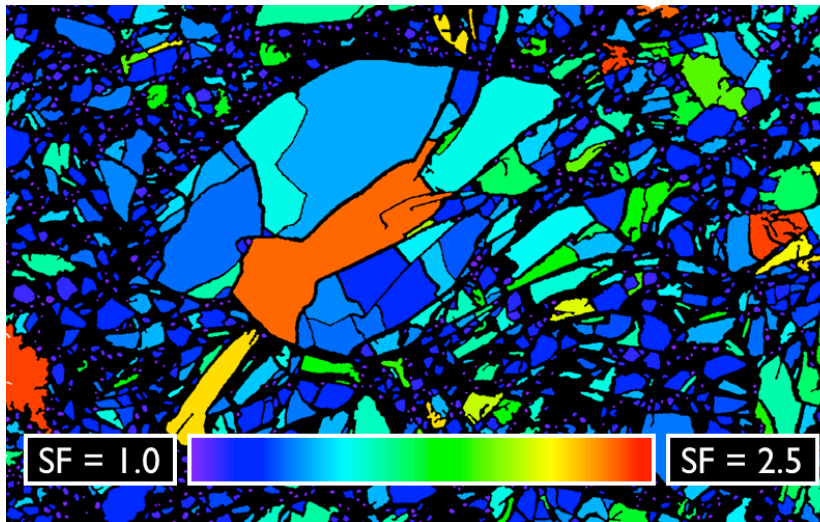


TSK19 HALLE

Workshop 2

Tuesday, March, 7, 2022



Selected topics in image analysis of deformed rocks

Shape analysis

Grain size distributions

Spatial distributions

renee.heilbronner@unibas.ch

(3)

Schedule

Renée – Lectures

- 10:00-10:30 shape analysis
- 10:30-11:00 discussion & break
- 11:00-11:30 grainsize
- 11:30-12:00 discussion & break
- 12:00-12:30 **phase distributions & correlations**
- 12:30-14:00 discussion & lunch

