore size bin array
USERDEFINED01
10000*5.5
!
! Perturbation to distribute between 0.5 and 10.5
USERDEFINED01 EQUALS PERTURB 5.0 LINEAR
!
! get to nearest integer
USERDEFINED01 EQUALS NINT USERDEFINED01
!
! Now, we have a pore size bin array from 1 to 10 (integers)
!
```

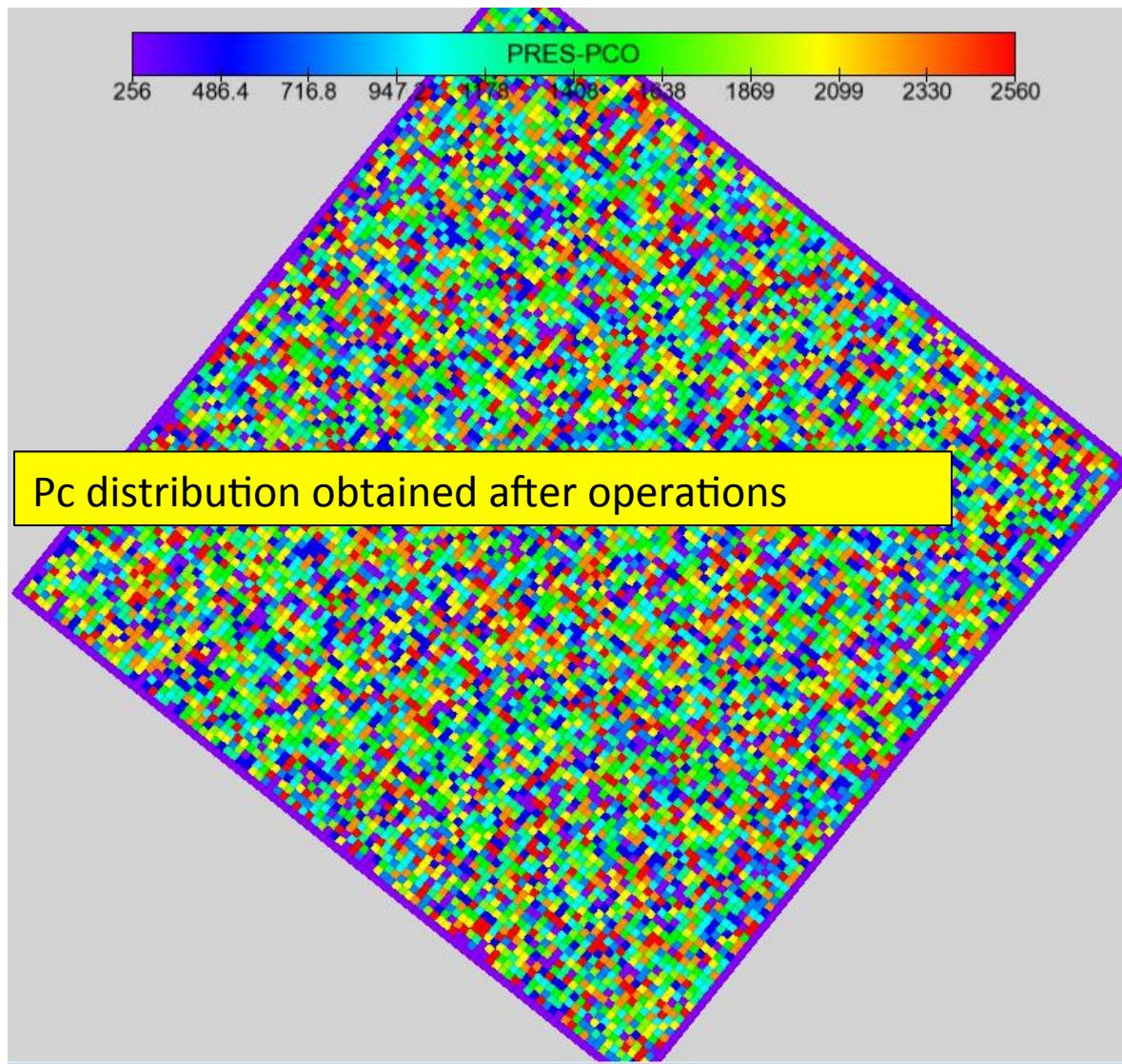
Simulation Examples



Simulation Examples

```
!
! initialize pore size array
USERDEFINED02
10000*1.0
!
! set the pore size for each bin
USERDEFINED02 EQUALS DIVIDEBY USERDEFINED01
! Now, we have a normalized pore size array between 0.1 and 1.0
!
! initialize PcMult array to be Pe
PCMULTGO
10000*256.0
!
! finalize multiplier
PCMULTGO EQUALS DIVIDEBY USERDEFINED02
!
! Now, we have Pc (or PcMult) for each block
!
```

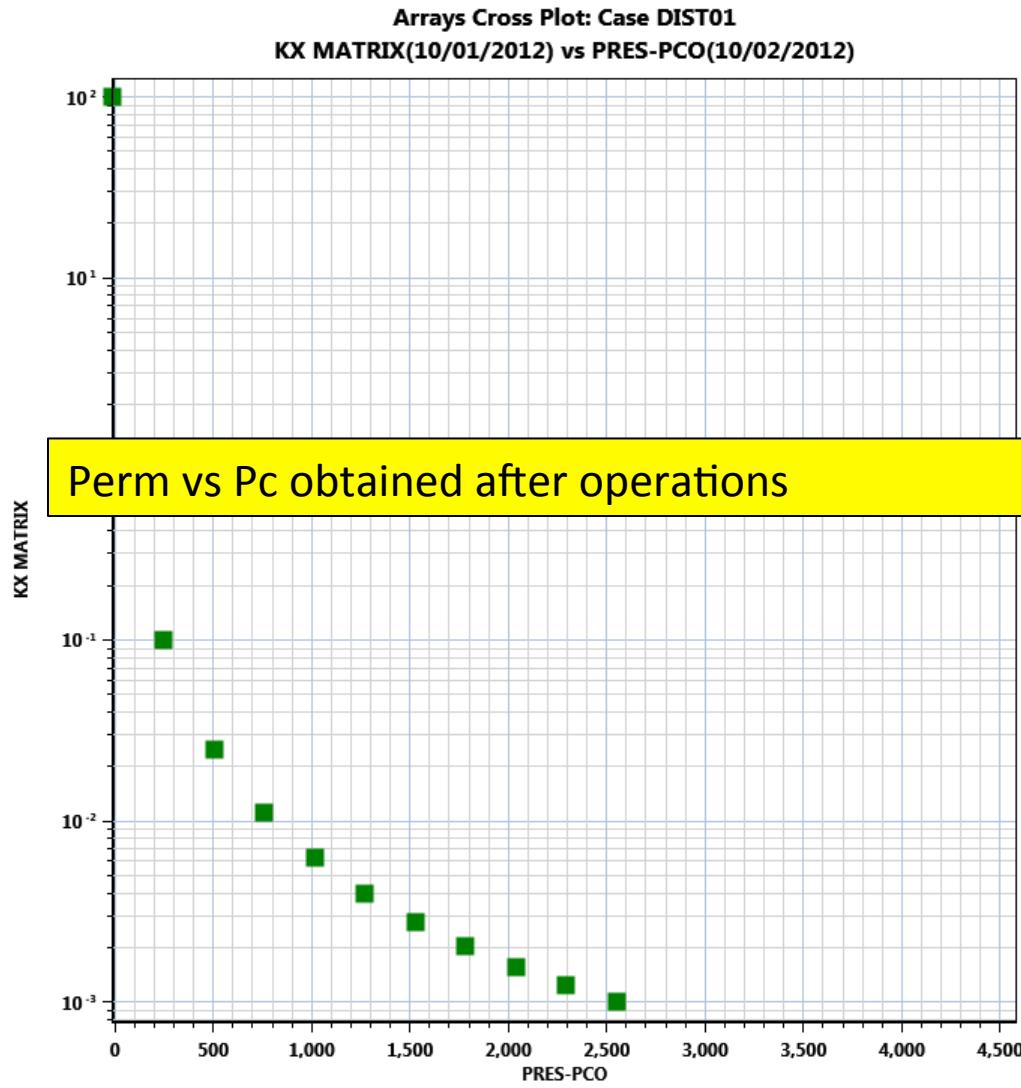
Simulation Examples



Simulation Examples

```
!
! initialize X permeability array
PERMEABILITY-X
10000*0.100
!
! Set permeability proportional to pore size (squared)
PERMEABILITY-X EQUALS TIMES USERDEFINED02
PERMEABILITY-X EQUALS TIMES USERDEFINED02
!
! initialize remaining permeability arrays
PERMEABILITY-Y
10000*1
PERMEABILITY-Z
10000*1
!
