

Generalized network modelling of two-phase flow

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Outline

□ Background

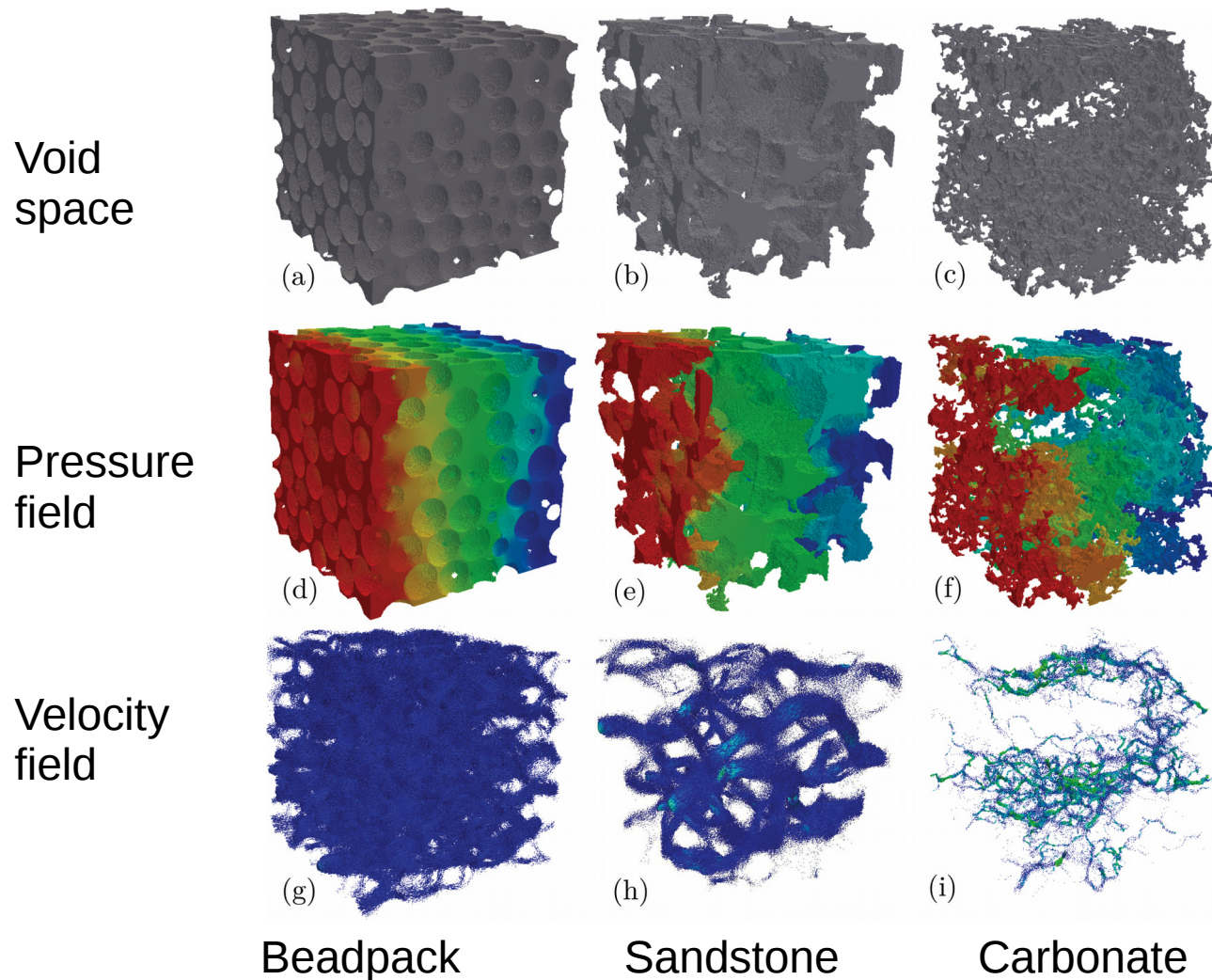
- Direct simulation and conventional network modelling

□ Generalized network modelling

- Network extraction
- Flow simulation
- Validation

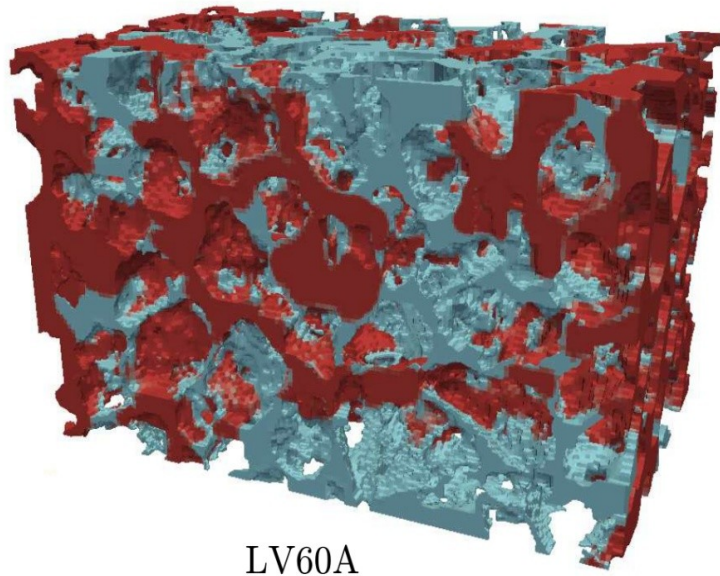
□ Conclusion and future work

Single-phase flow computations

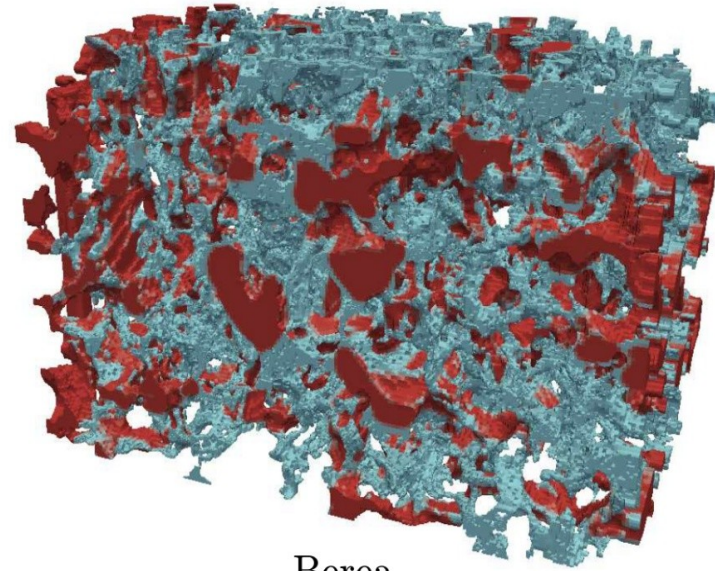


Simulation time: 12 hours on a single workstation (16 processors)

Direct simulations are computationally expensive



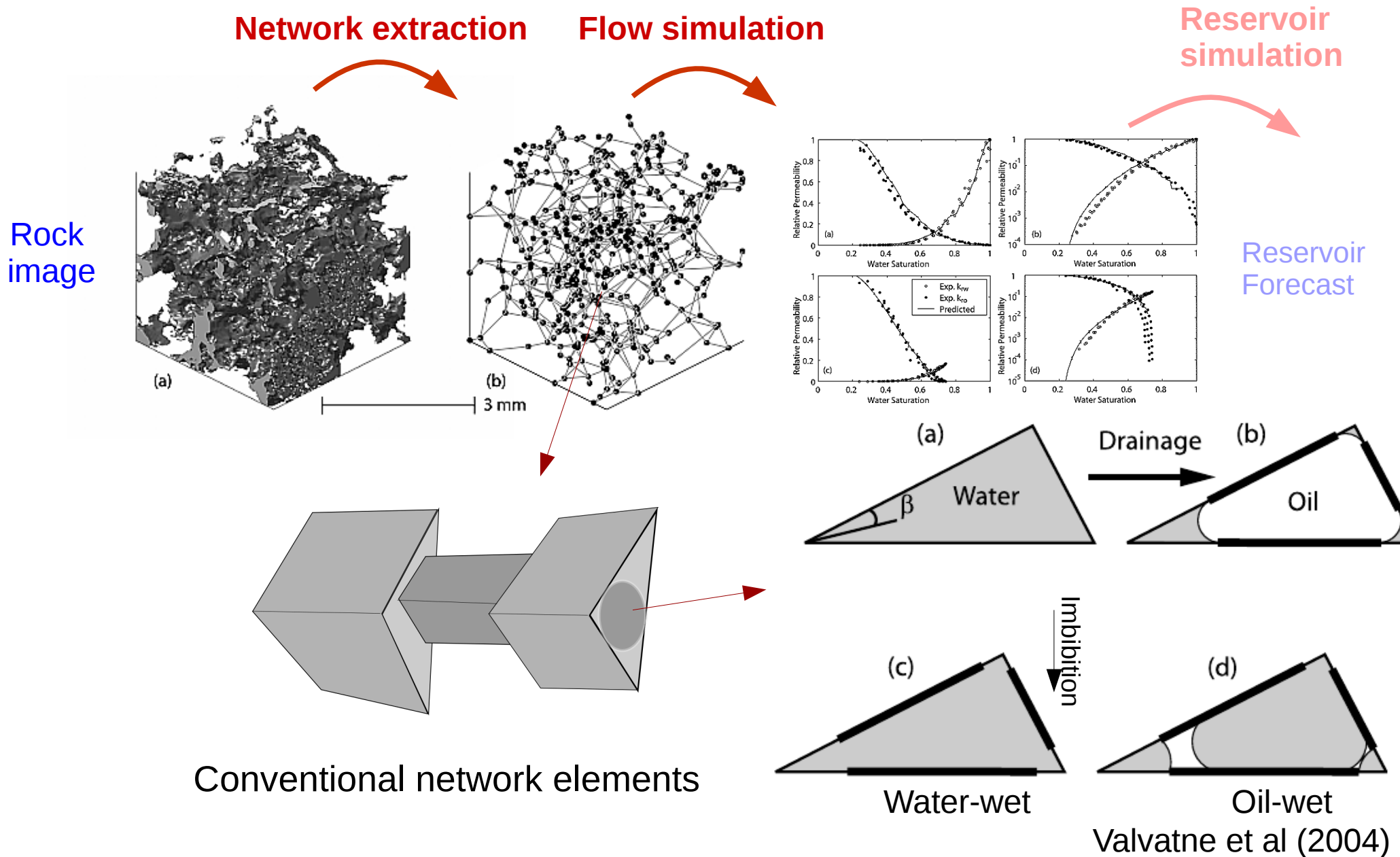
LV60A



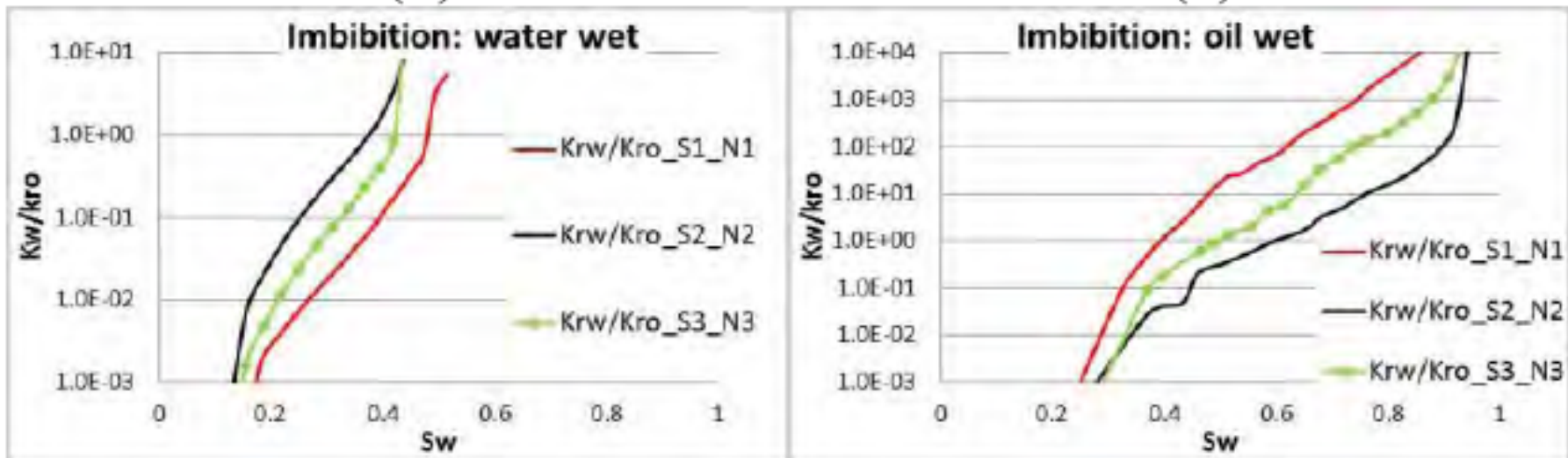
Berea

- Simulations on 1million cells took ~ 1 week on 24 processors, for a relatively high capillary number of 10^{-5}
- Simulation time increases as we decrease the capillary number
- Two-phase flow REV-size is about 1 billion grid-blocks

Background: upscaling from pore-scale to Darcy-scale:



Background – conventional network model drawbacks



Igor Bondino (Total) et al. SCA (2012):

- different network extraction algorithms give significantly different results for the same fluid/rock system

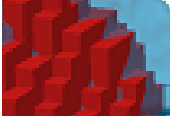
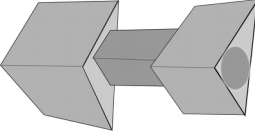
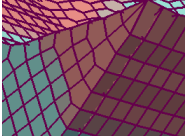
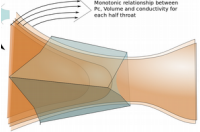
Source of the problem:

Oversimplification of the pore-space

Summary: **conductances** and **volumes** may not be assigned correctly to **capillary radii** and hence causing uncertainty in the flow modelling predictions

Background: Sources of uncertainty

Description of the pore space:

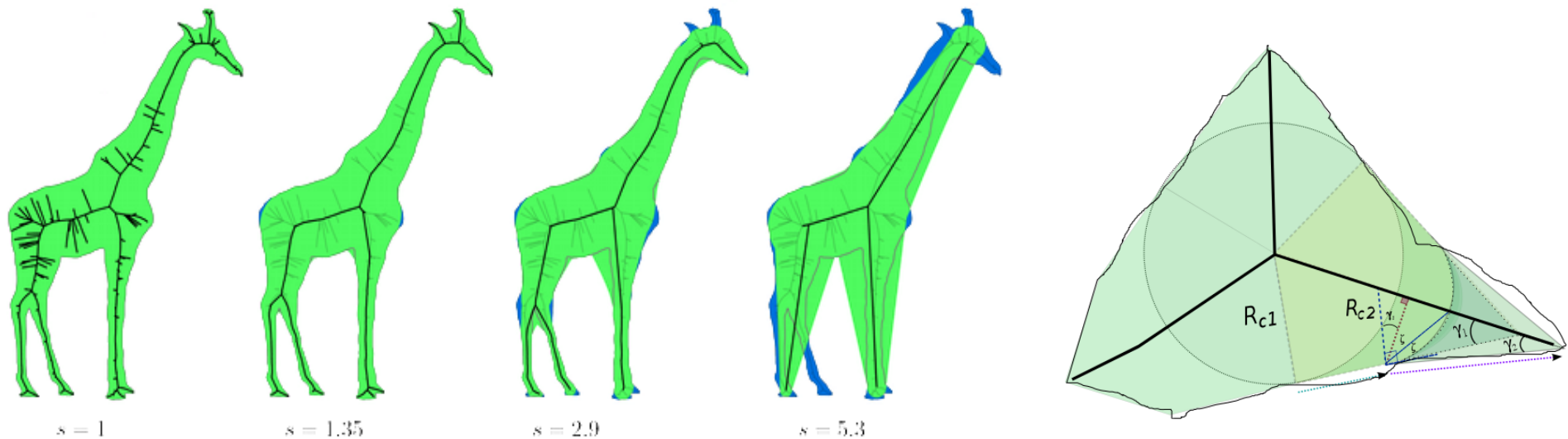
	Discretization	resolution	Error	uncertainty
→ Segmentation / noise	stair-case 	high	$O(\delta x)$ to $O(\delta x^2)$	Low to high
→ Conventional pore network	stair-case 	low	$O(\delta x)$, Refinement not possible	High
→ Direct simulations	linear 	high	$O(\delta x^2)$	Low to medium
→ Generalised networks	linear 	medium	$O(\delta x^2)$, Refinement not implemented	Low to medium

Fluid/rock properties:

- Contact angle distribution
- Clay / micro-porosity identification

Mathematics behind network extraction

There is a one-to-one relationship between any 3D geometry and its medial surface → no information loss → no additional uncertainty due to oversimplification of the pore space

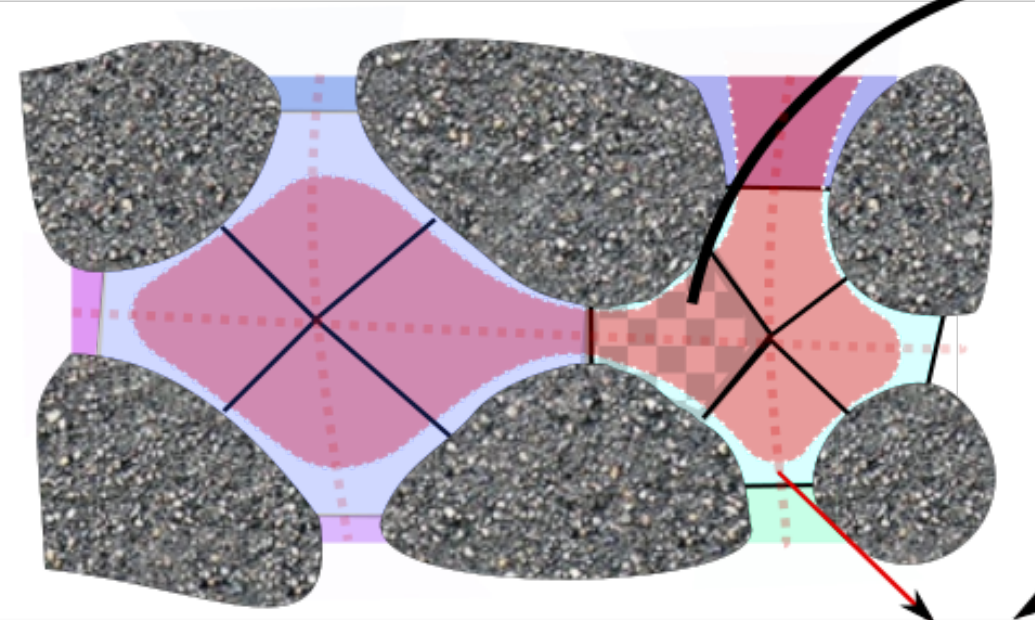


Picture from: http://www.agg.ethz.ch/research/medial_axis

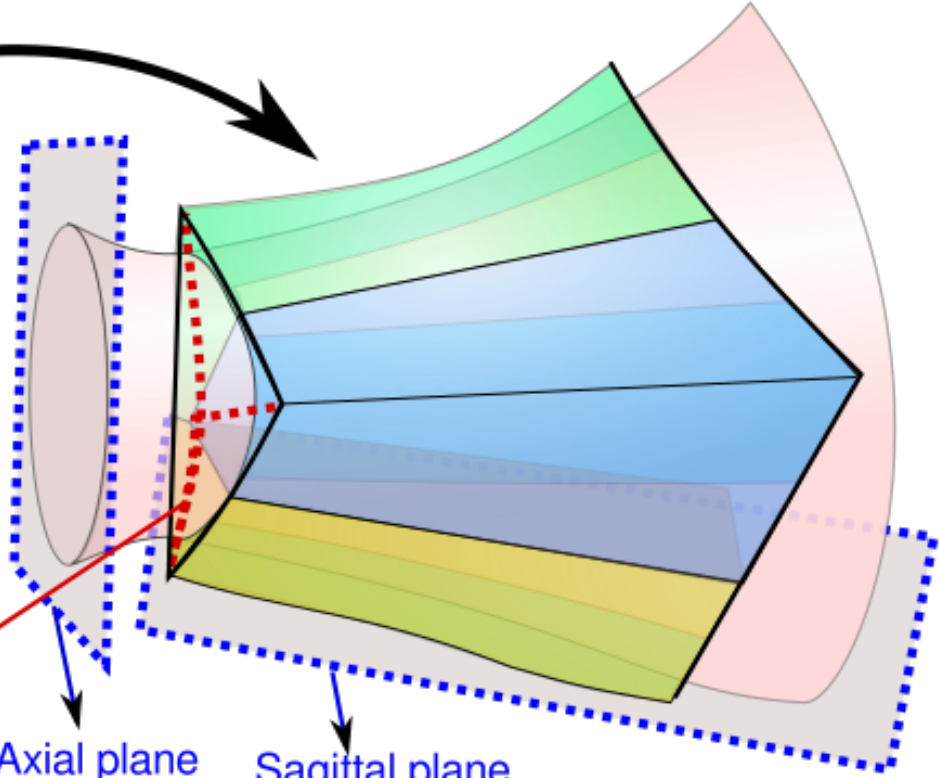
- **Generalized pore network can be viewed as a coarse medial-surface representation of the pore space**

Background – generalised network model

Half-throat



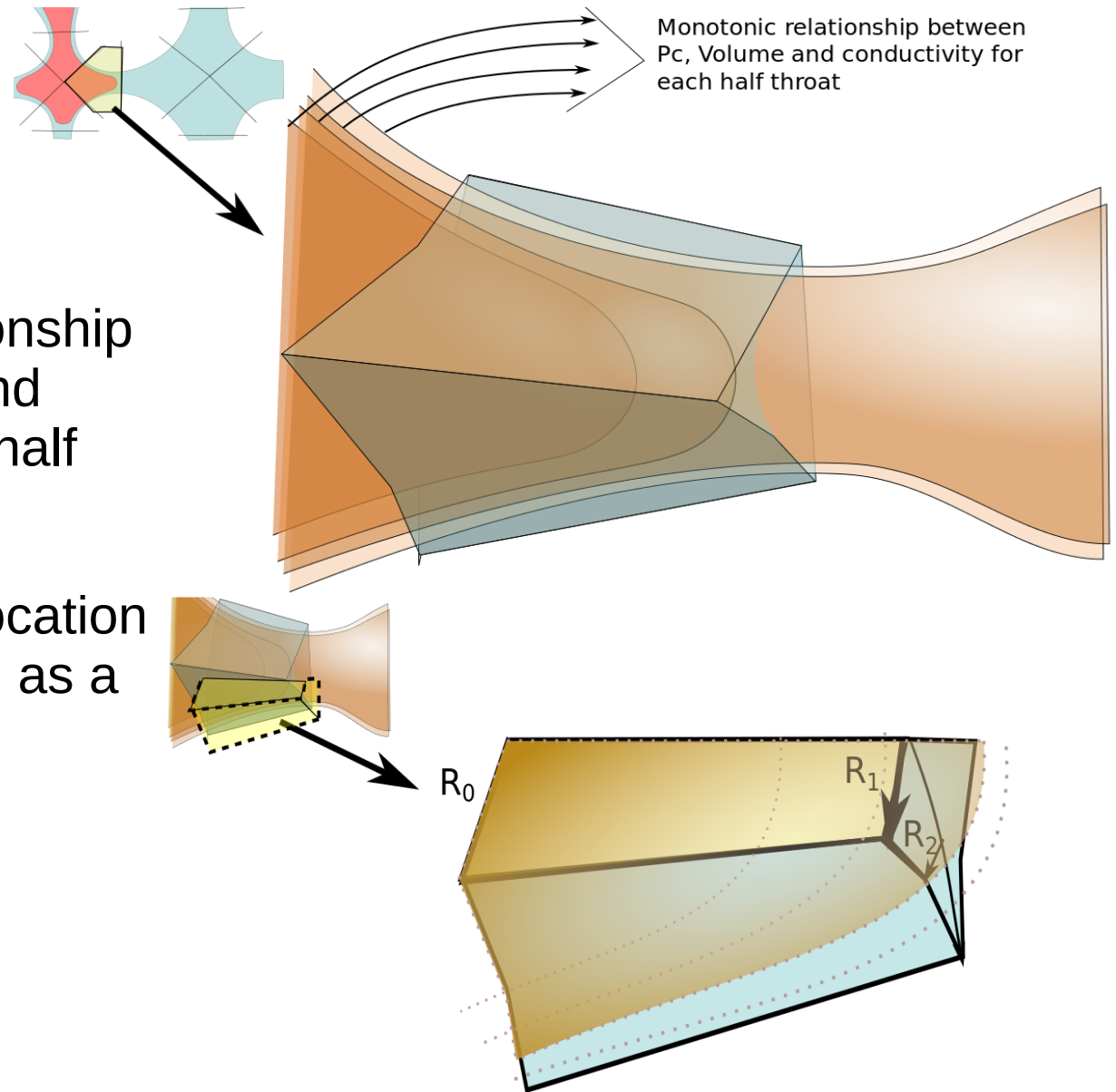
(a)



(b)

Half-throat corners

Flow simulation: interface tracking + post-processing



There is a one-to-one relationship between capillary radius and interface location for each half throat

→ allowing the interface location to be recorded and tracked as a single scalar variable

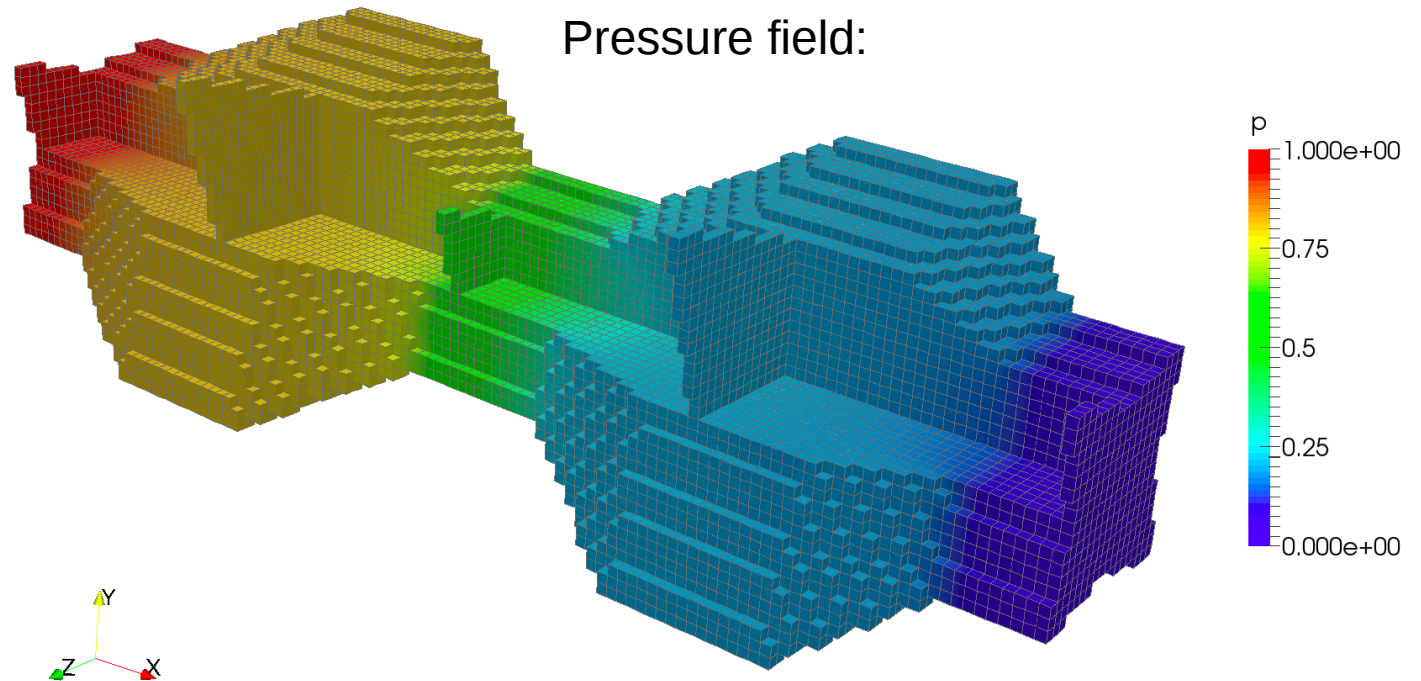
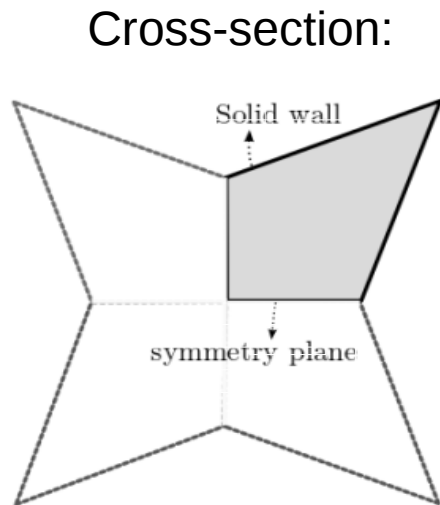
Generalized network modelling - summary

- **The necessary parameters are extracted directly from the micro-CT image as tabulated functions and used in a generalized network flow simulator:**
- - ✓ MS-P theory is used to **relate Pc to Volume (Sw)** for each pore/throat
 - ✓ The necessary parameters: pore/throat radius and corner angles.
→ extracted directly from the micro-CT image
 - ✓ Direct simulation is used to compute throat **conductivities**
 - ✓ Effect of contact angle and other fluid properties is added during flow simulation.