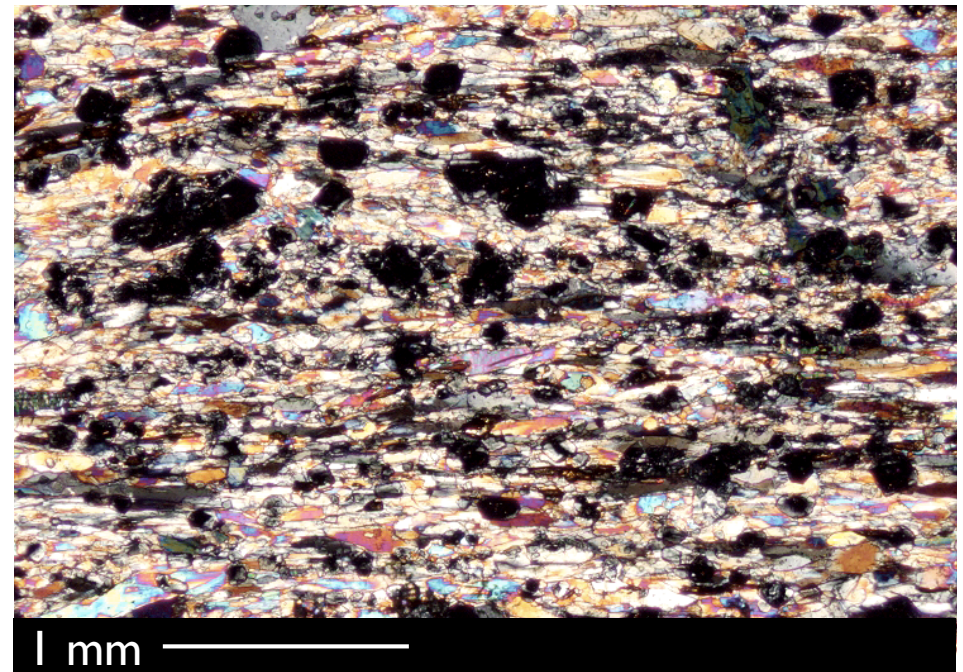
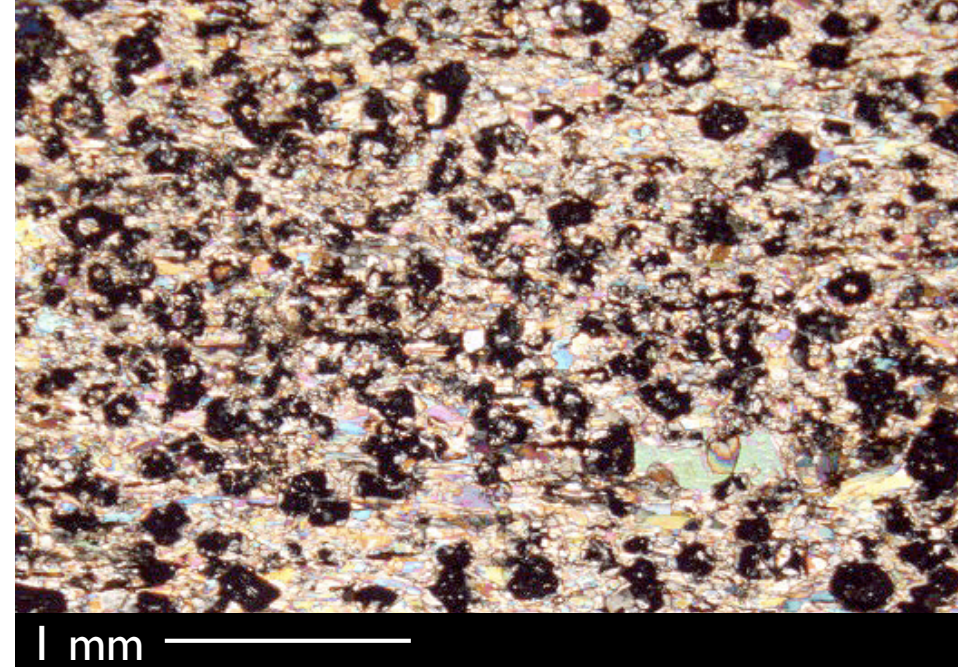
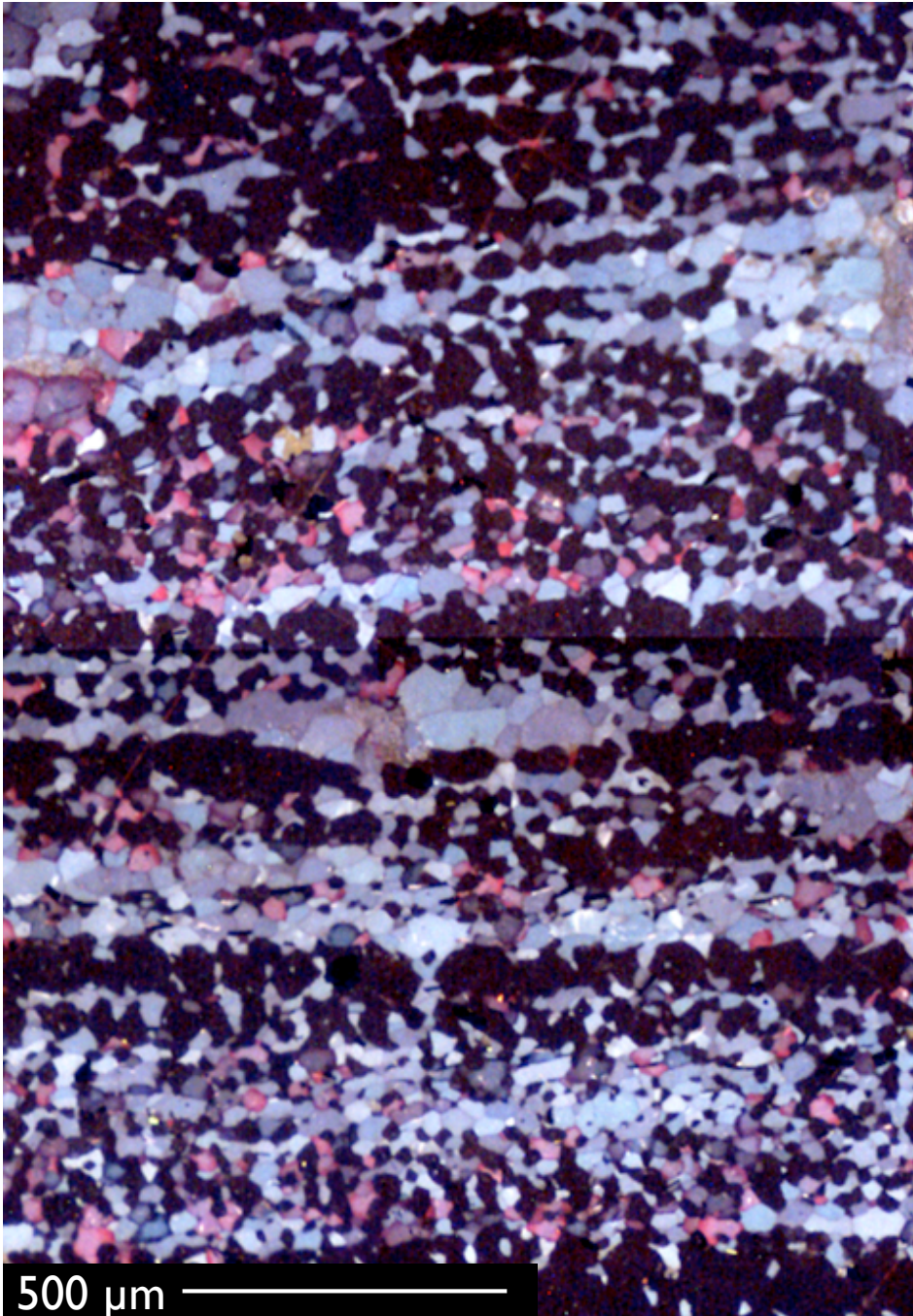
Rüdiger – Lab