! set equal to KX
PERMEABILITY-Y EQUALS PERMEABILITY-X
PERMEABILITY-Z EQUALS PERMEABILITY-X
!
! Now, we have permeability for each block
!
```

Simulation Examples



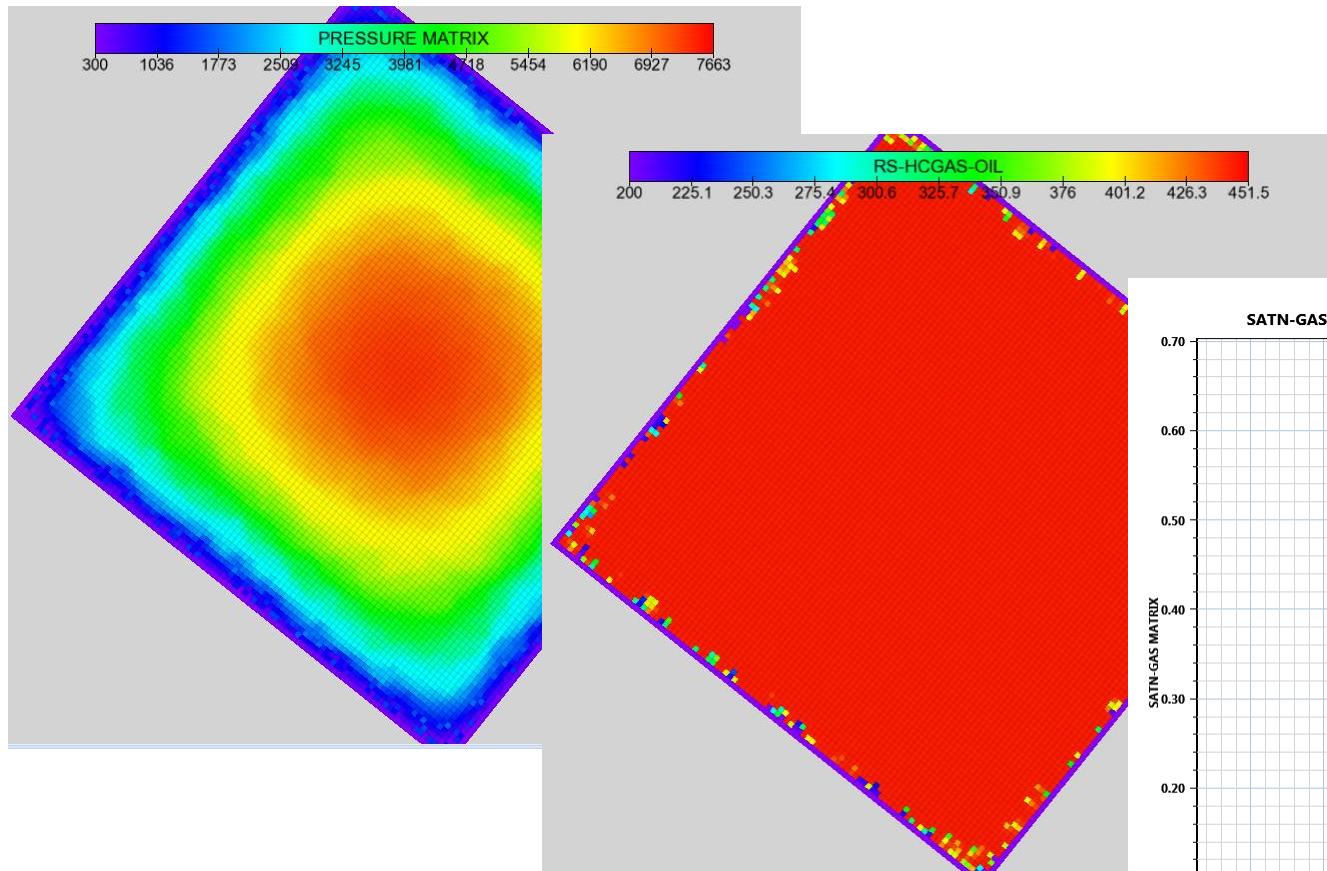
Simulation Examples

```
!
! initialize water film thickness array (half of 0.1 - which is the minimum of USERDEFINED02)
USERDEFINED03
10000*0.05
!
USERDEFINED03 EQUALS TIMES    -1.0
USERDEFINED03 EQUALS PLUS     USERDEFINED02
USERDEFINED03 EQUALS DIVIDEBY USERDEFINED02
USERDEFINED03 EQUALS POWER    3.0
!
! swcrit update
SWCRIT EQUALS 1.0
SWCRIT EQUALS MINUS USERDEFINED03
!
! avoid illogical values
IF SWCRIT GT 0.95
    SWCRIT EQUALS 0.95
ENDIF
!
SWCONN EQUALS SWCRIT
!
```

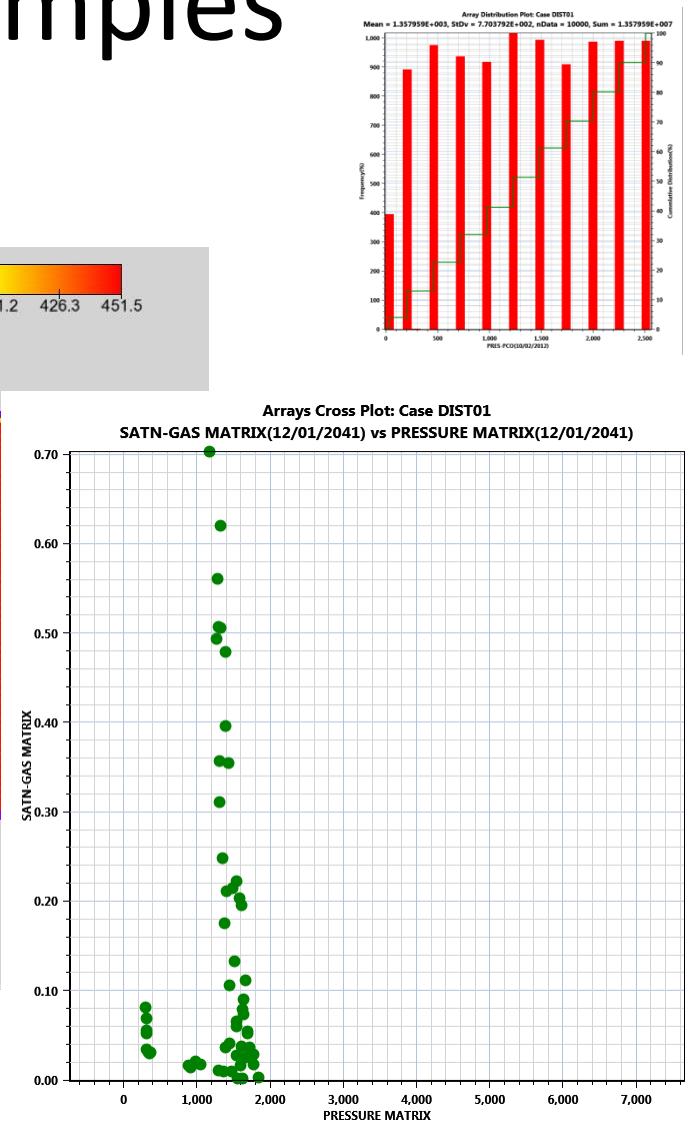
Simulation Examples

```
!
USE-EXCESSPC 1                                ! to turn on excess pressure calculation
!
! MODIFY THE FRACTURES
!
BOX-SETUP
1 100 1 1 1 1
PERMEABILITY-X EQUALS 100      !00
PERMEABILITY-Y EQUALS 100      !00
SWCRIT      EQUALS 0.02
SORW        EQUALS 0.02
SORG        EQUALS 0.02
SGR         EQUALS 0.02
SWCONN      EQUALS 0.02
SLCONN      EQUALS 0.04
BOX-CLEAR
!
BOX-SETUP
1 100 1 1 1 1
PCMULTGO    EQUALS TIMES 0.0
POROSITY     EQUALS TIMES 0.25
COMPRESS-PV  EQUALS TIMES 10.0
COMPRESS-KX  EQUALS TIMES 10.0
COMPRESS-KY  EQUALS TIMES 10.0
COMPRESS-KZ  EQUALS TIMES 10.0
BOX-CLEAR
!