Generalised pore-network **extraction**

Parametrisation of pore space

Pore-space converted into voxelated (micro-CT) format, on which Navier-Stokes equations are solved: and converted into generalised network format

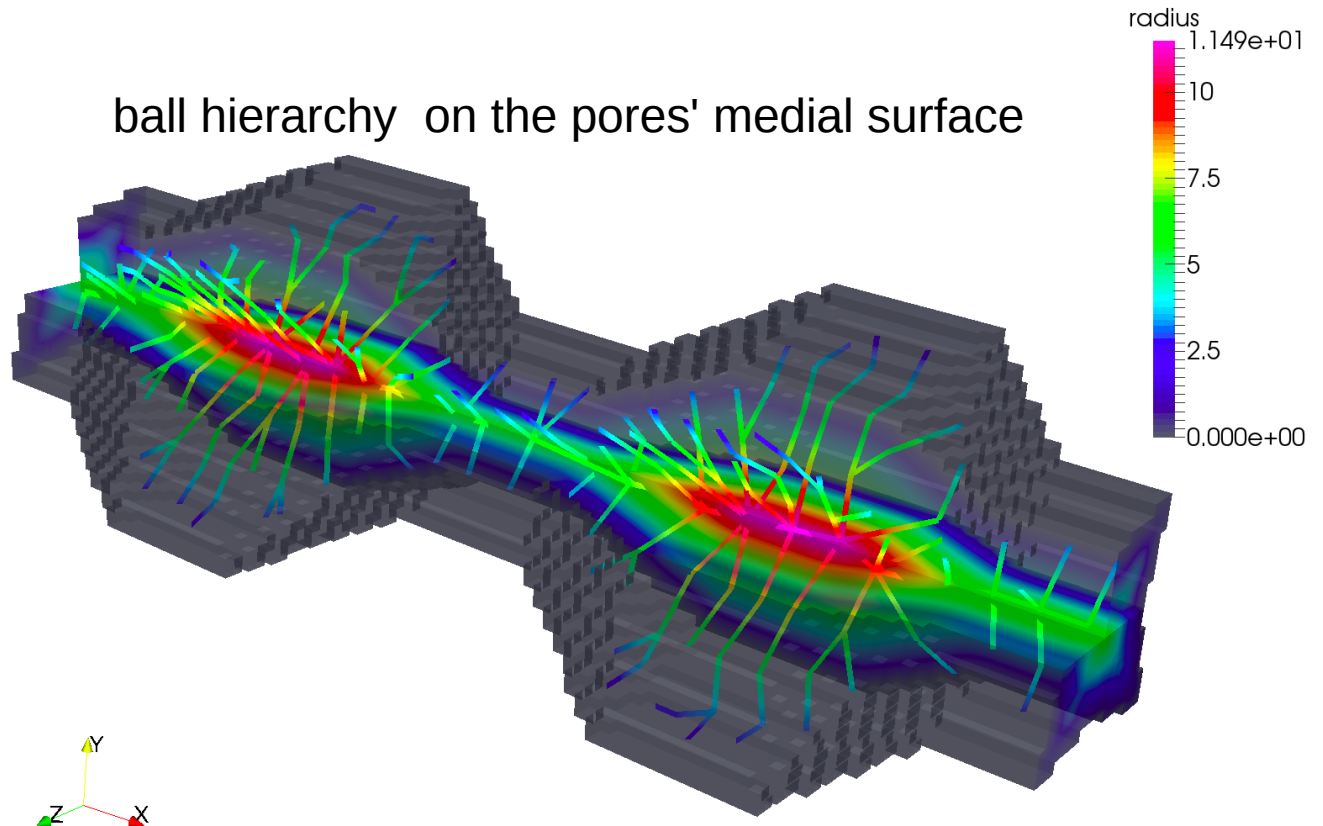


Parametrisation of pore space

Step2.

- 2.1. Extract ball hierarchy (medial surface),
- 2.2. find local maxima of the radius and
- 2.3. Assign a pore index to each of the local maxima

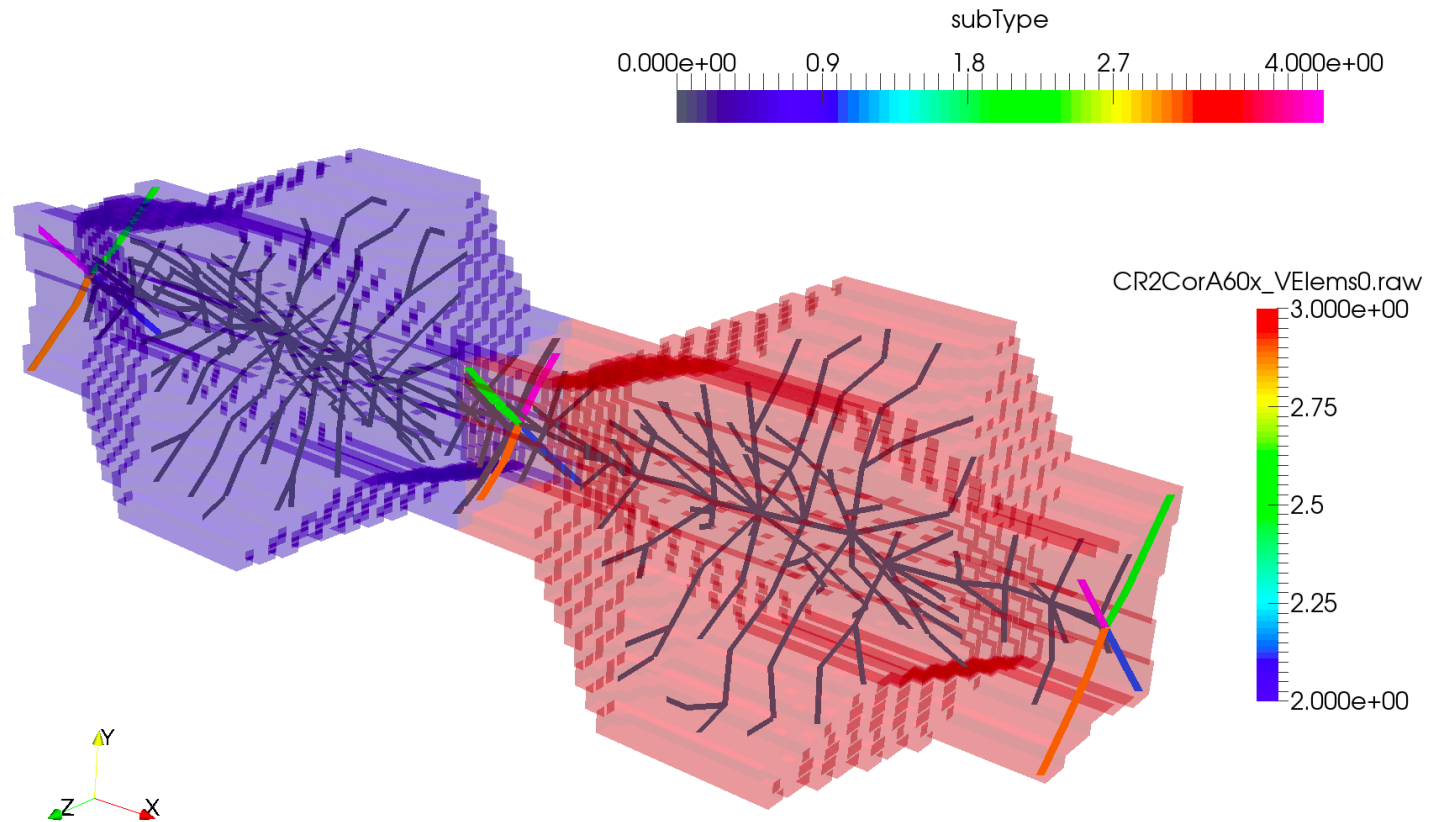
ball hierarchy on the pores' medial surface



Parametrisation of pore space

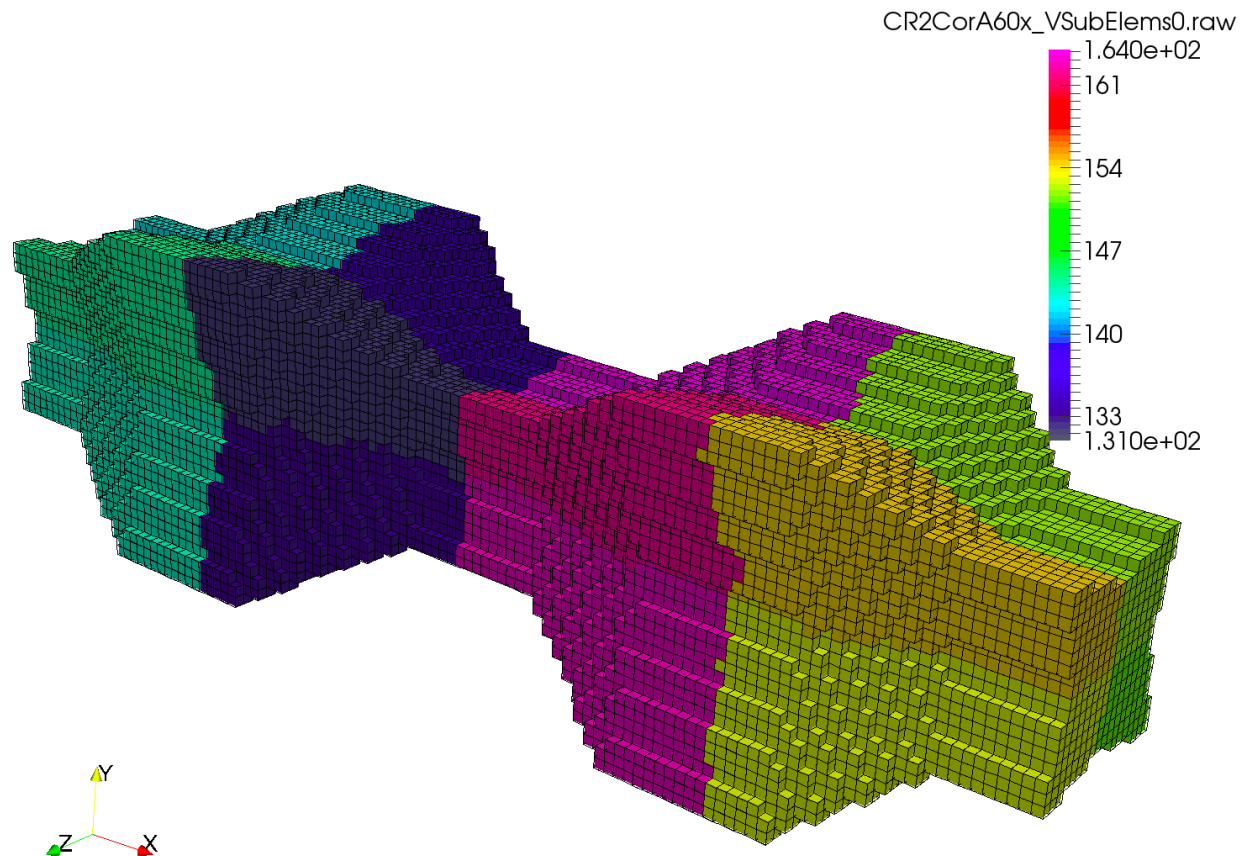
Step 3.0. Find pore-to-pore connections (throats)

Step 3.1. Identify and label individual corners on the throat medial axis



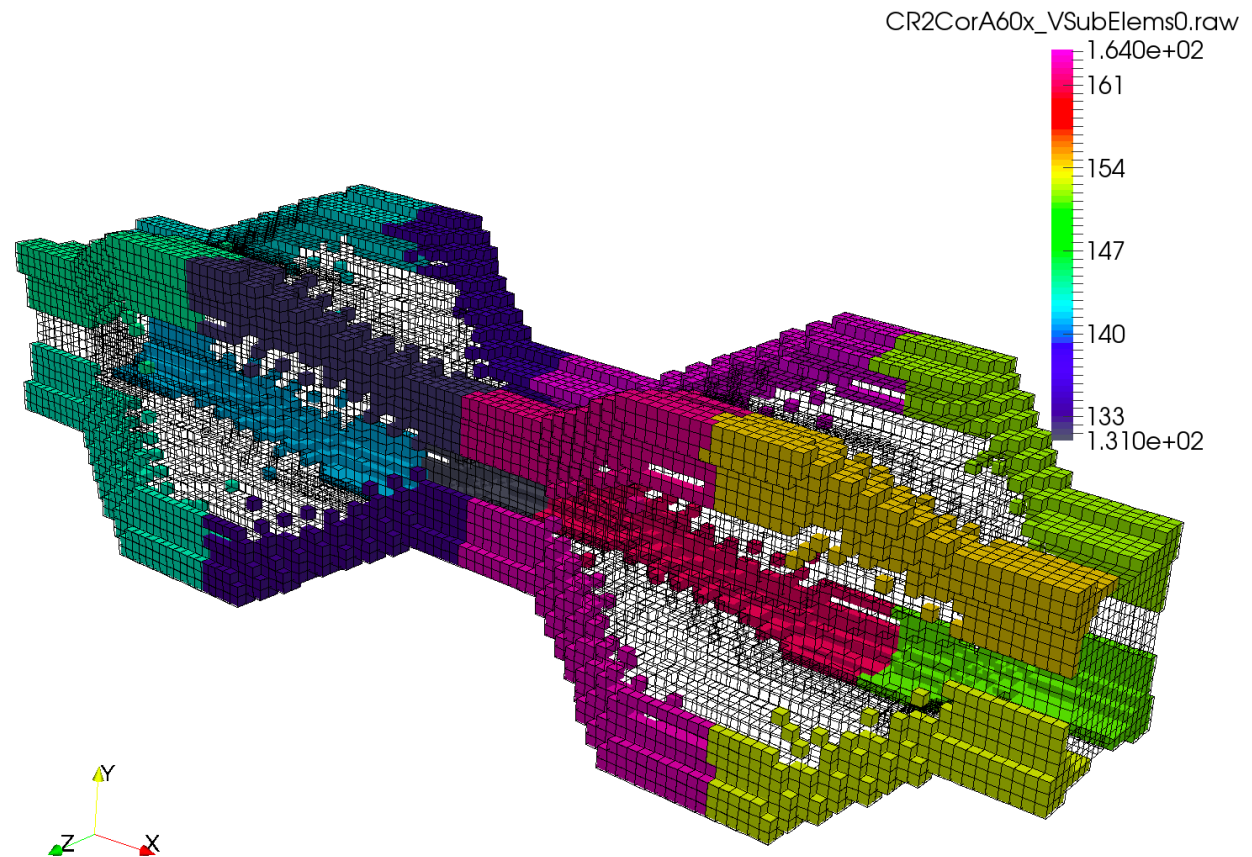
Parametrisation of pore space

Step 3.1. Map corner labels to the original image and analyse direct simulation results for each corner at Level 0