- 14:00-15:30 using Fiji / imagej
- 15:30-16:00 break
- 16:00-17:00 playtime (with your own data)

|

spatial distributions

spatial dispersion

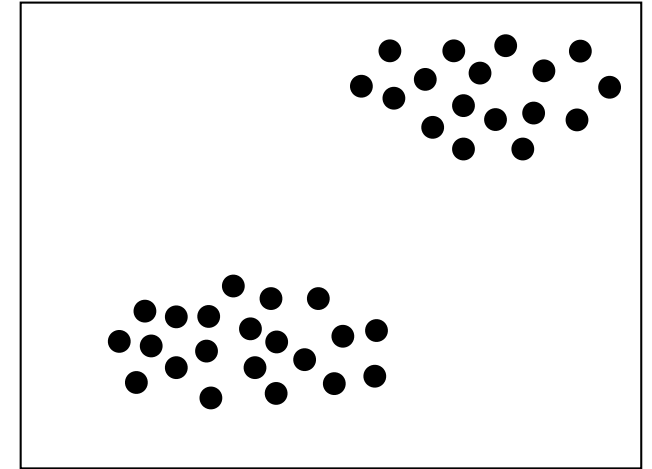
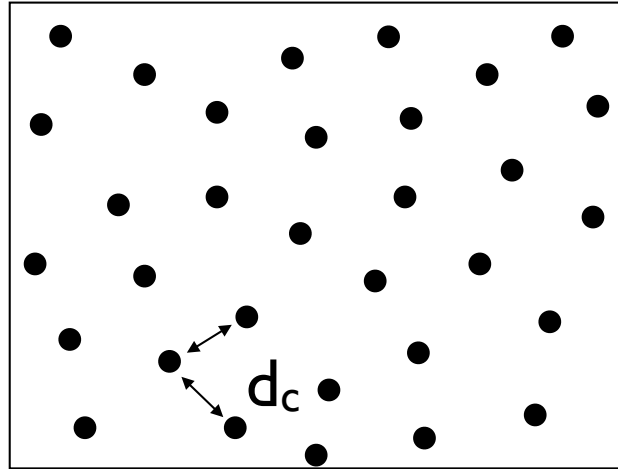
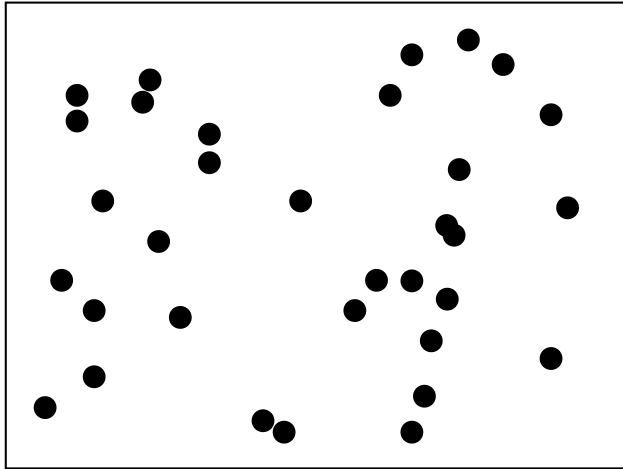