```

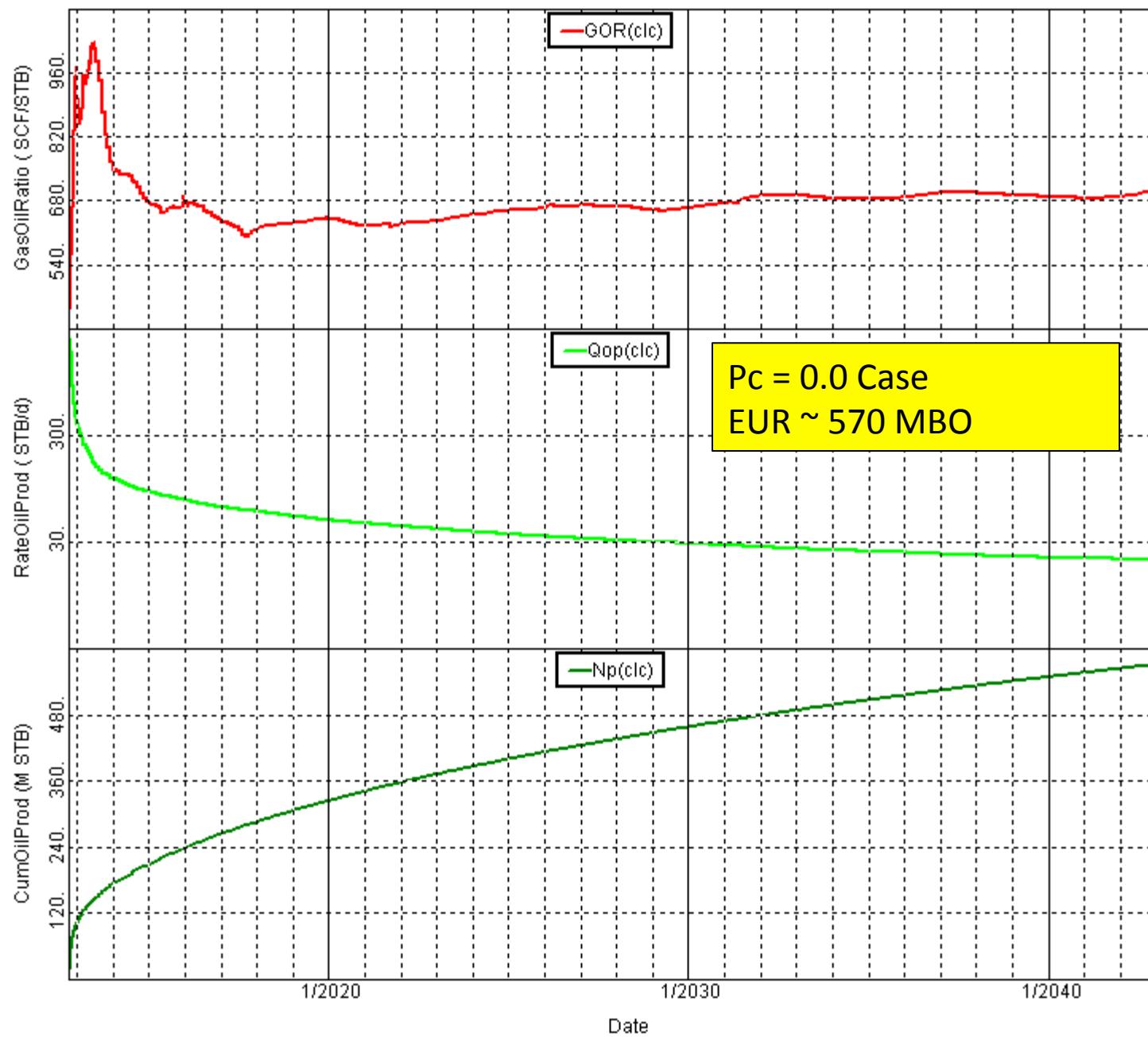
Simulation Examples

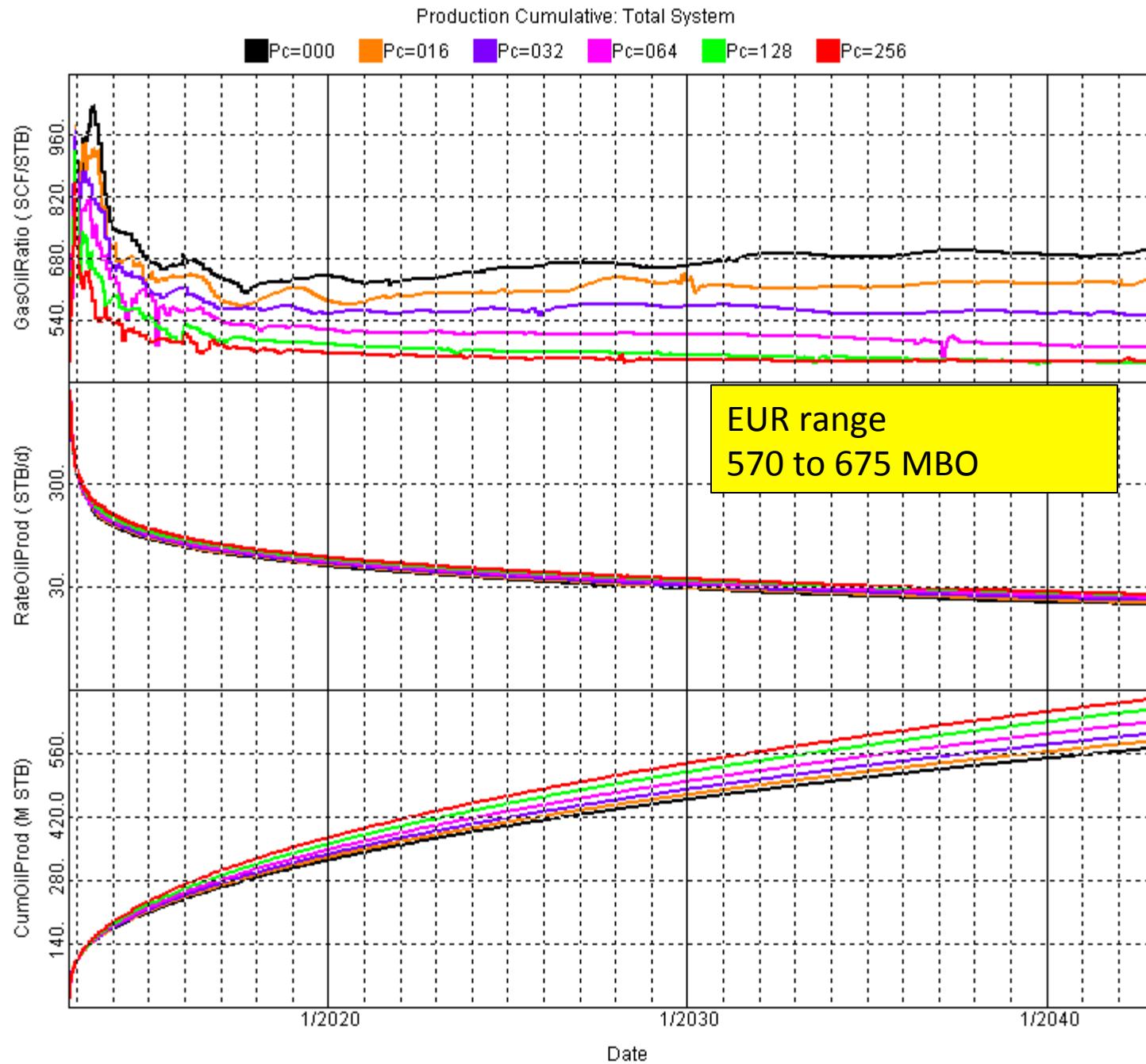


Pe=256 case after 30 years production



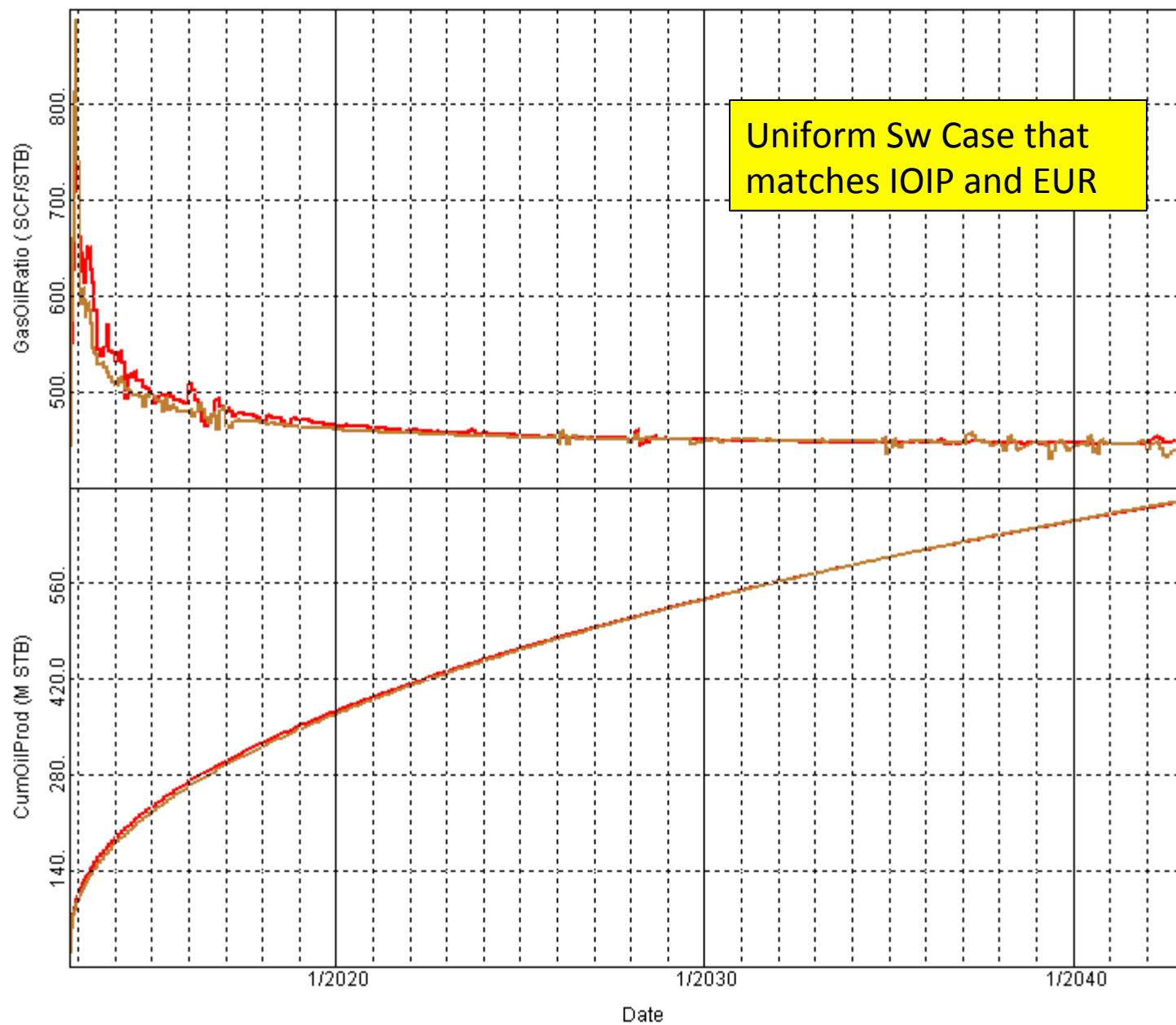
Production Cumulative: Total System





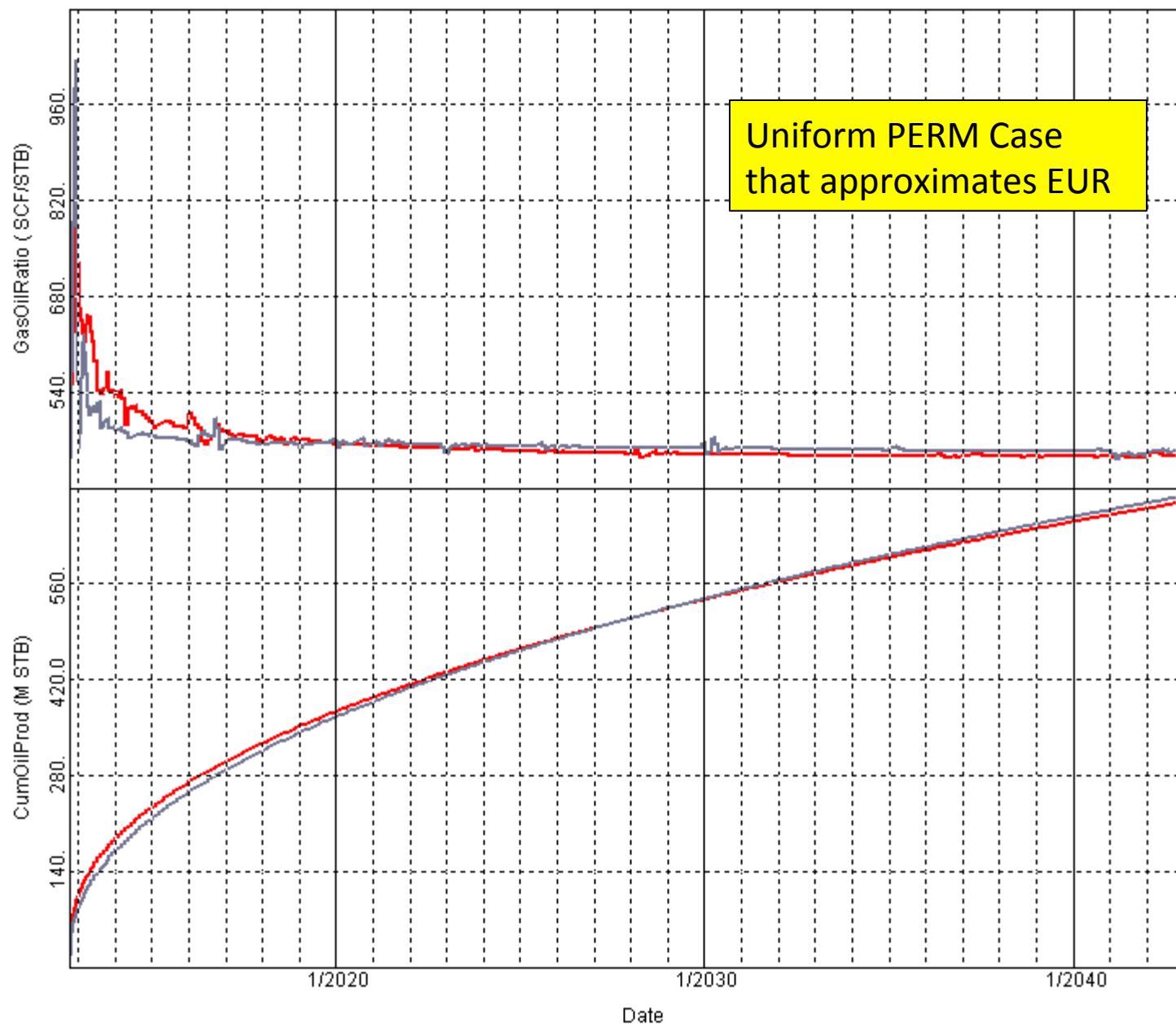
Production Cumulative: Total System

Pc=256 SwAvg



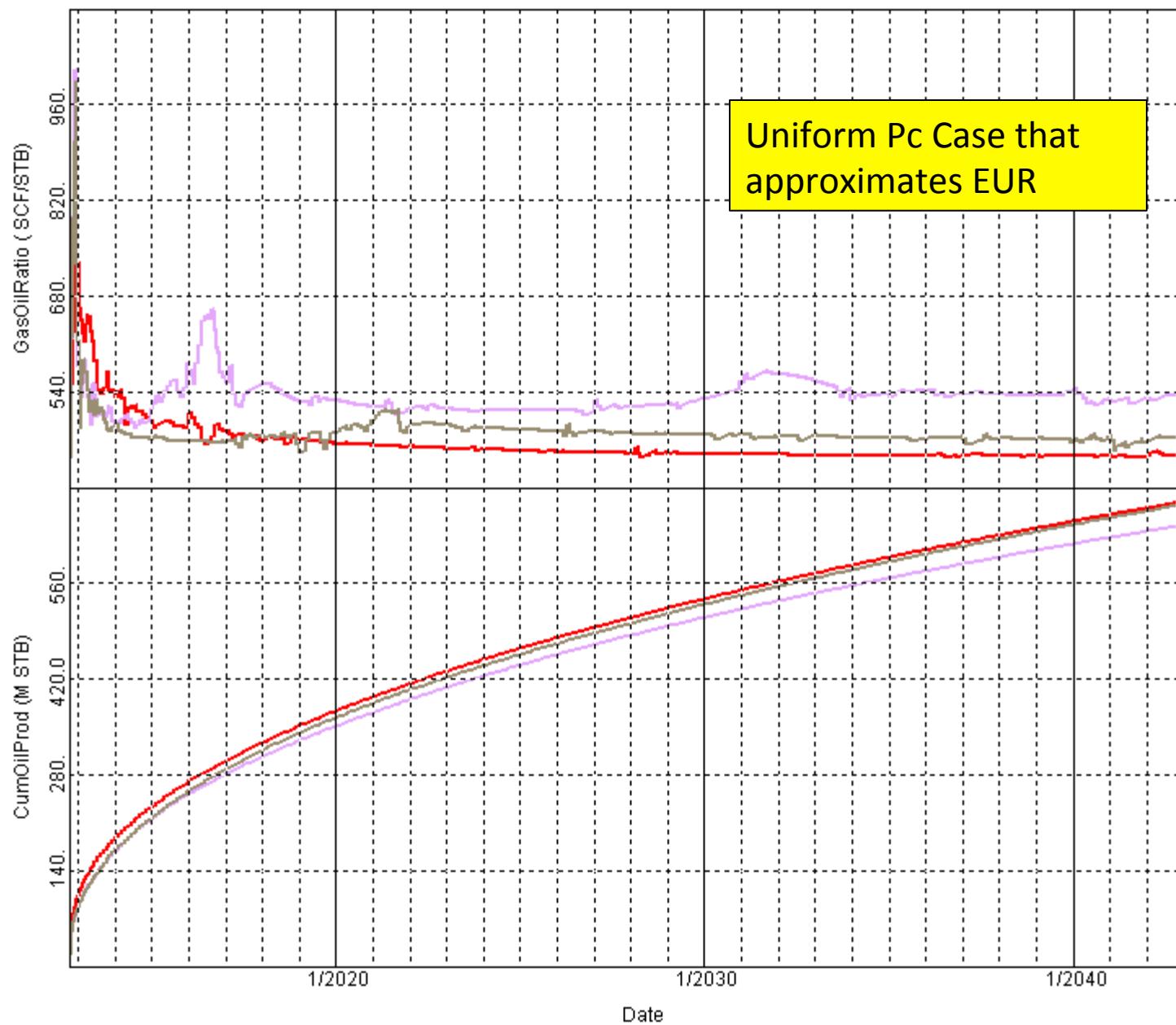