Parametrisation of pore space

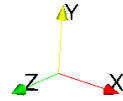
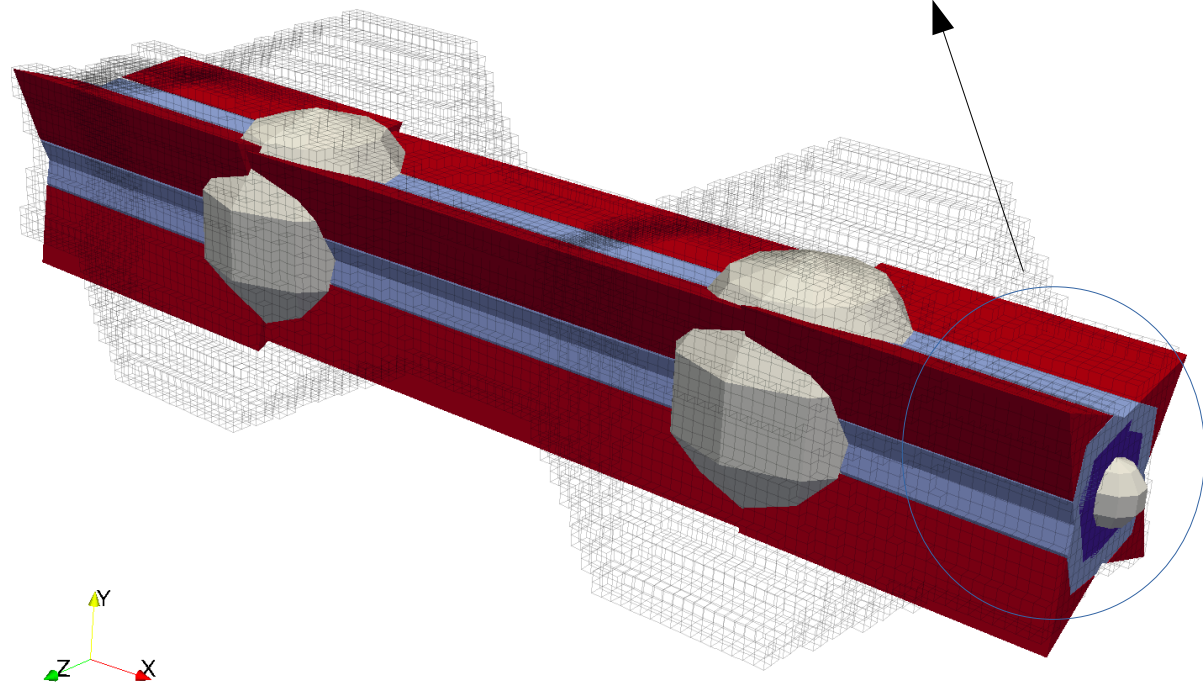
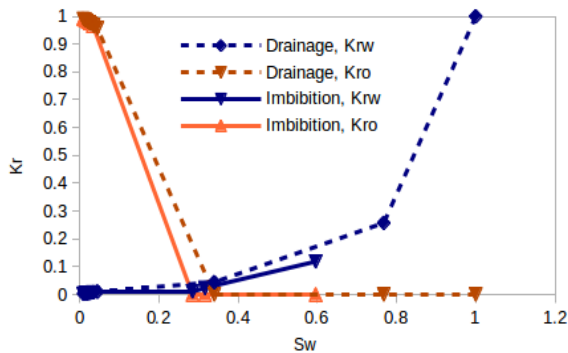
Step 3.3. More balls from throat centres are deleted from corners to analyse direct simulations at Level 2



Comparison - generalized network extraction

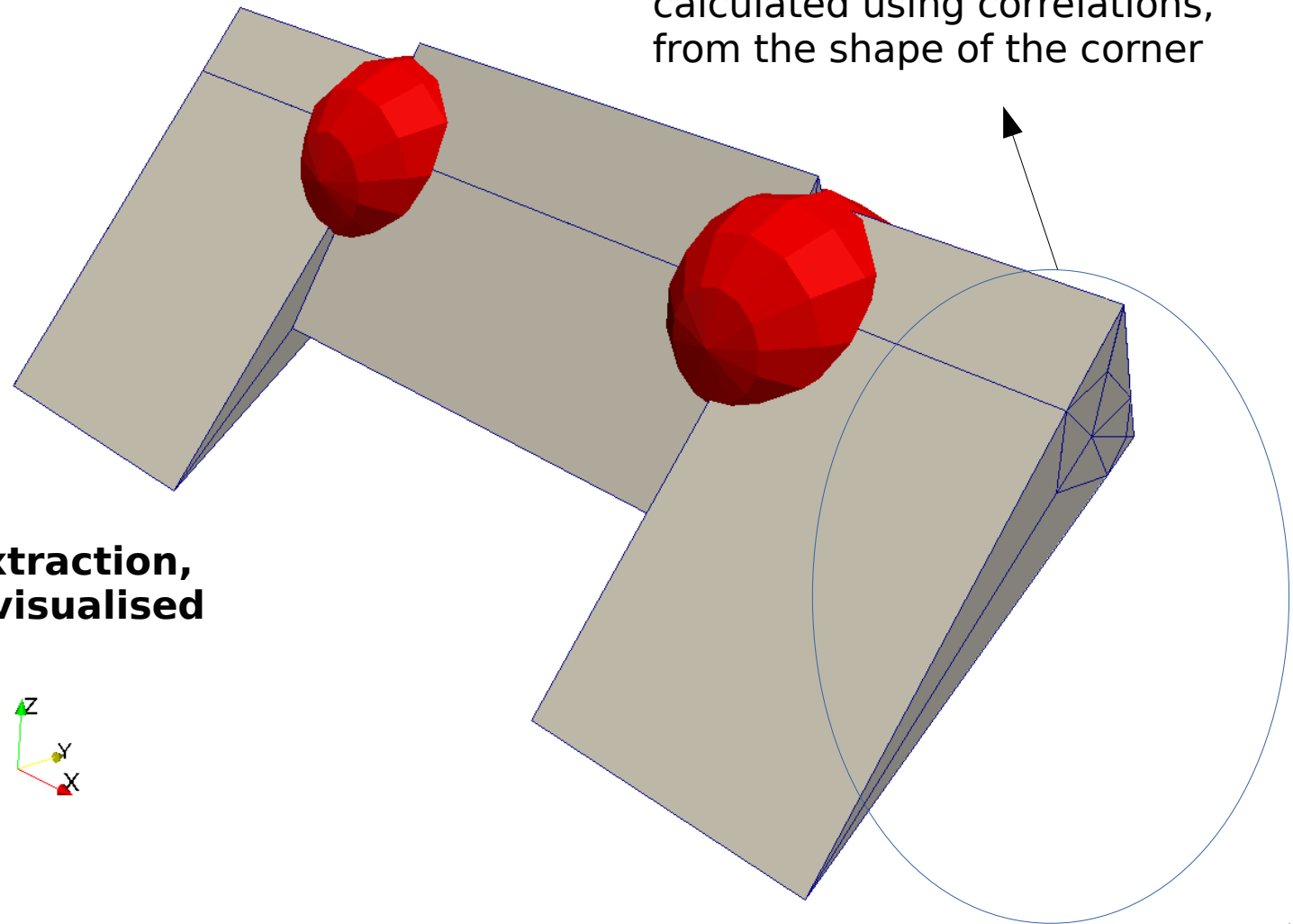
Corner conductivities are calculated from tabulated functions extracted from direct single-phase flow simulations

Generalised network simulation results - validation in progress



Comparison - conventional network extraction

Corner conductivities are calculated using correlations, from the shape of the corner



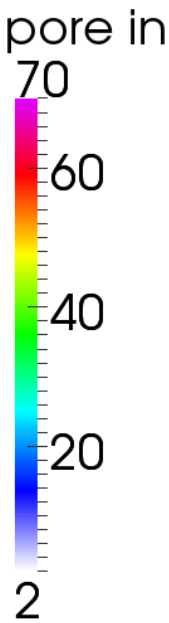
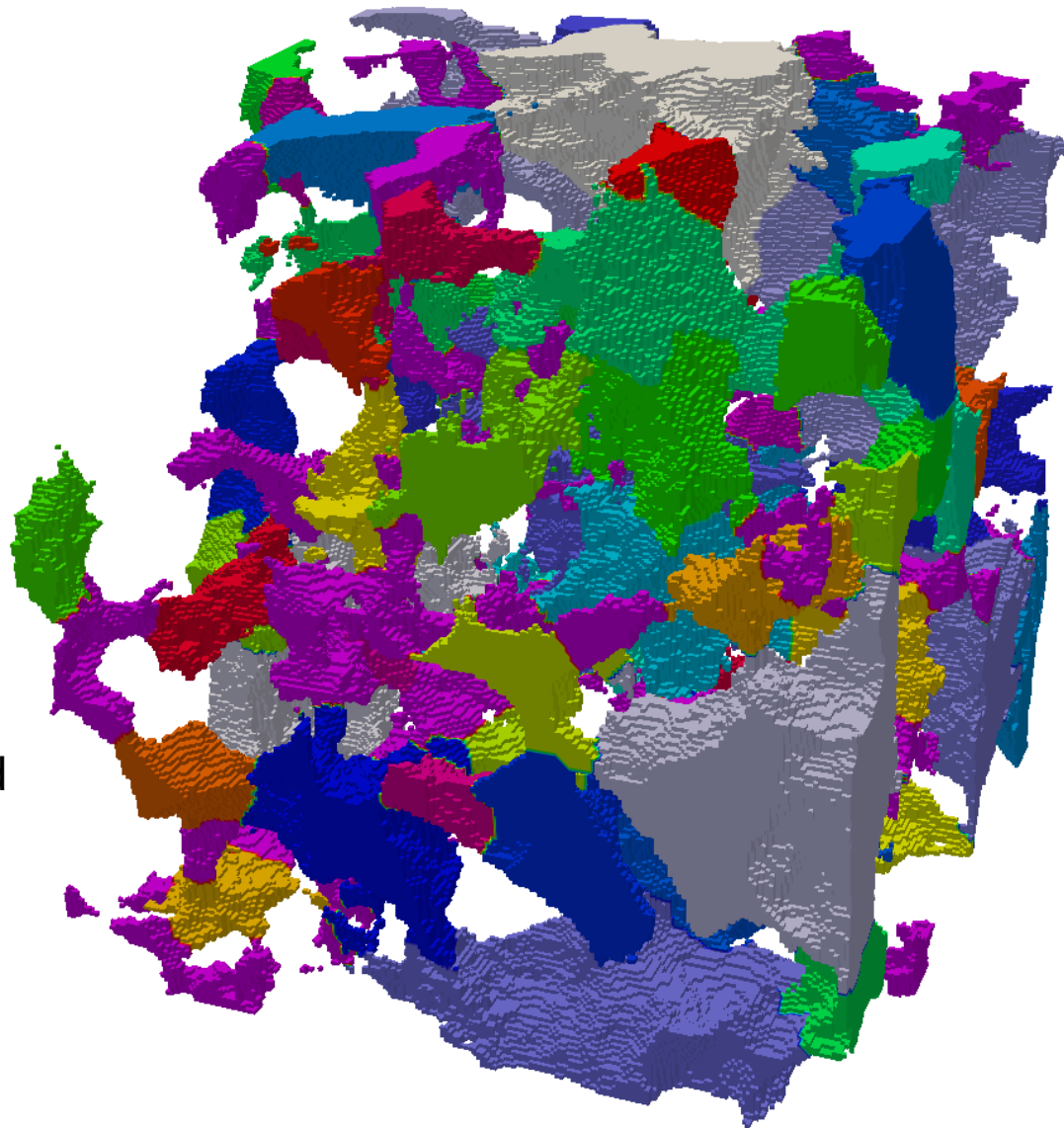
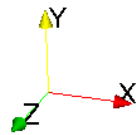
**Conventional network extraction,
Only throat corners are visualised**