spatial distributions

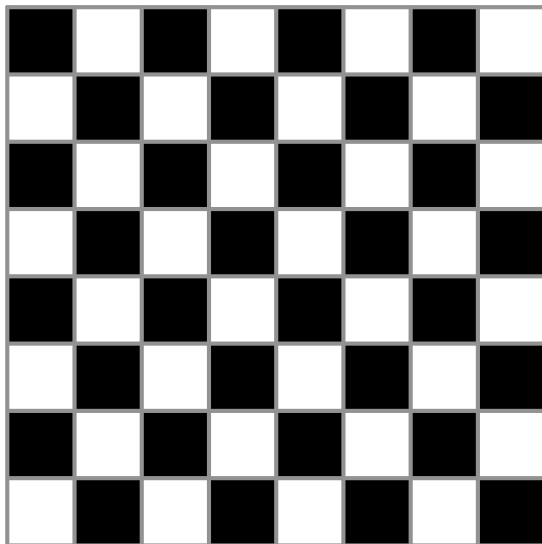
- random
- clustered
- ordered ('anti-clustered')

spatial distributions

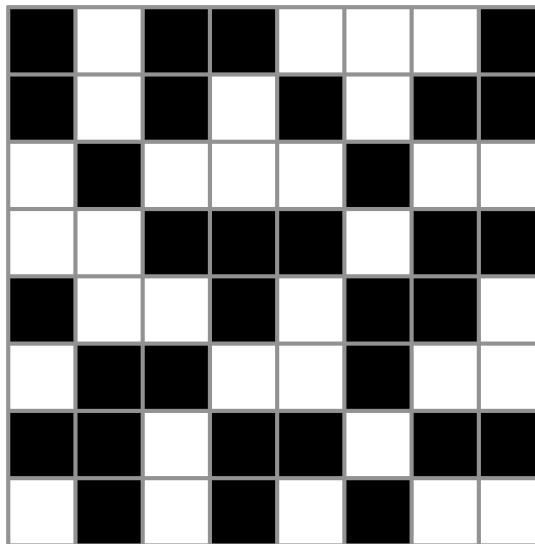
spatially dispersed center points



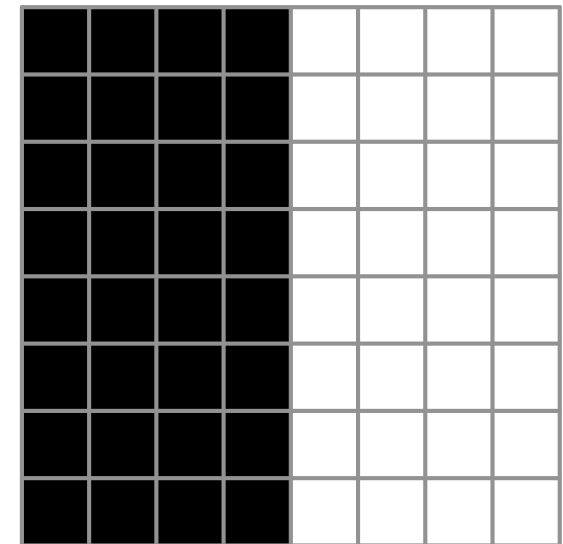