Production Cumulative: Total System

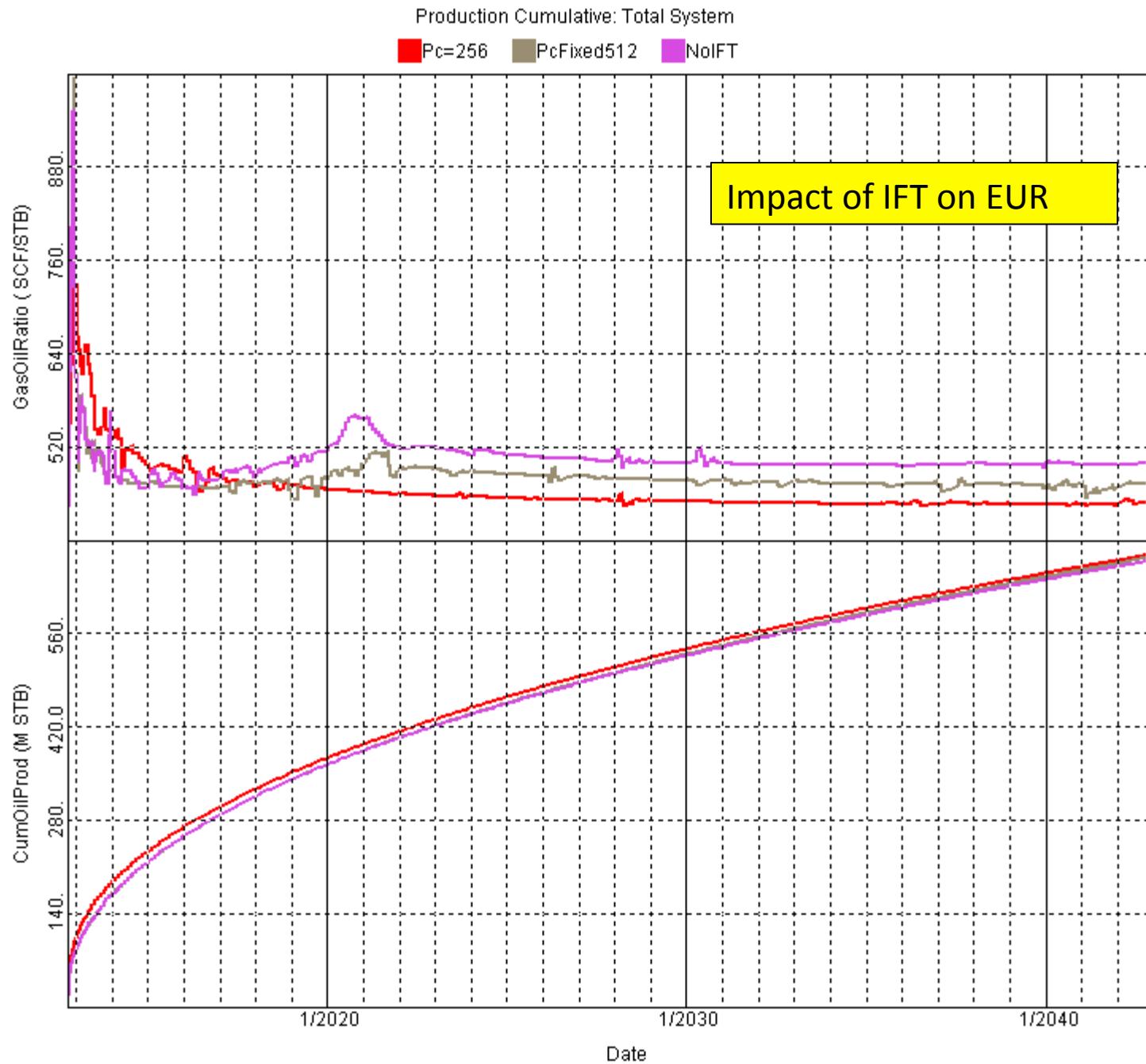
Pc=256 PermAvg



Production Cumulative: Total System

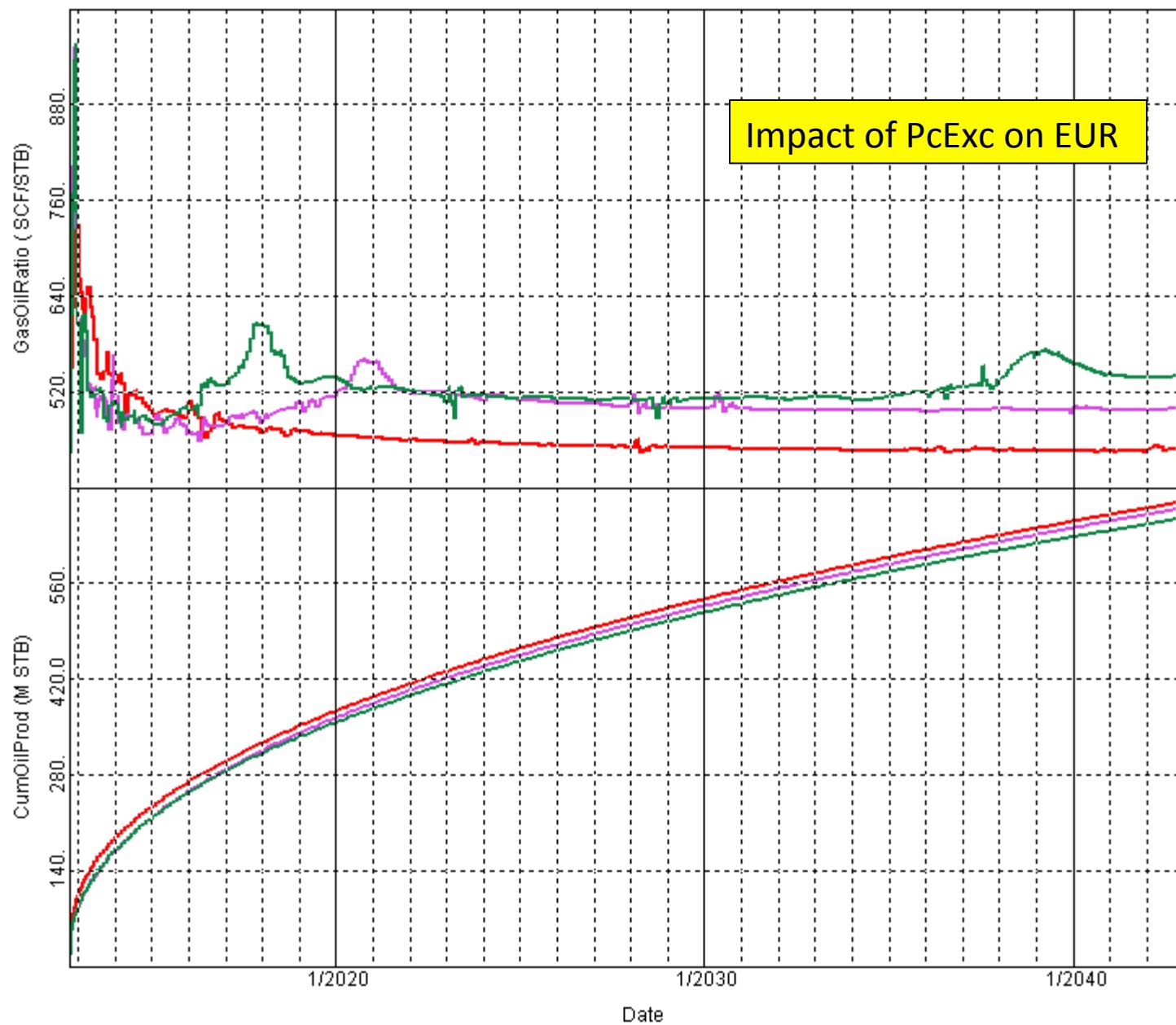
Pc=256 PcFixed256 PcFixed512

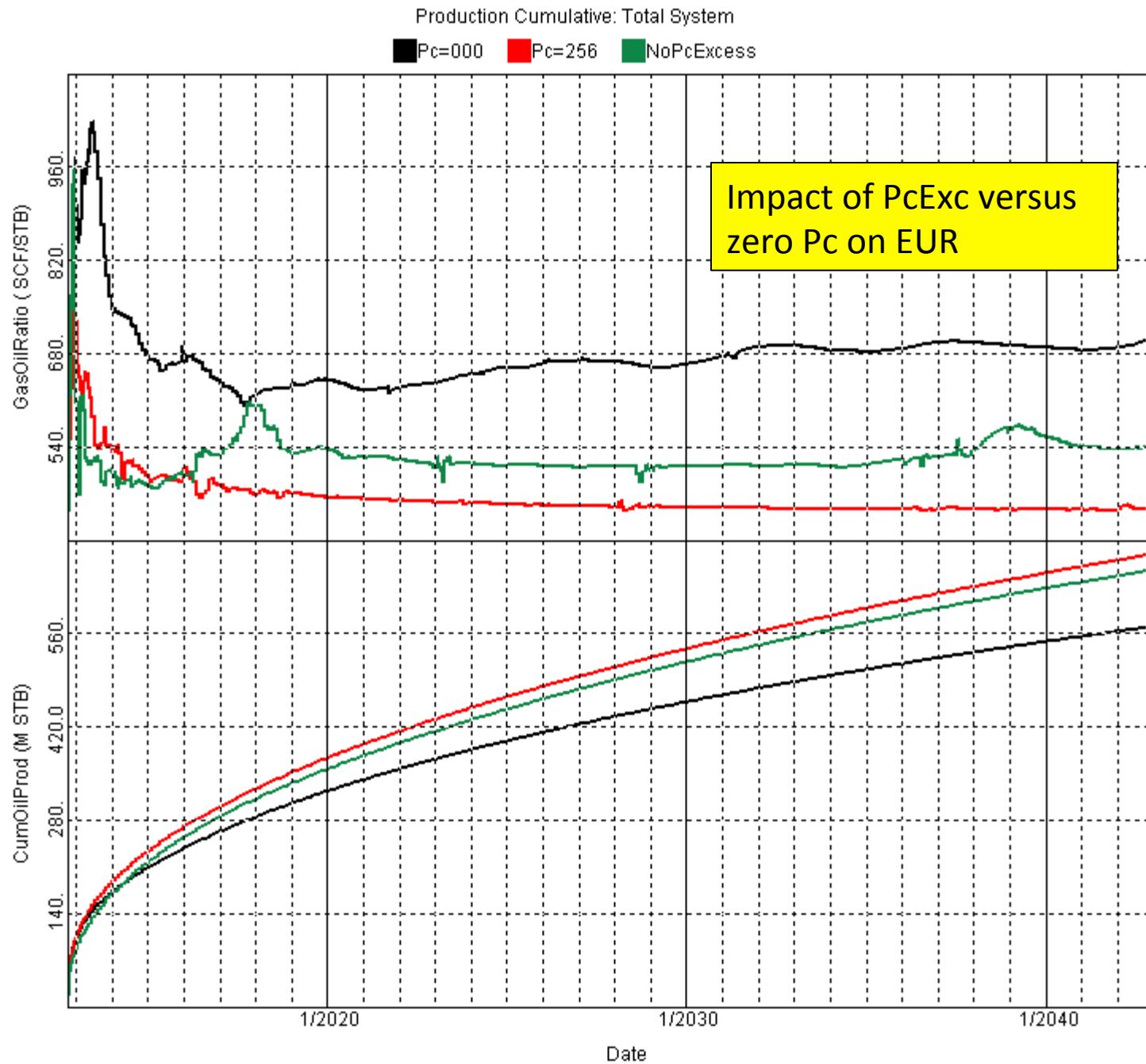




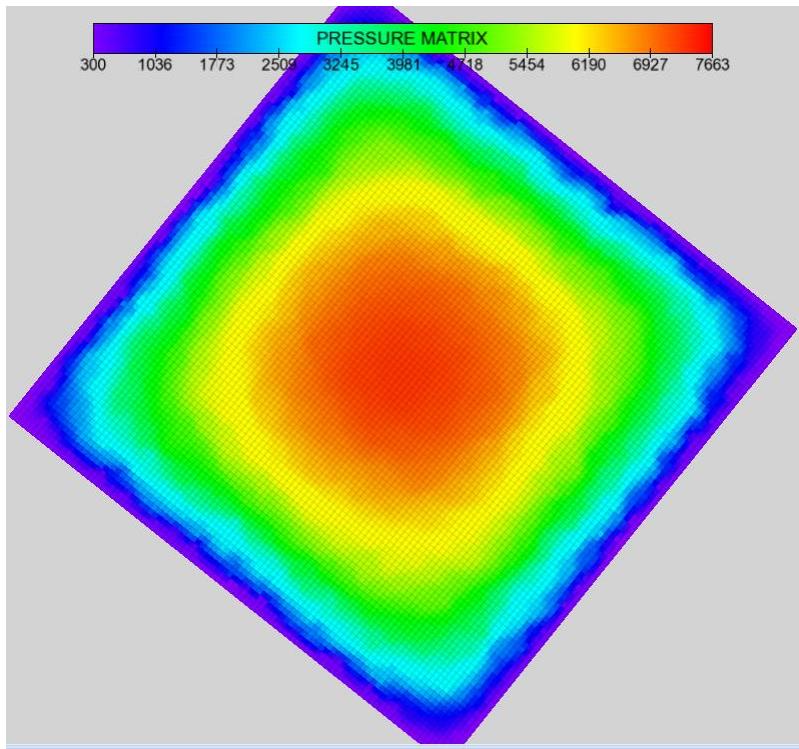
Production Cumulative: Total System

Pc=256 NoIFT NoPcExcess

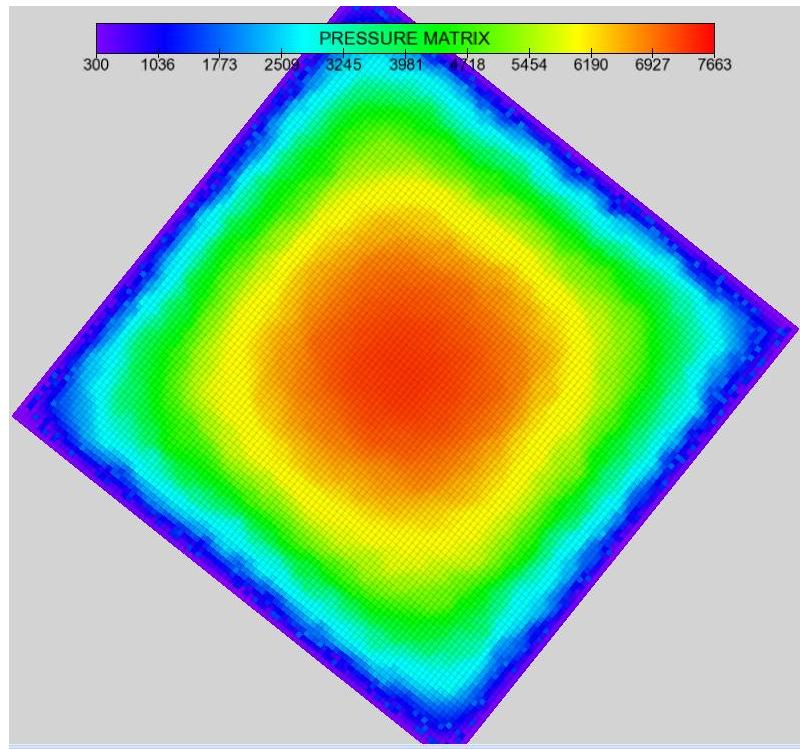