Network extraction demonstration – Berea sandstone

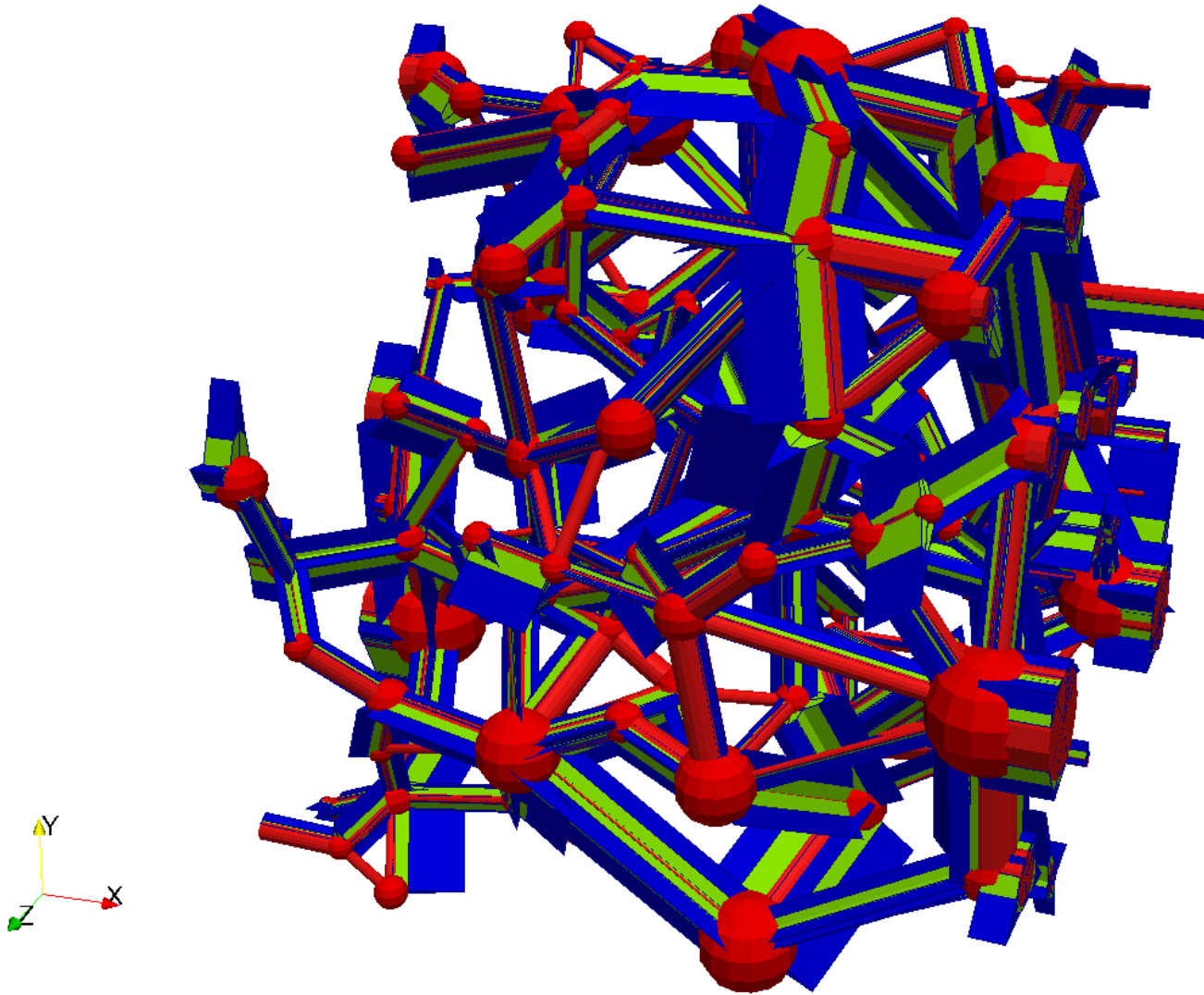
Element Indices

✓ High resolution

➔ Important for detecting and analysing corners accurately

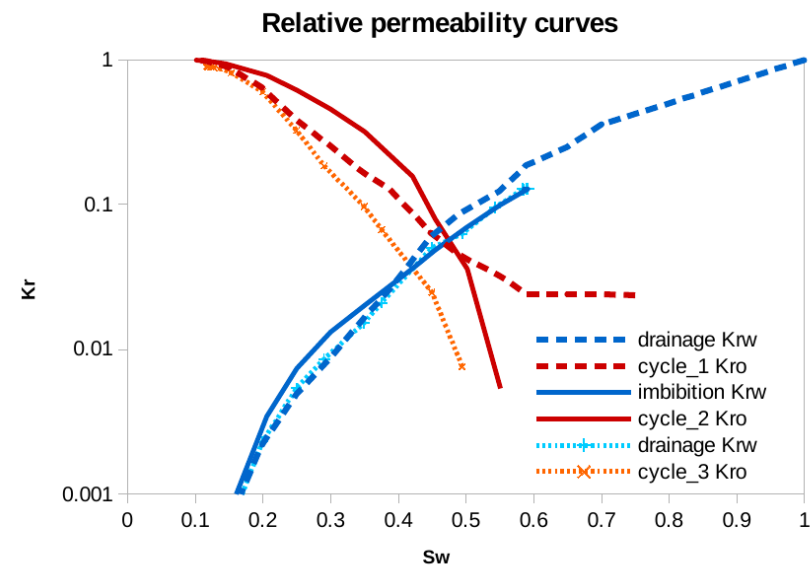
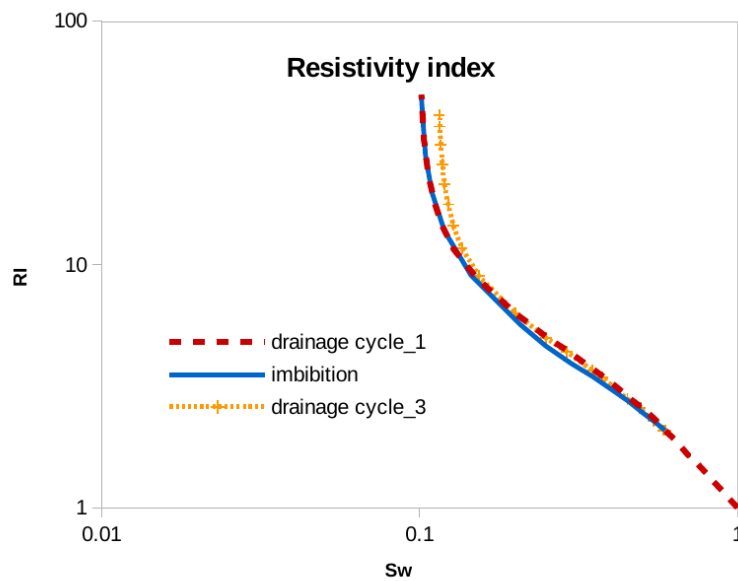
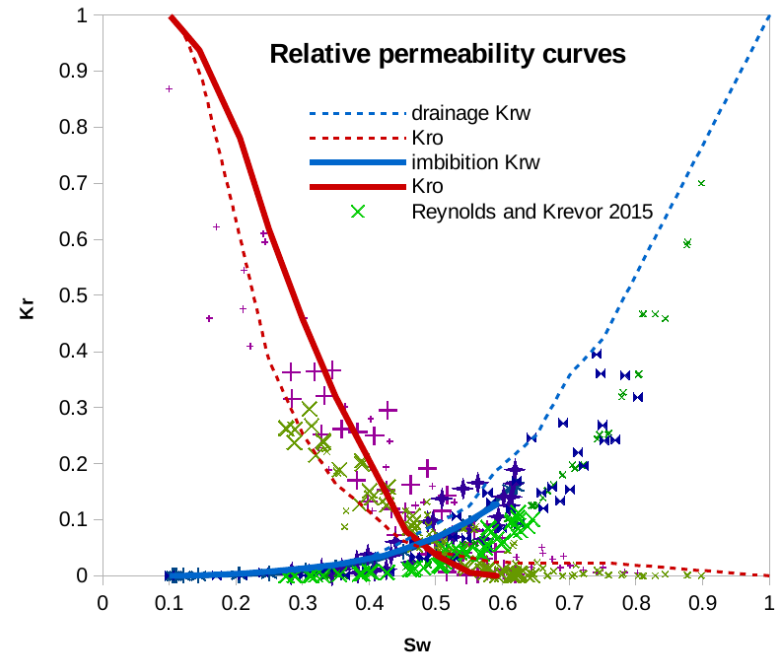
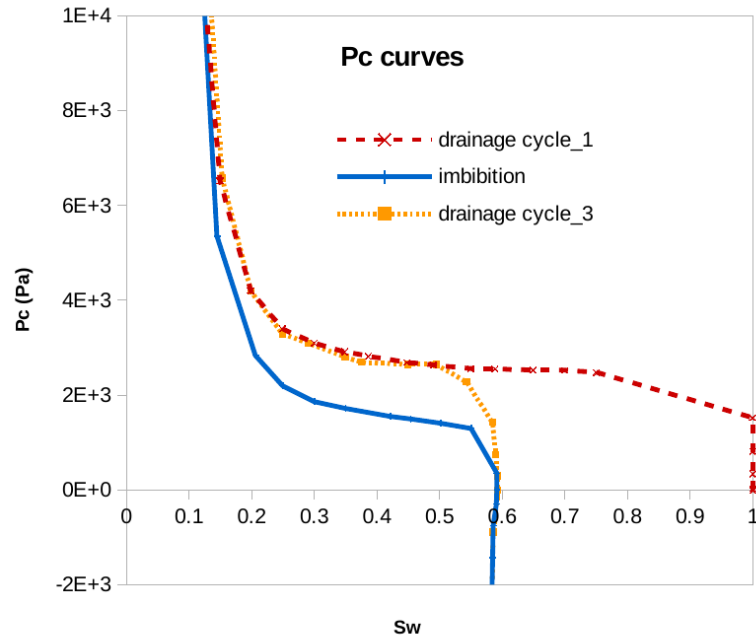


Generalised network in flow simulator, a primitive visualization



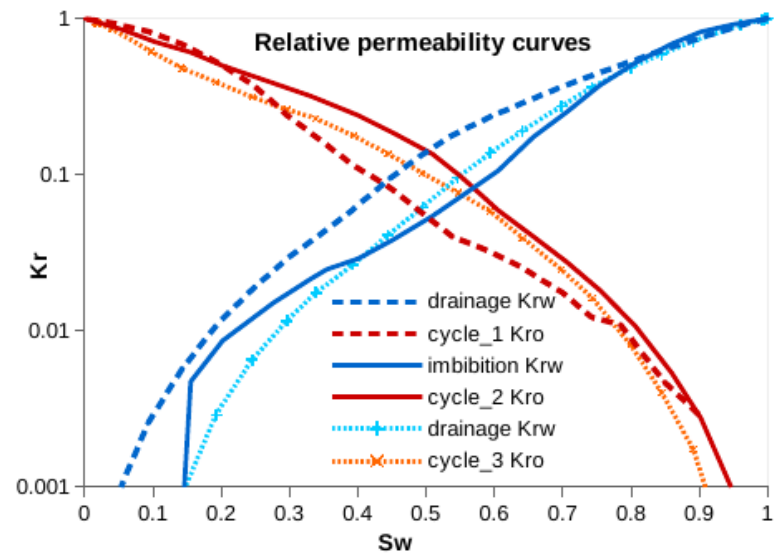
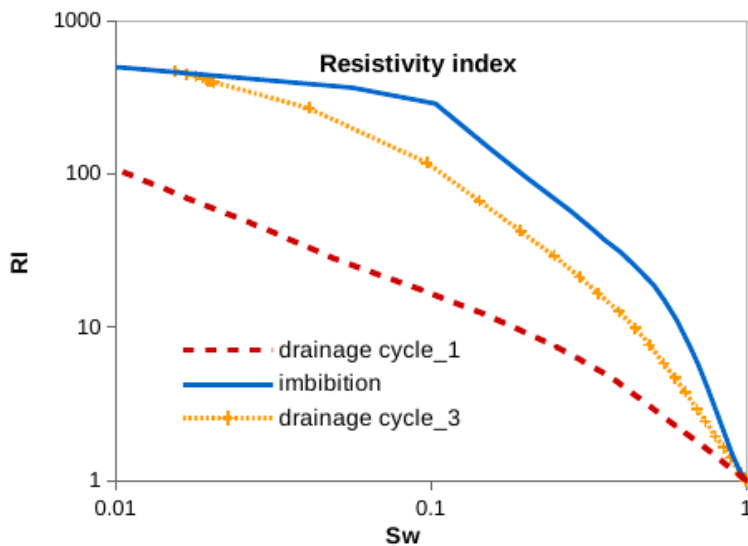
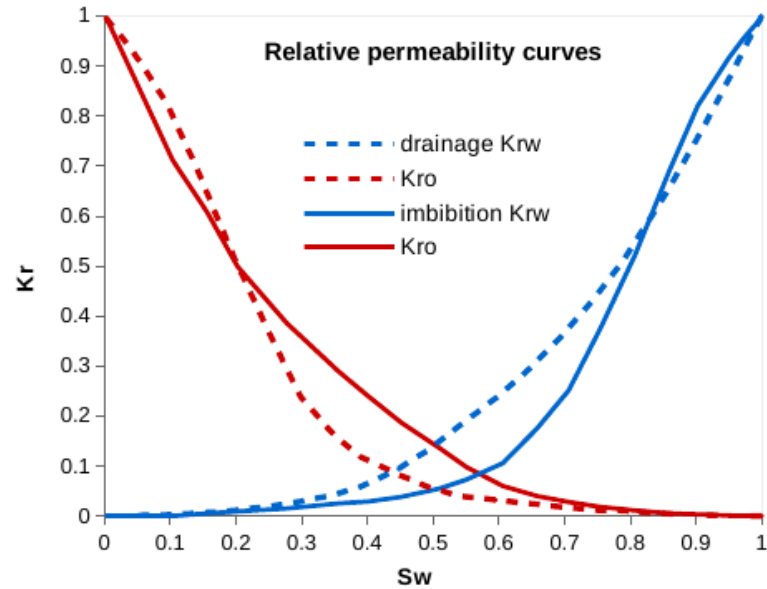
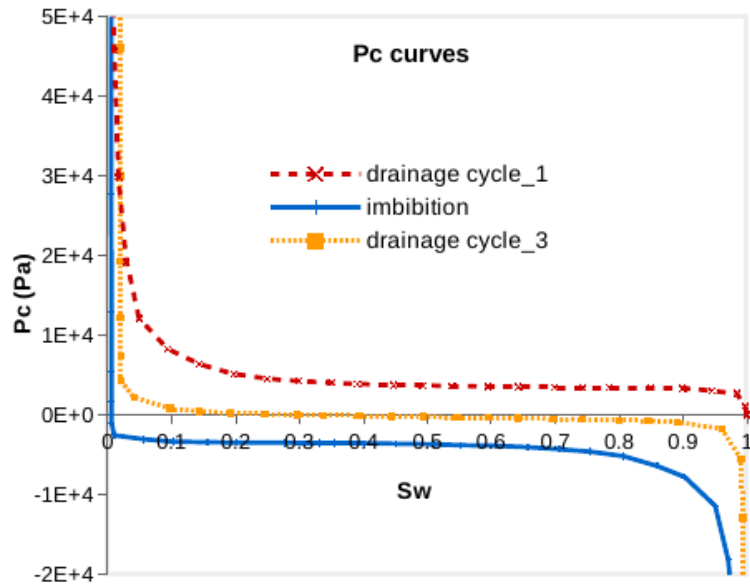
Relative permeability predictions – water-wet

Bentheimer sandstone (from IC-PSM website, 1000^3 @ 3um)



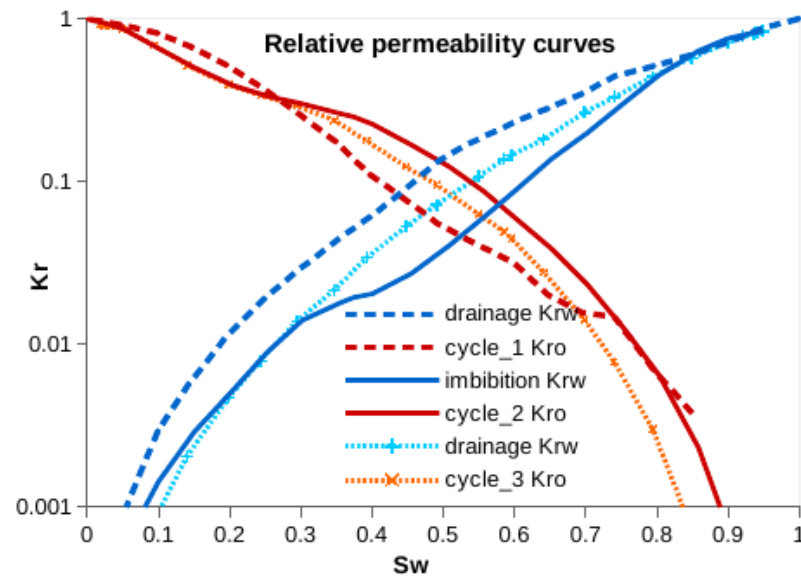
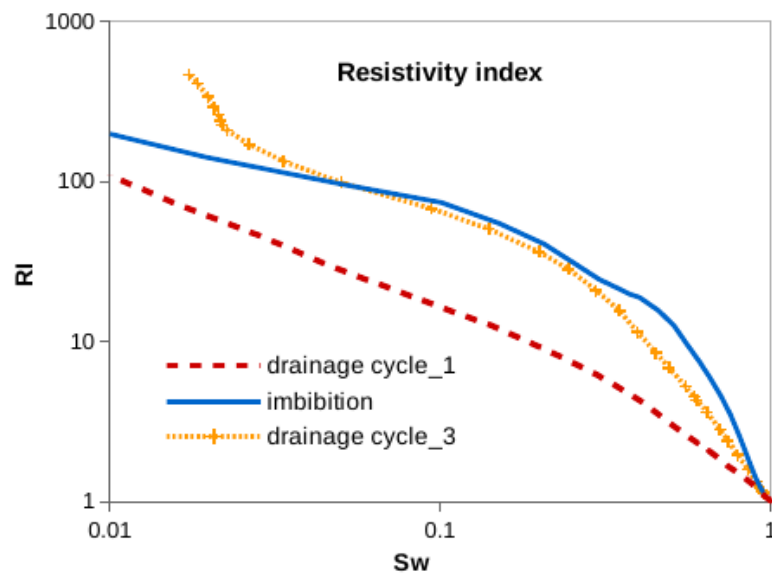
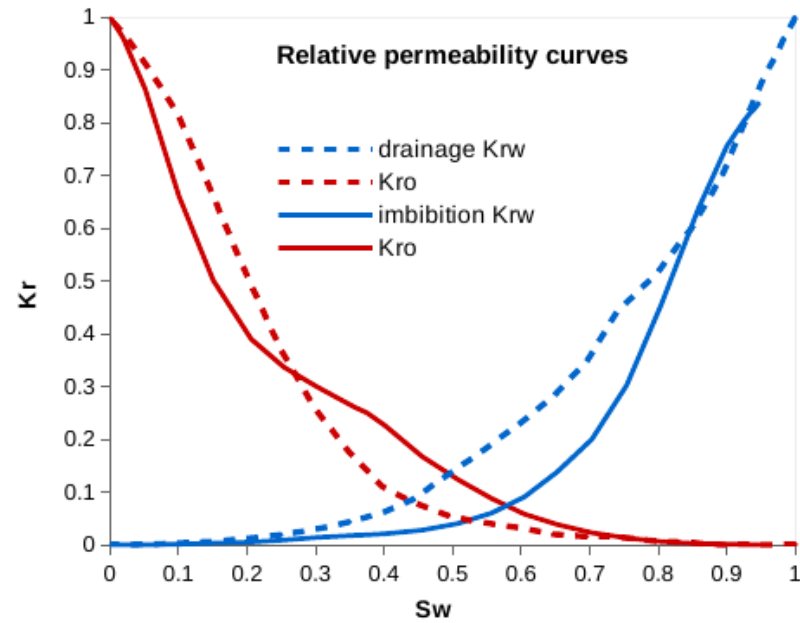
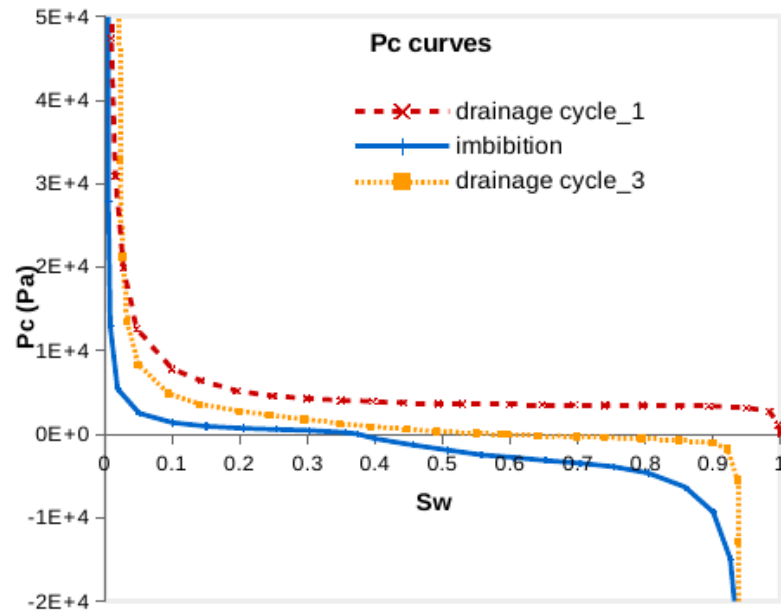
Relative permeability predictions – oil-wet