chessboard model



ordered

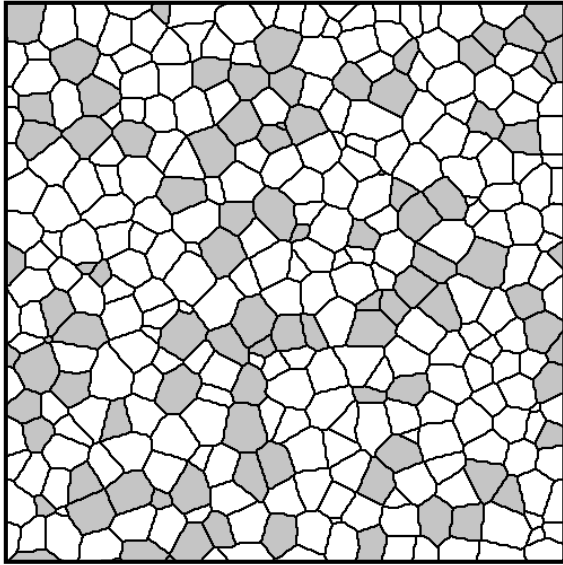


random

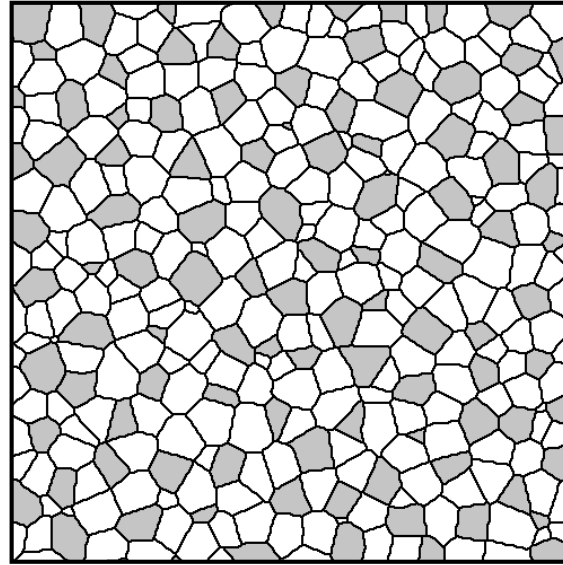


clustered

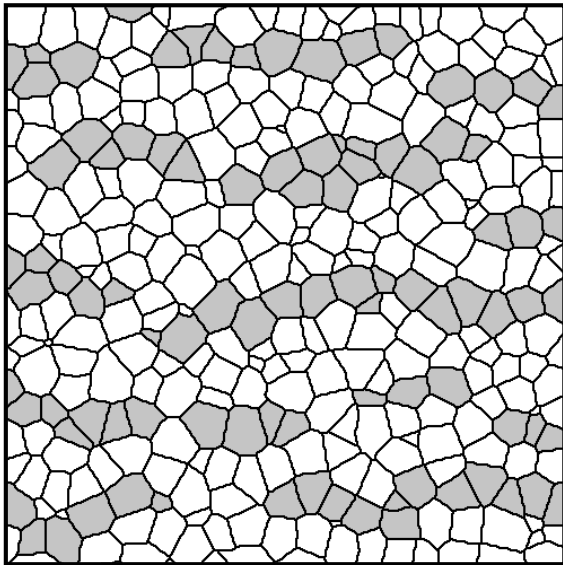
spatial distribution



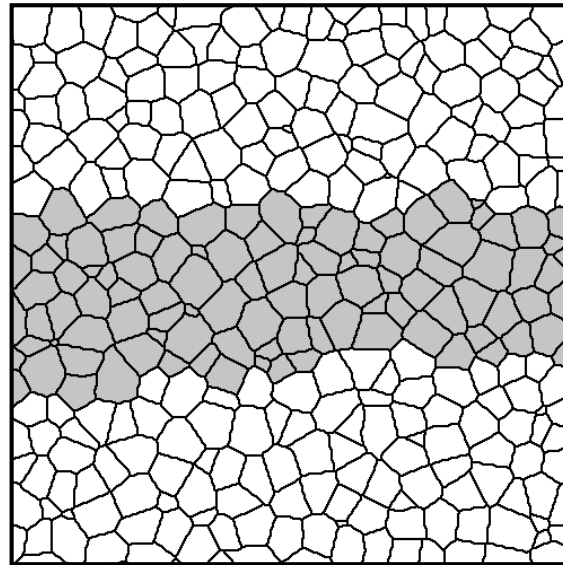
random