Simulation Examples

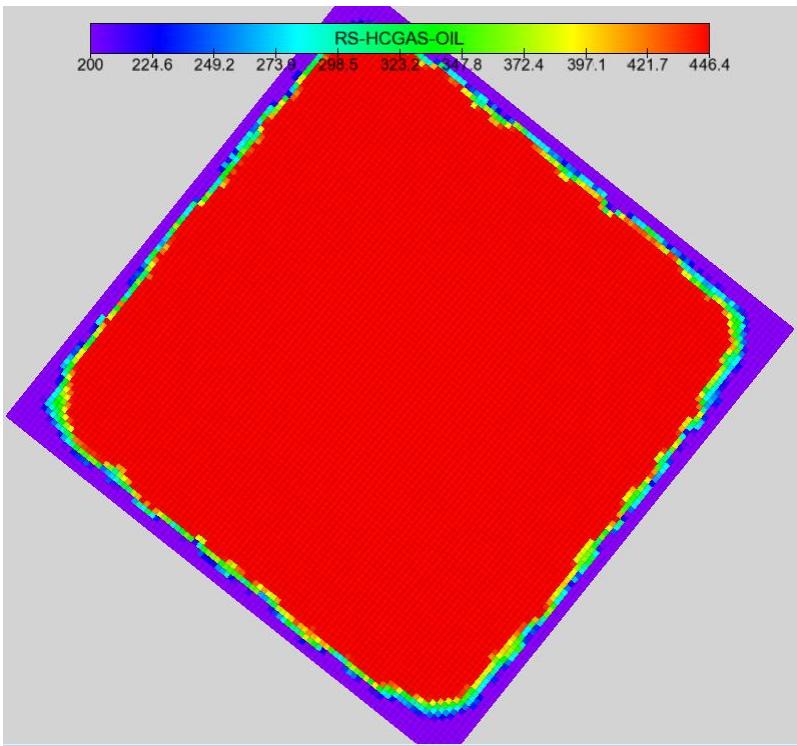


Pc=0.0 case after 30 years production

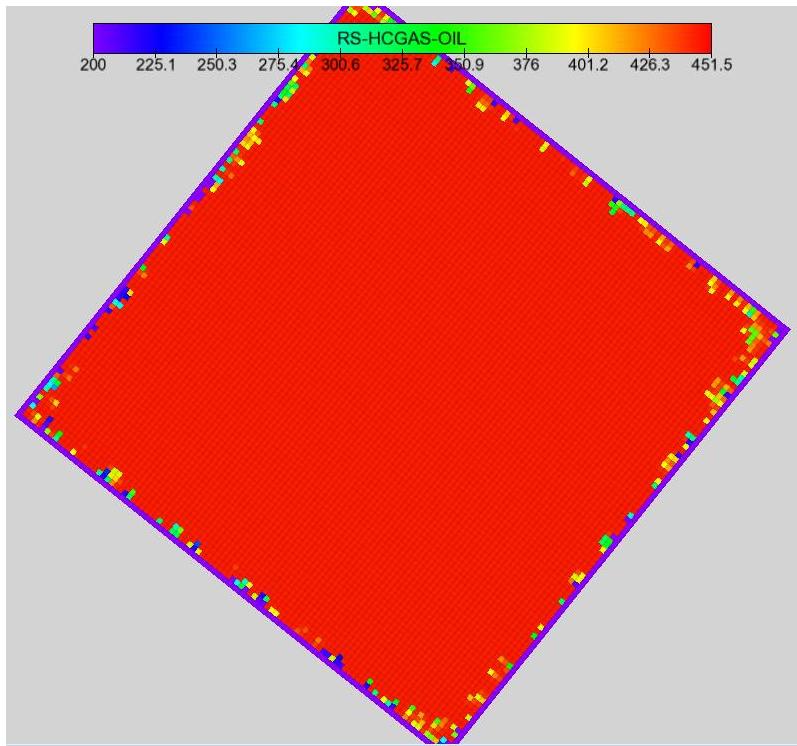


Pe=256 case after 30 years production

Simulation Examples



Pc=0.0 case after 30 years production



Pe=256 case after 30 years production

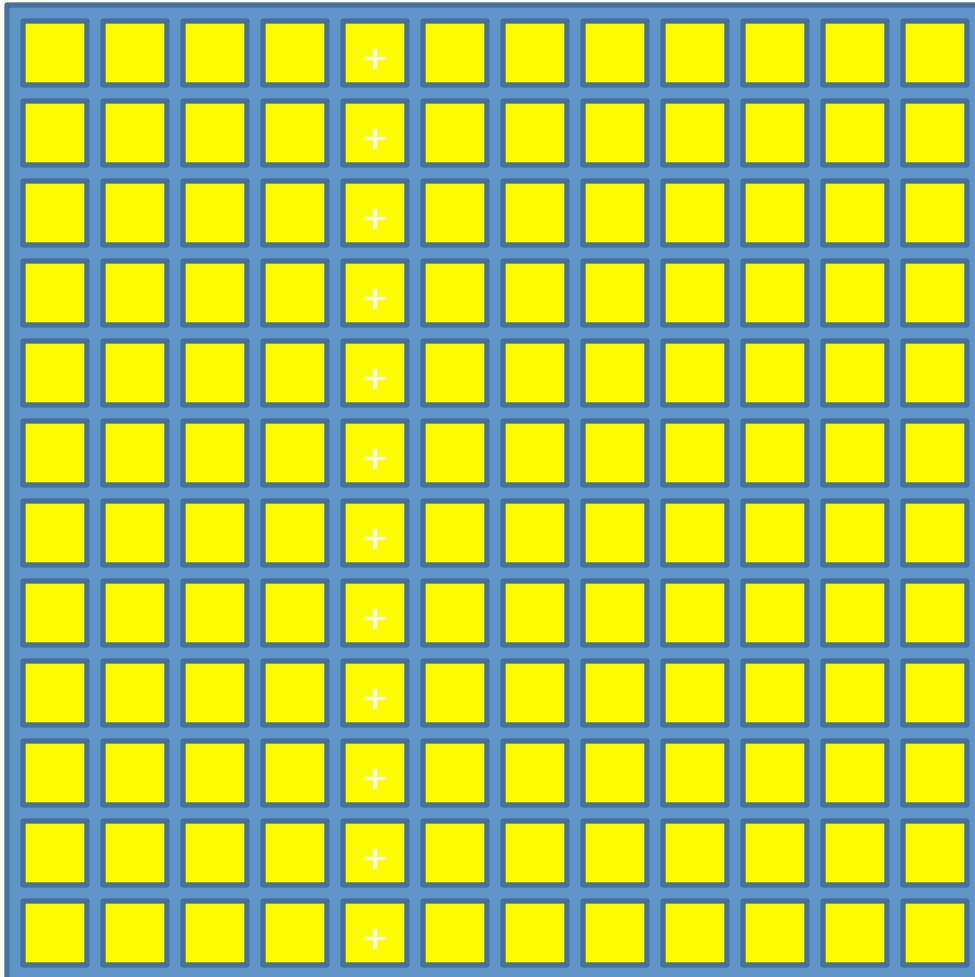
n-Porosity Simulator