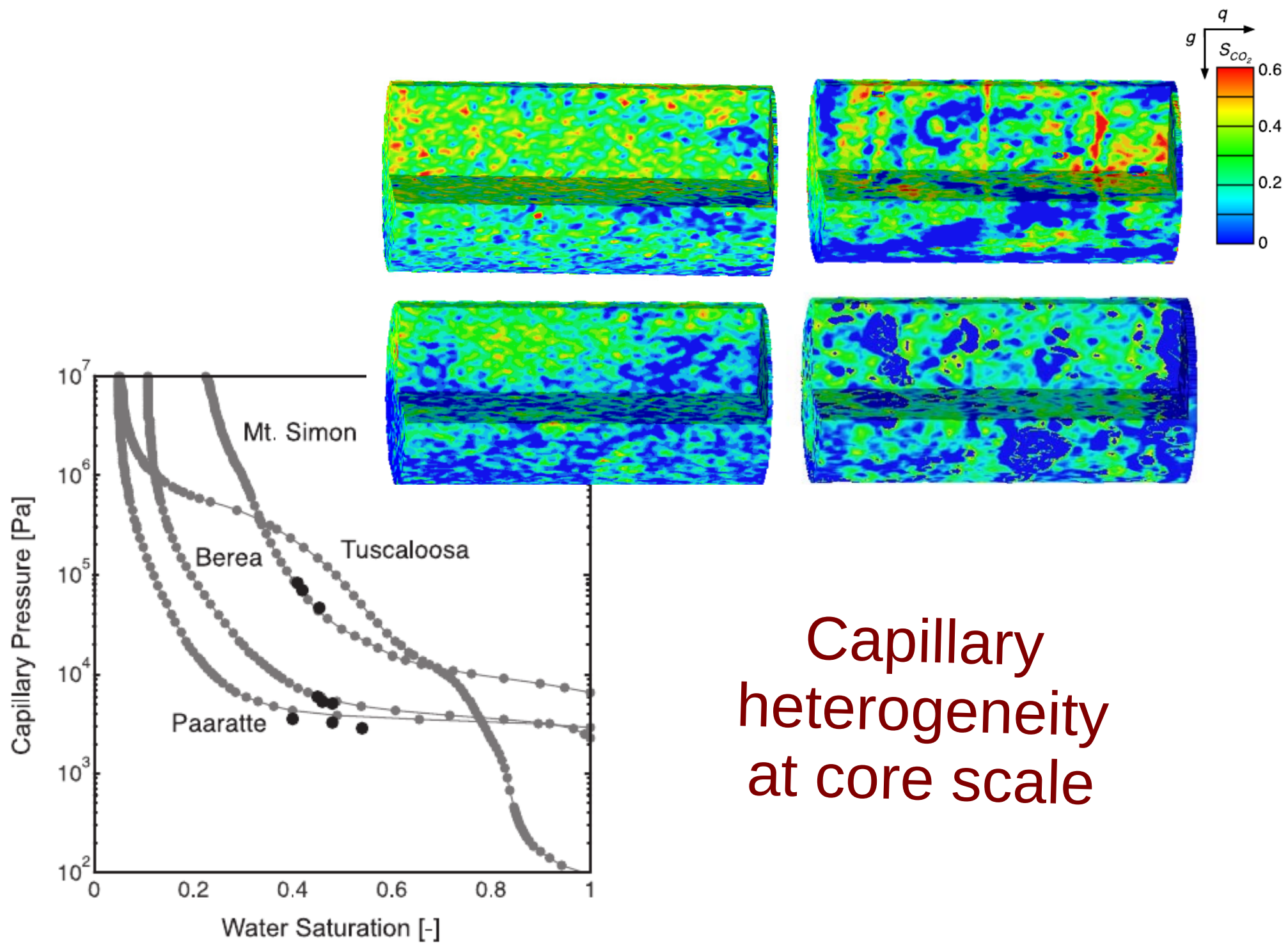
Bentheimer sandstone (from IC-PSM website, 1000^3 @ 3um)



Relative permeability predictions – **mix-wet**

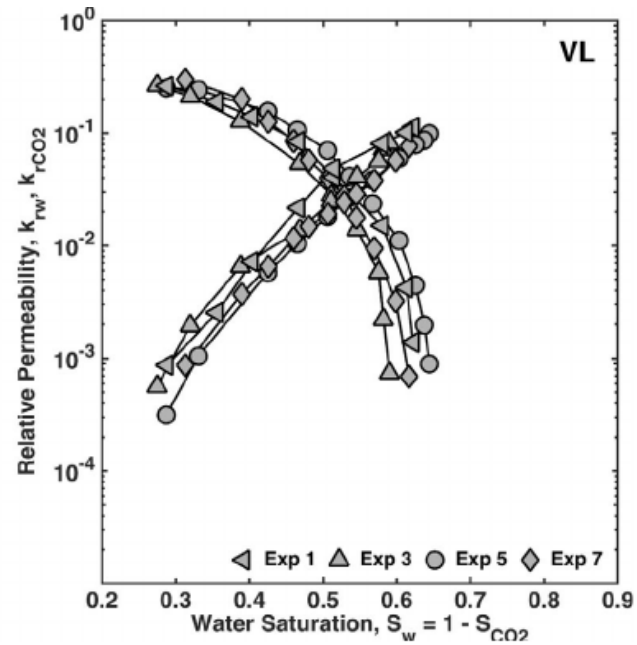
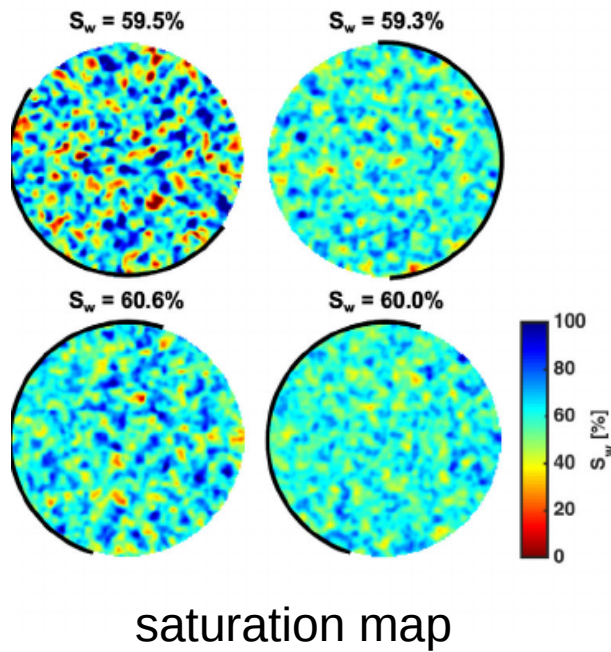
Bentheimer sandstone (from IC-PSM website, 1000^3 @ 3um)





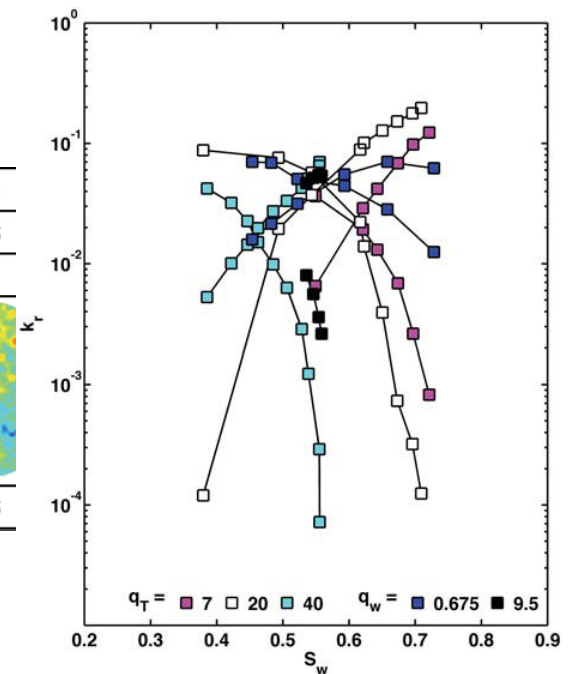
Capillary
heterogeneity
at core scale

Experimental relative permeability measurement

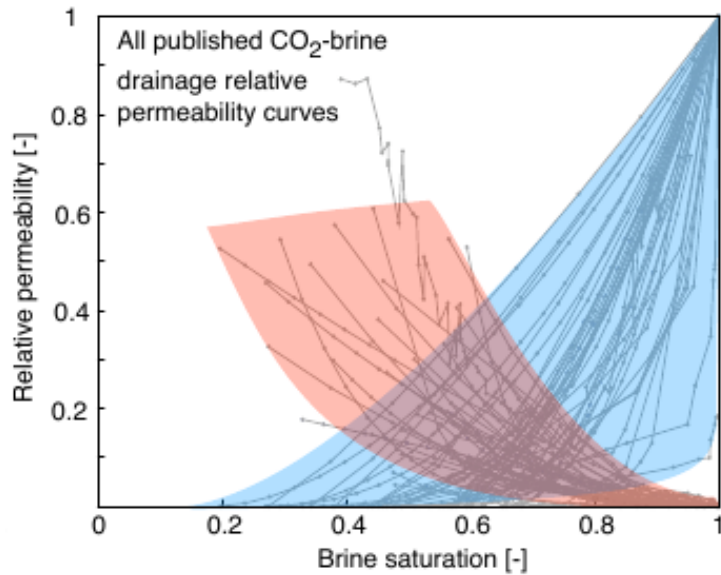


f_{N_2}	0.943	0.511	0.54762	0.62	0.72857	0.1	0.81
q_{N_2} (ml/min)	6.6	19.489	11.5	15.5	25.5	4.0	40.5
q_T (ml/min)	7	20	21	25	35	40	50
S_w (%)	54.7	54.3	55.8	55.6	53.3	55.1	52.5

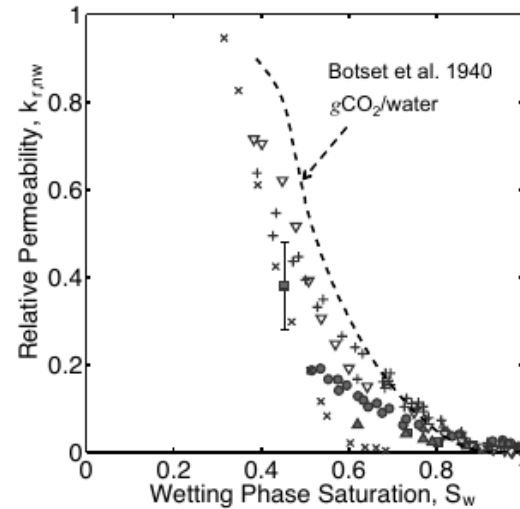
Reynolds and Krevor (2015)



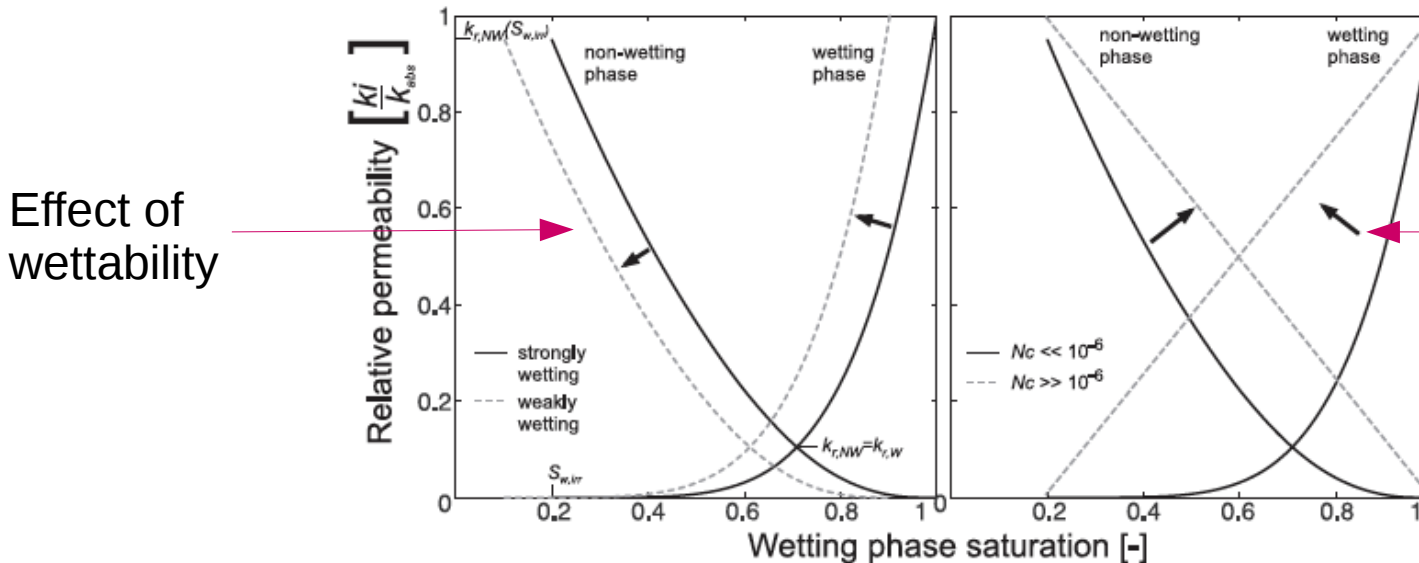
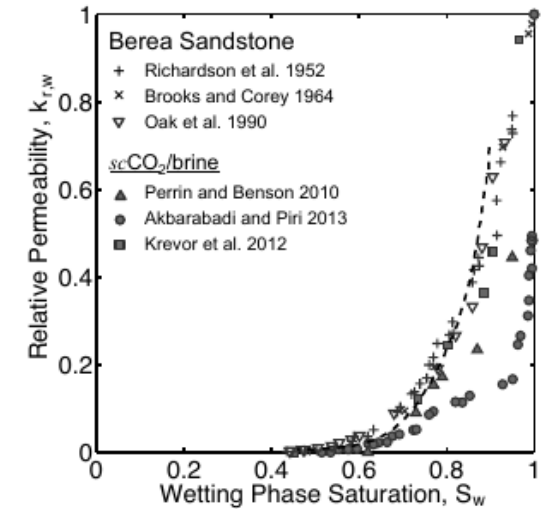
Experimental relative permeability measurement