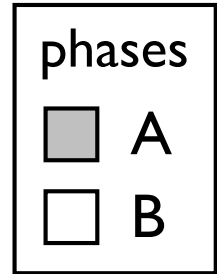
ordered



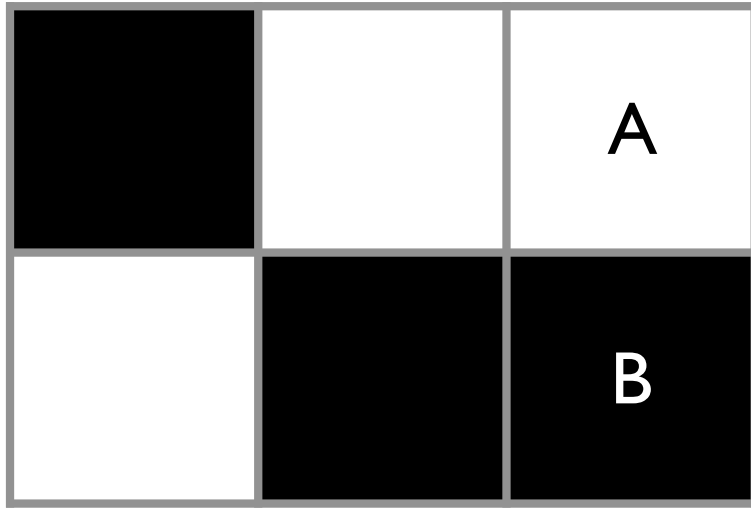
clustered



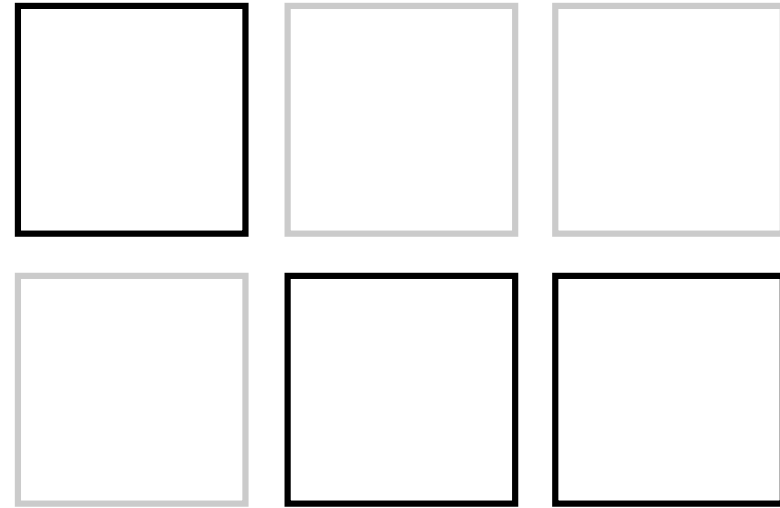
highly
clustered



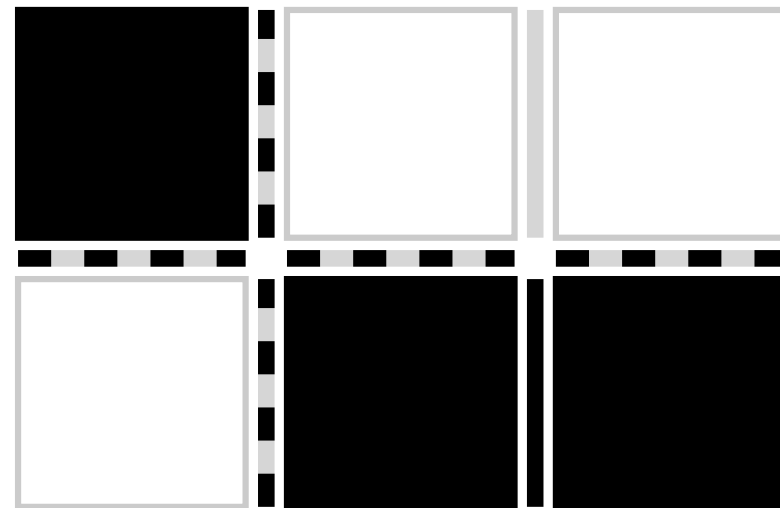
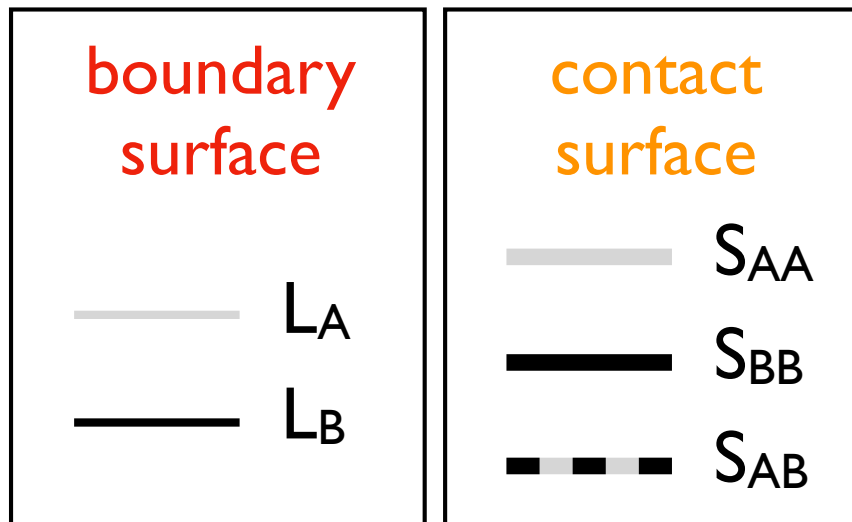
phases, boundaries and contacts