- COZSim Enhancements
 - Memory allocation, new arrays
 - Dual-Porosity formulation (not tested)
 - n-Porosity formulation (not tested)
 - Automated pore size bin creation based on P_c
 - Connectivity map of pore size bins

n-Porosity Simulator

- COZView Enhancements
 - Simplified simulation model building for dual-, and n-Porosity models
 - Memory allocation, database issues for dual-porosity models
 - Post-processing for dual-Porosity models (not tested)
 - Memory allocation, database issues for n-Porosity models
 - Post-processing for n-Porosity models

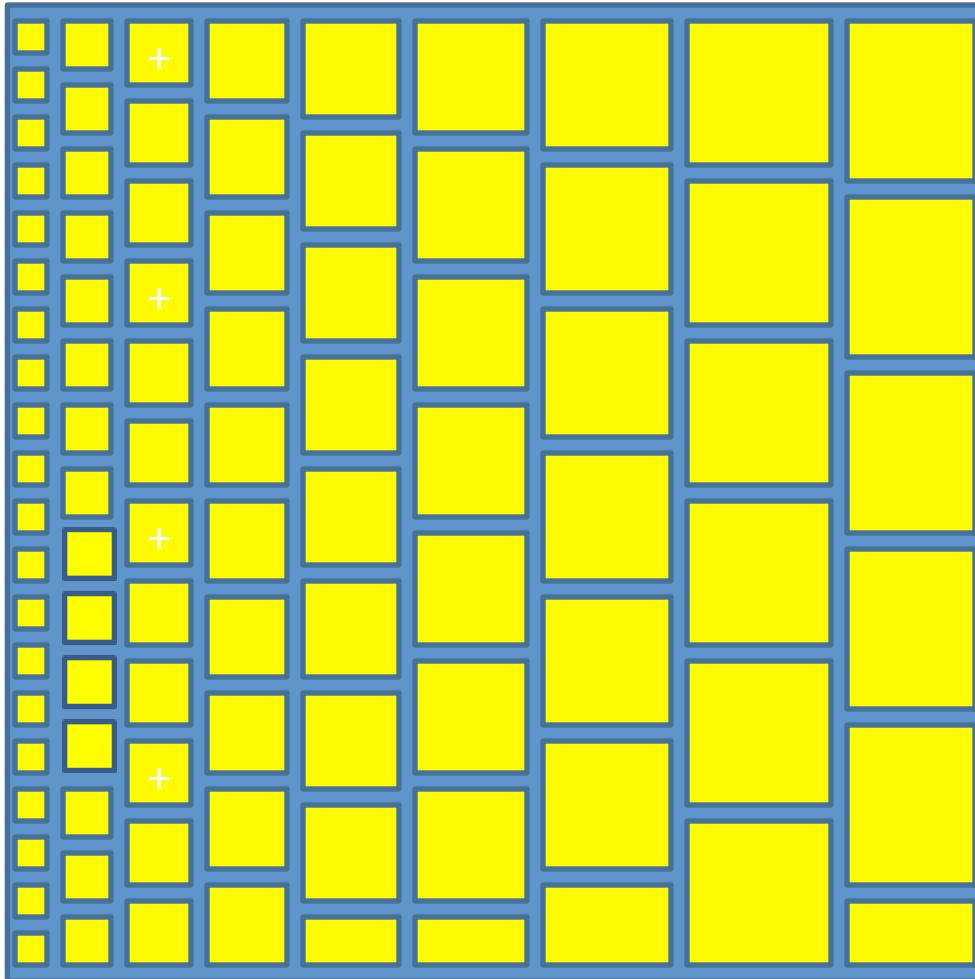
Dual-Porosity Systems



In existing simulators, dual porosity formulation assumes that ALL of the matrix can be represented as uniform blocks in continuum.