Different rocks, CO₂-brine



Berea sandstone, CO₂-brine



Effect of wettability

Effect of capillary number

Validation with direct simulations - **star-shaped** pore-throat system

Star-shaped geometry parameters:

Corner Angle : 45 degrees

$$R_{\text{pore}}/R_{\text{throat}} = 2.25$$

$$R_{\text{pore}}/L_{\text{throat}} = 1.25$$

Direct simulation

simulations run on 12 processors

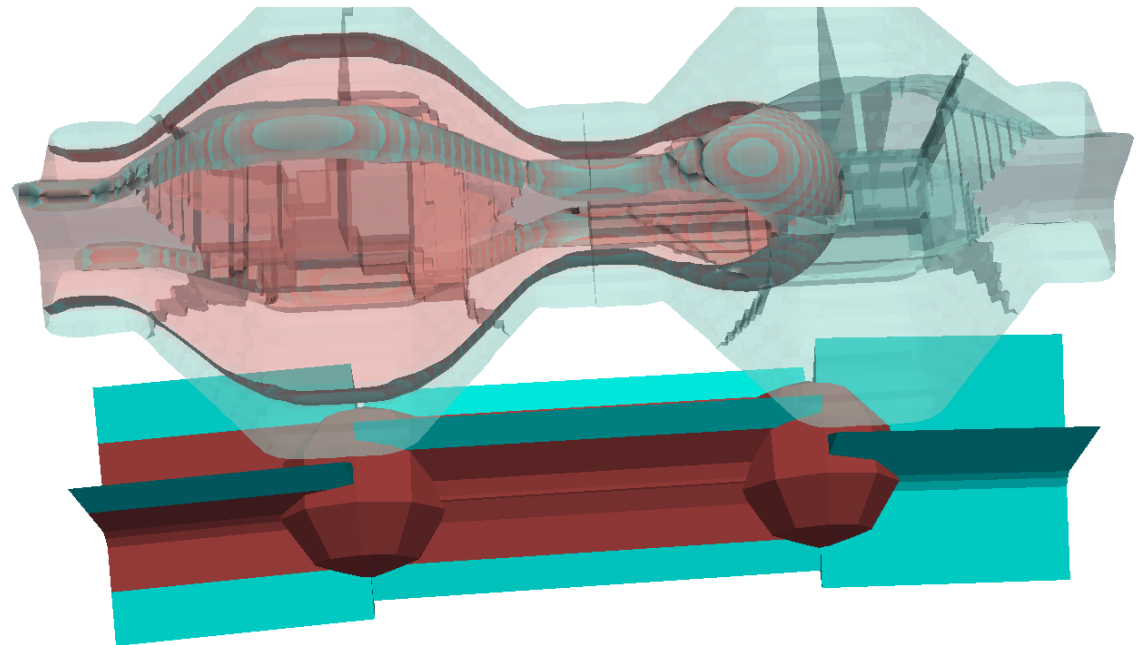
Simulation time:

3-4+ hours for $R_{\text{throat}}/\delta x = 5$

2-3+ days for $R_{\text{throat}}/\delta x = 10$

Network model

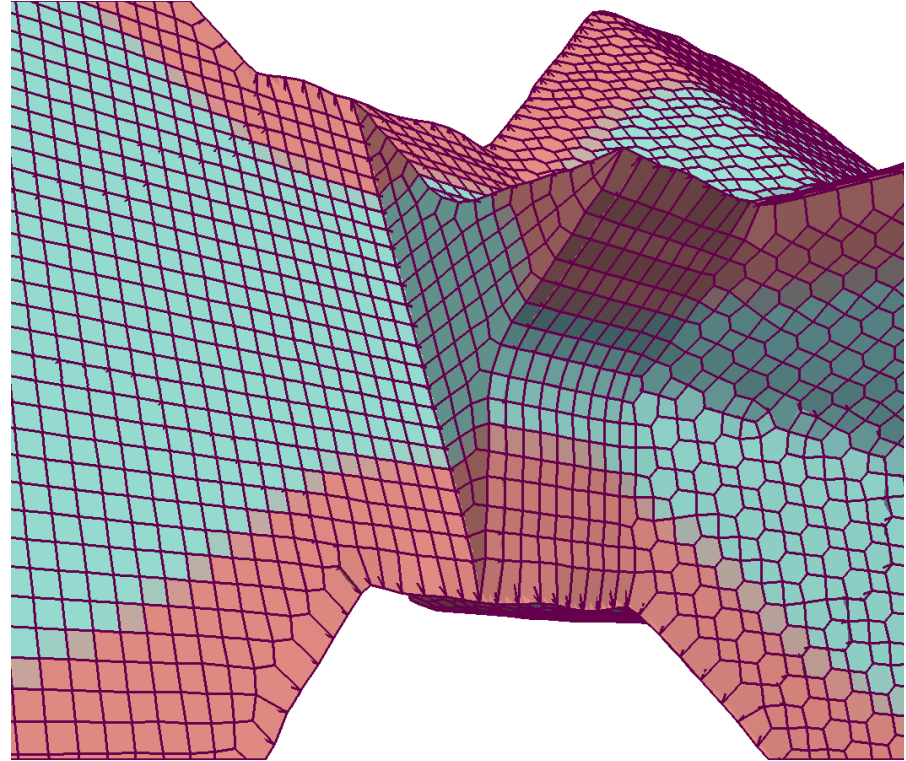
Simulation time: **less than 1 second**



Note: network model visualization is primitive, the actual calculations are done based on a fixed 4-segment linear interpolation scheme for each half corner

Improvements to direct simulations:

- Mesh quality improvements:



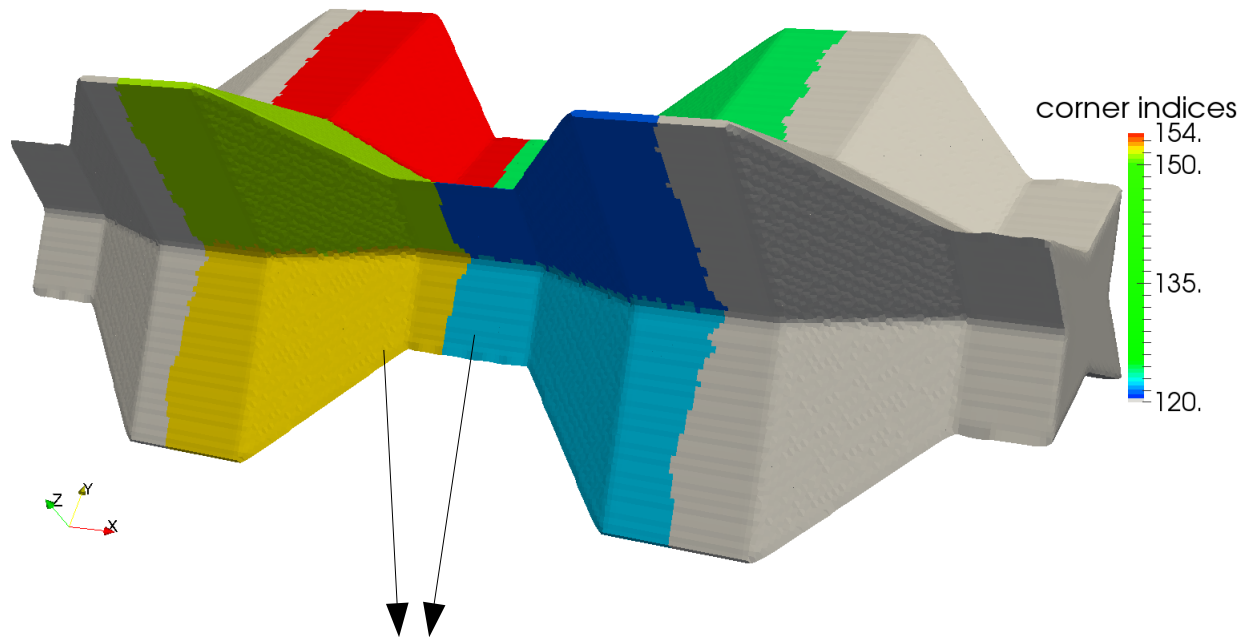
- Improved surface tension model using explicit interface reconstruction

Thanks to **Mosayeb Shams** for sharing his code, **Arthur Moncorgé** and **Stephane Zaleski** for their feedback and advice

Post-processing direct simulations

Corner Indices from network extraction are used directly as input control-volumes during direct simulations,

The code that is used for computing relative permeability curves is used to analyse individual corners extracted from rock images.



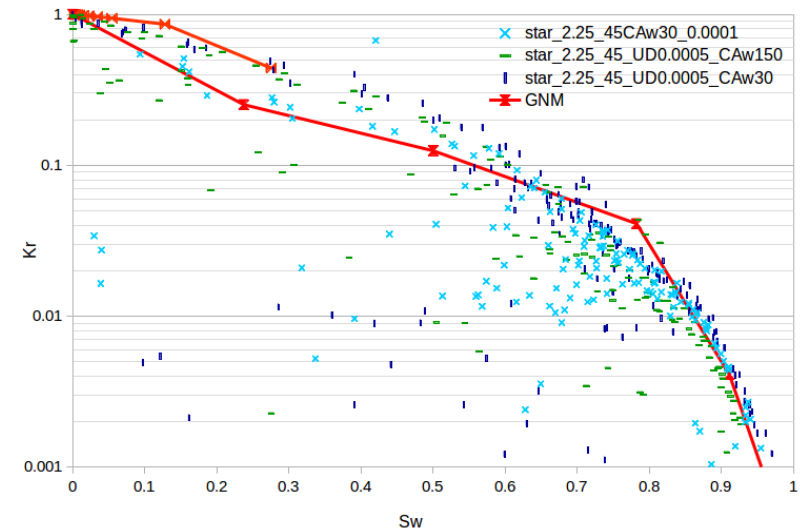
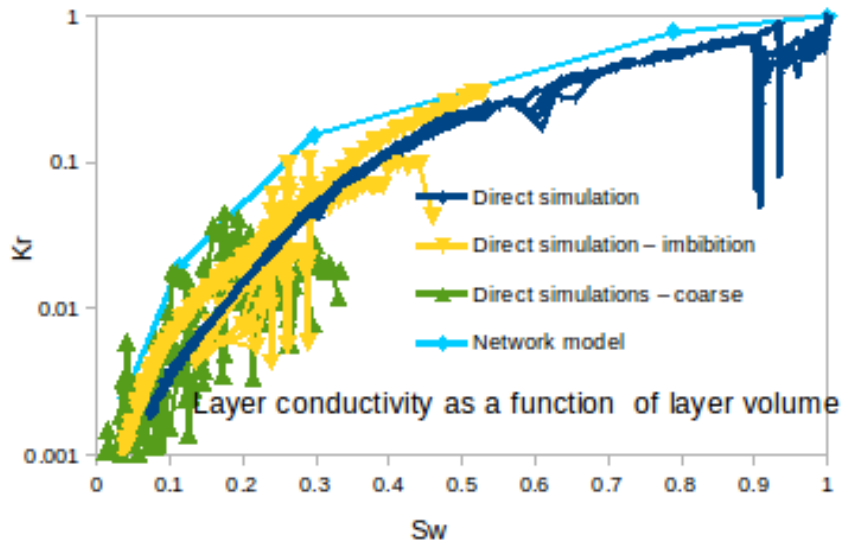
Each throat (pore-to-pore connection) is divided to 8 half throat corners during network extraction stage

Validation of network flow simulator code saturation/conductivity computations

Layer saturation/conductivities depend on:

- **Corner cross-sectional shape**
- **Variation of the cross-sectional area along the pore-to-pore connection**
- **Flow pattern / rates**

Direct simulations, star-shaped, drainage, piston-like invasion followed by layer flow,

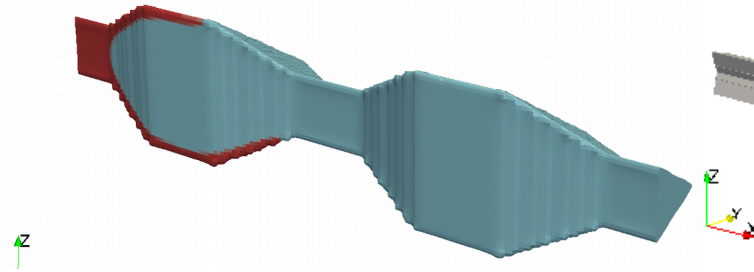


Parameters to be studied:

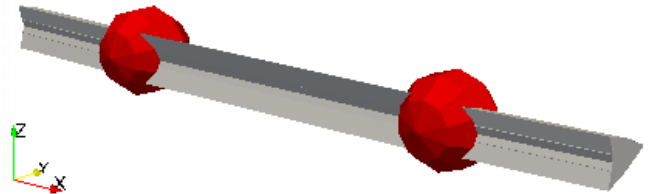
(Compute their Pc-volume-conductivity relationship)

- Contact angle
- Corner angle
- Longitudinal interface curvature
- Different pore geometries

Triangular geometry:



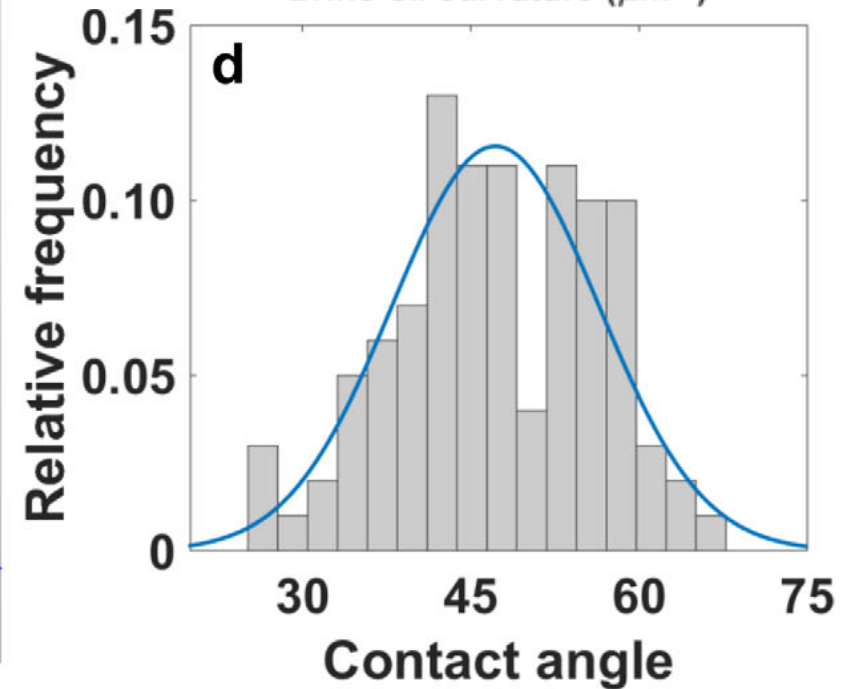
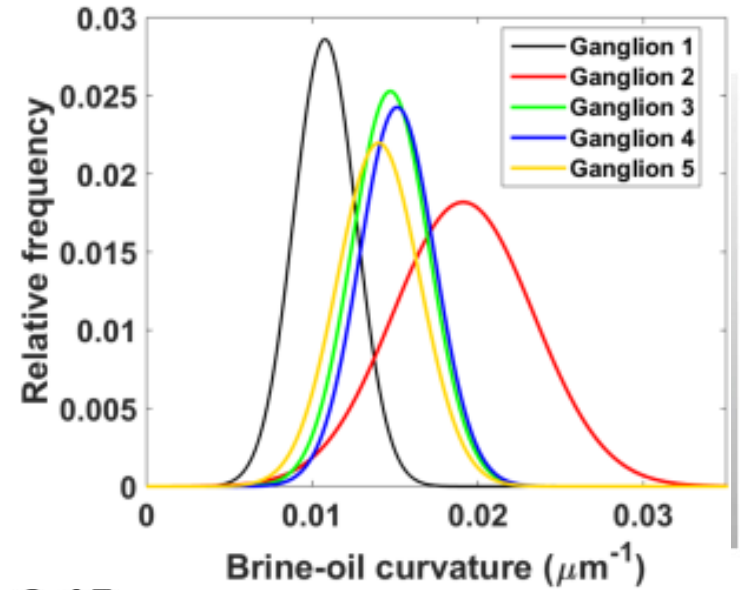
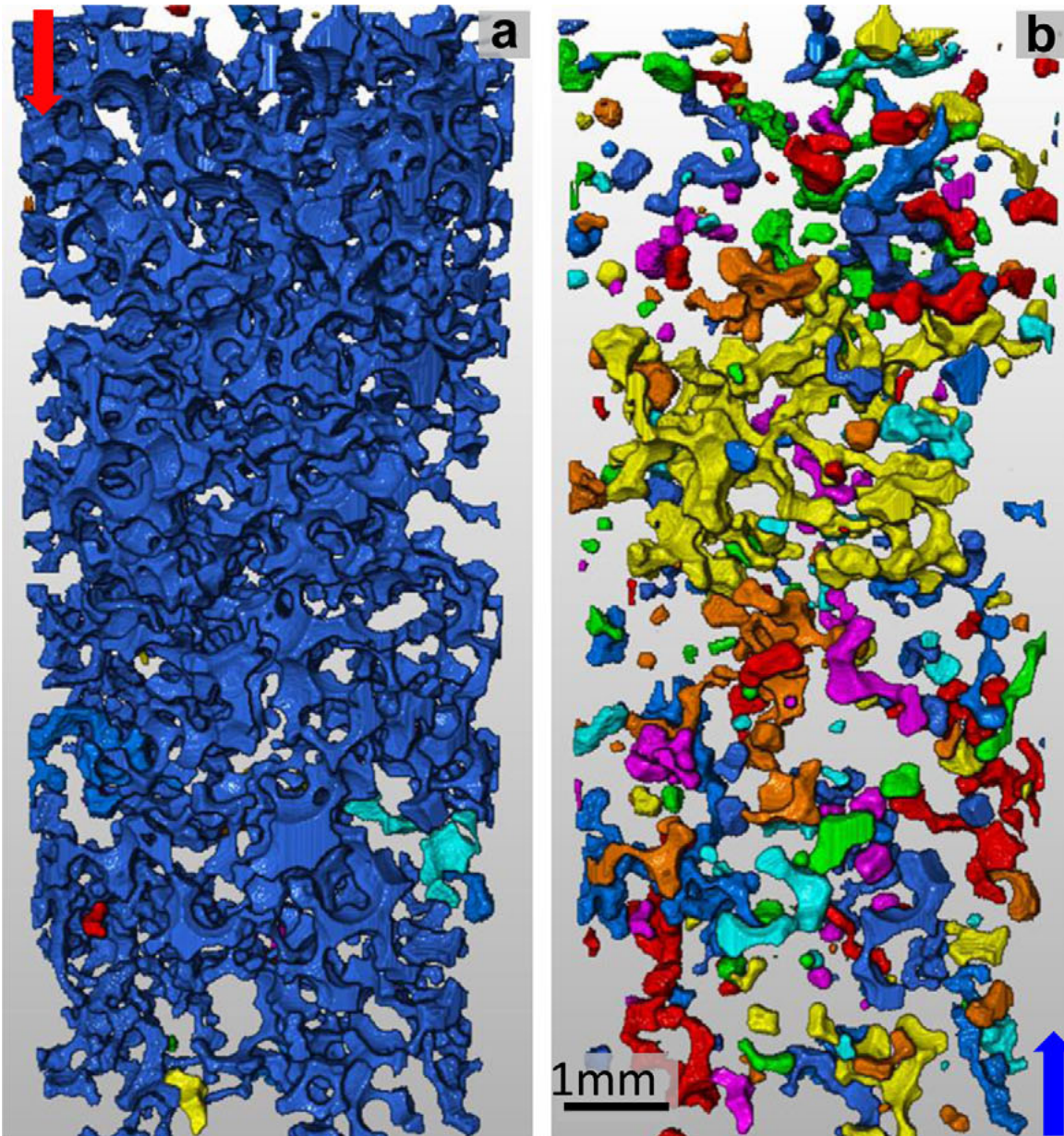
Direct simulation:



Network model:

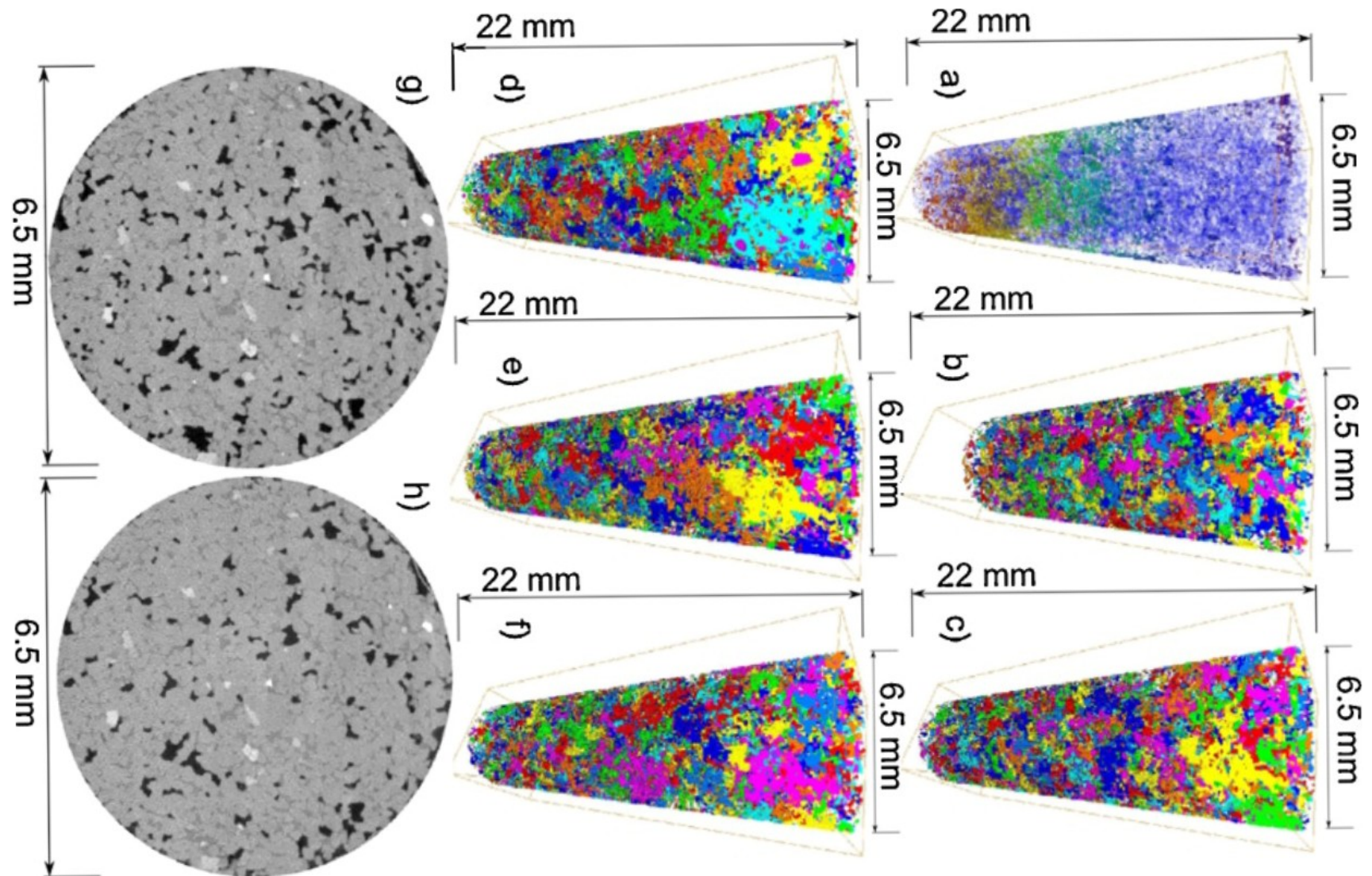
- More Complex test cases to verify cooperate-pore filling algorithm
-
- Run direct simulations on small, but high resolution micro-CT images

Curvature and contact angle measurements using high resolution micro-CT imaging



Validation using experimental measurements

Multi-phase micro-CT imaging experiments
each experiment repeated 5 times → uncertainty quantification

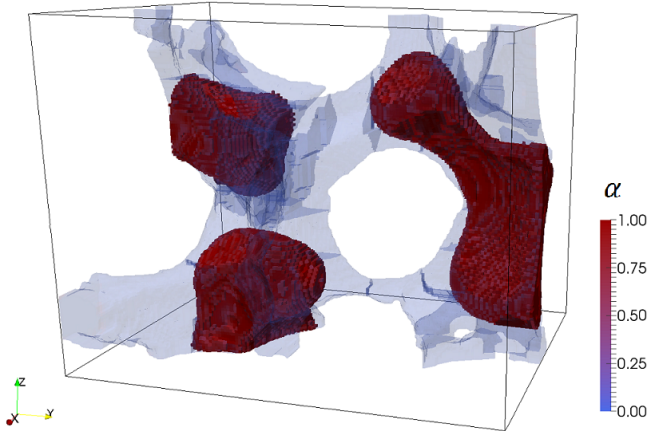
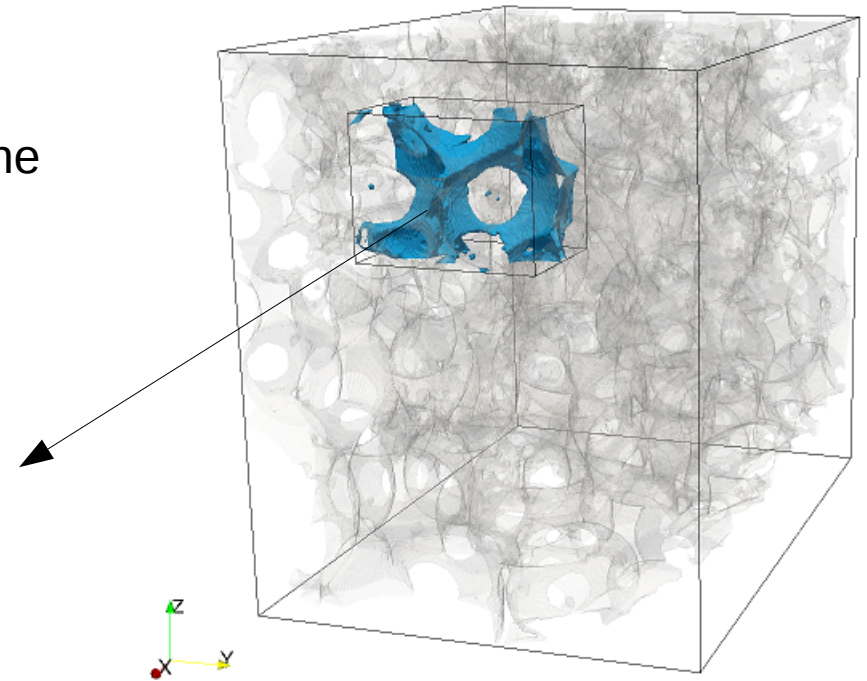


(Andrew et al, 2014)

Simulations on High resolution micro-CT images

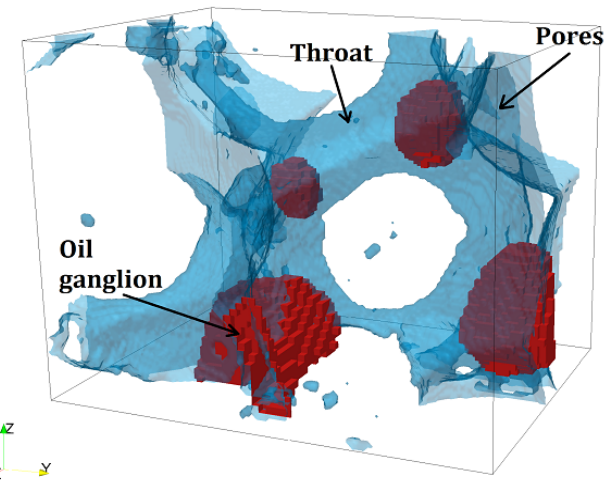
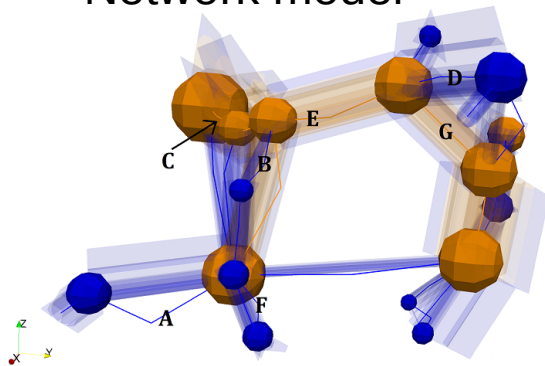
Small sample taken from the larger-scale image:

Direct simulation
(Lingru Zheng, MSc thesis, July 2016):



Multi-phase micro-CT image:

Network model



Conclusion, network modelling

- The computations are comparable in accuracy with direct simulations and in speed with network models (a gain of about 6 orders of magnitude in performance)
- ✓ Corner extraction and connectivity-tracking provides richer computations and predictions and gives us the opportunity to have a pore-by-pore comparison on complex micro-CT images
- Further work is needed to quantify the range of uncertainties/variations in experimental measurements and compare them with network model predictions

Acknowledgement

Thanks to Total for providing the
financial support

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Thank you for your attention