This formulation assumes that all of these uniform blocks have the same pore size distribution and pore throat size distribution, and can be represented using a single saturation value and a single capillary pressure value.

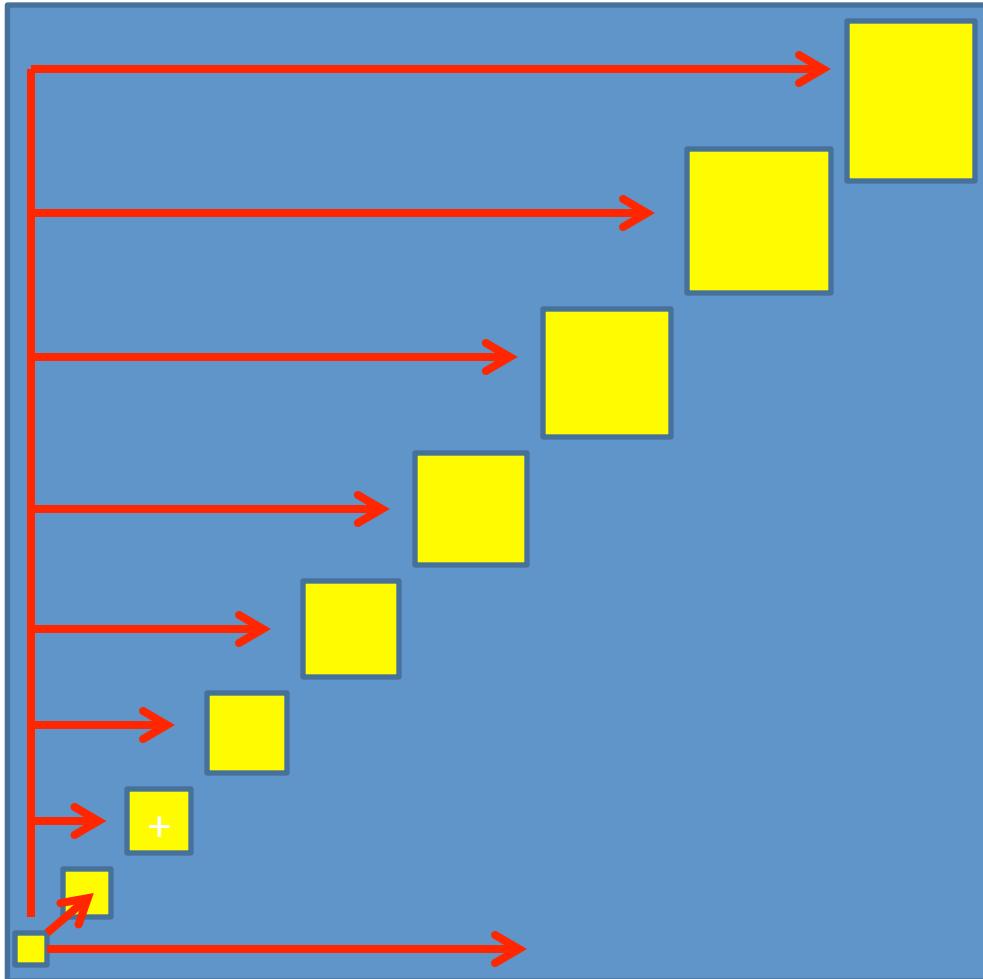
Multiple-Porosity Systems



Instead of assuming uniformly distributed matrix, what if we distribute the pores and the pore throats based on capillary pressure function?

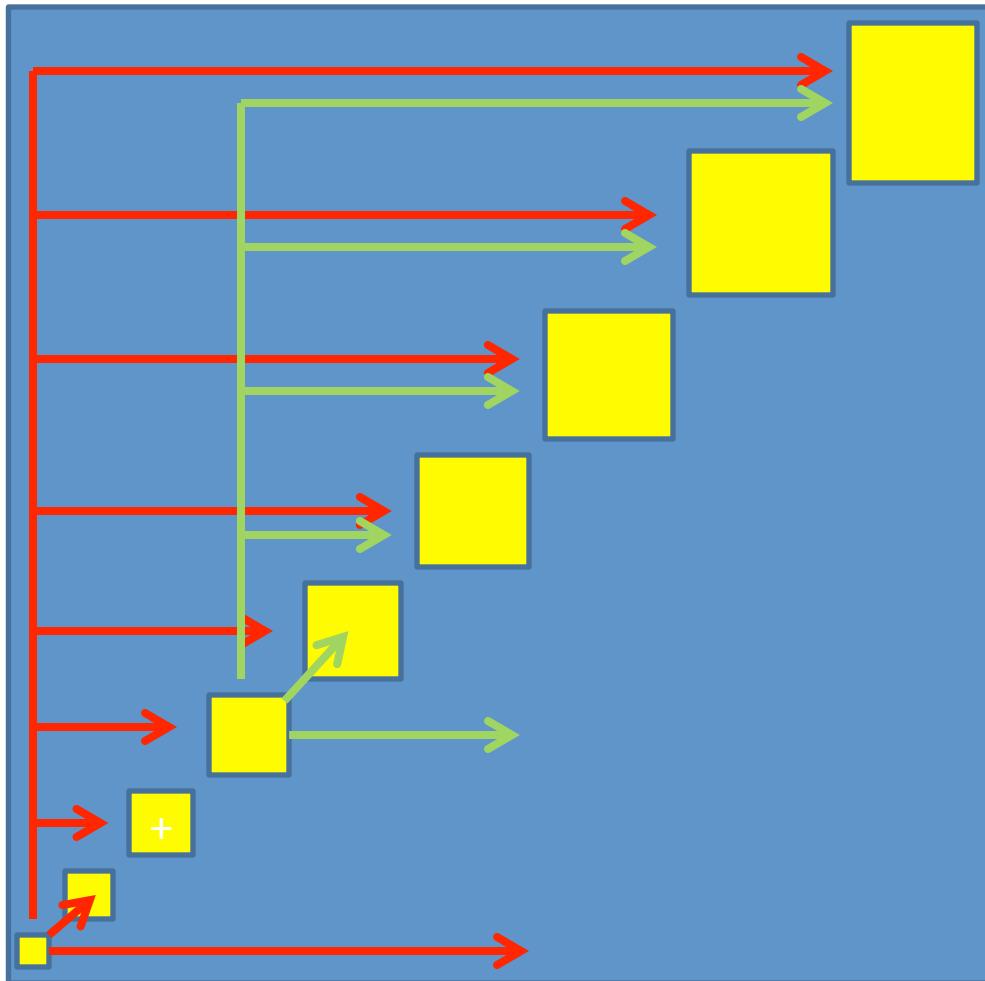
We should have many matrix blocks that represent small pores and few matrix blocks that represent large pores.

Numerical Solution



The new distribution will require the smallest pores to be connected to all of the larger pores, as well as to the natural fractures (or to the largest pores that dominate the flow).

Numerical Solution



Same connectivity logic applies to all of the intermediate pore sizes.

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Thank You