2 types of grains

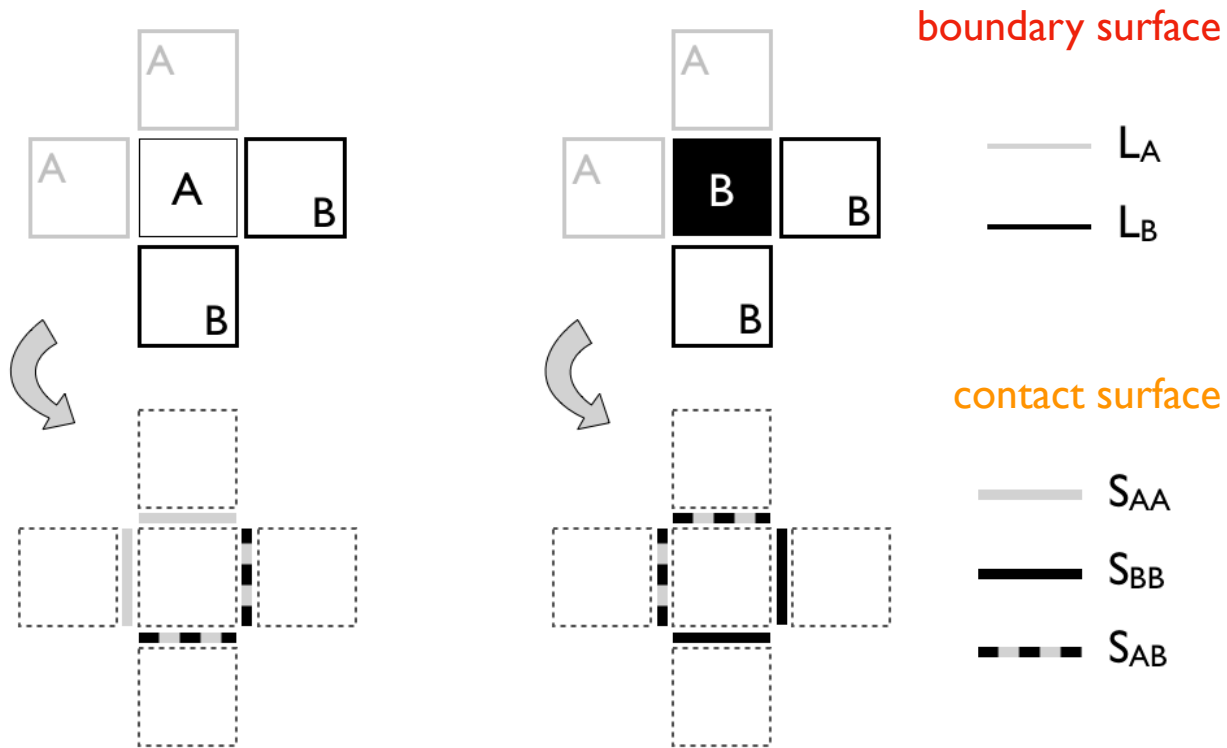


2 types of grain boundary surface



3 types of grain contact surface

conceptual model



For any given grain A, B:
the chance p_A , p_B , of sharing a boundary surface with a grain A or B is proportional to:

- volume fraction¹⁾ of A, B

or

- surface fraction²⁾ of A, B

$$AA = p_A \cdot p_A = p_A^2$$

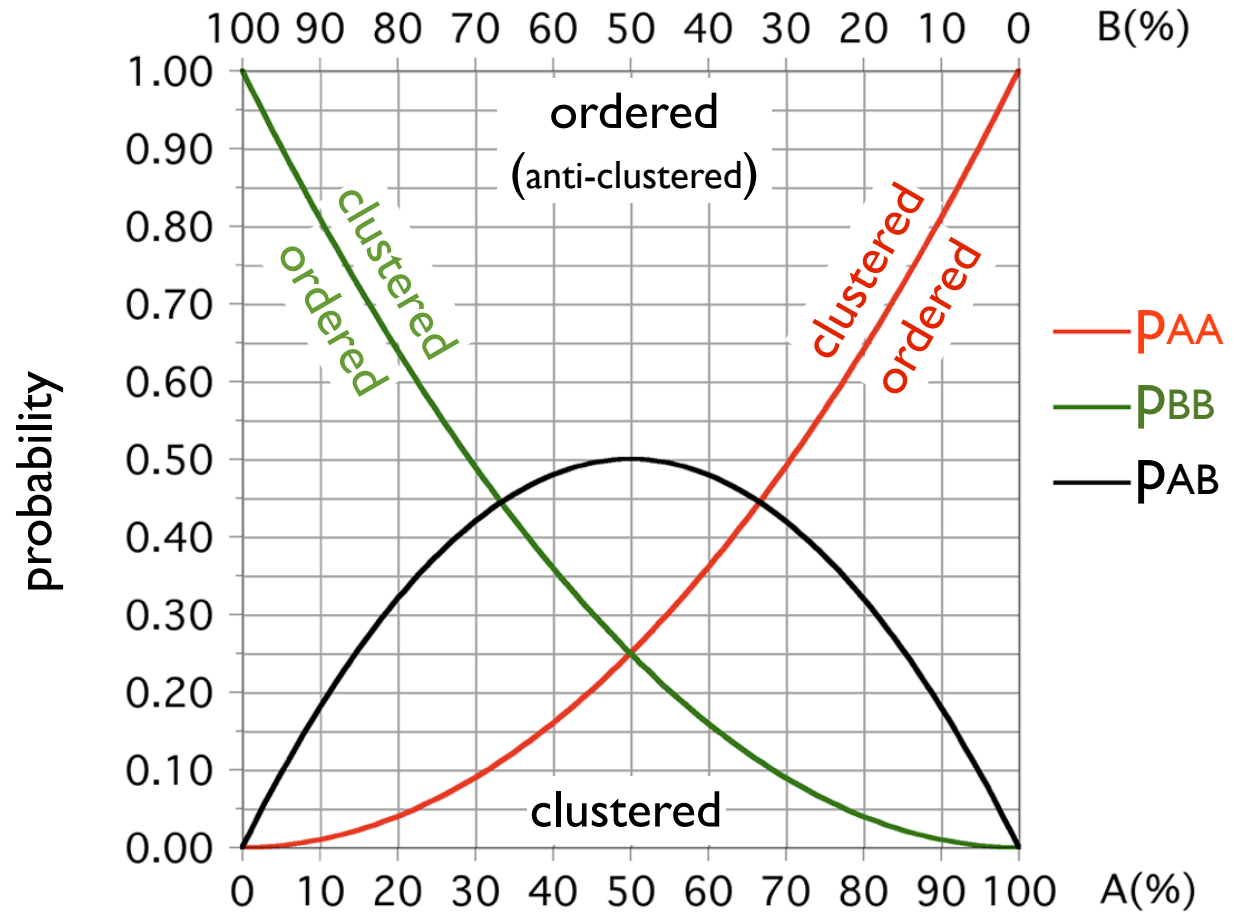
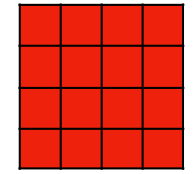
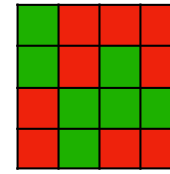
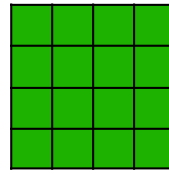
$$BB = p_B \cdot p_B = (1-p_A)^2$$

$$AB = BA = p_A \cdot p_B + p_B \cdot p_A = 2 \cdot p_A \cdot (1-p_A)$$

1) = area(phase)
/ (total cross sectional area)

2) = outline(phase)
/ (total boundary length)

binomial distribution



$$AA = p_A \cdot p_A = p_A^2$$

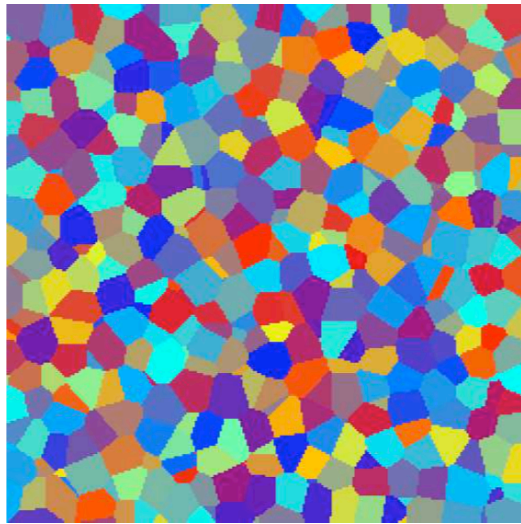
$$BB = p_B \cdot p_B = (1-p_A)^2$$

$$AB = BA = p_A \cdot p_B + p_B \cdot p_A = 2 \cdot p_A \cdot (1-p_A)$$

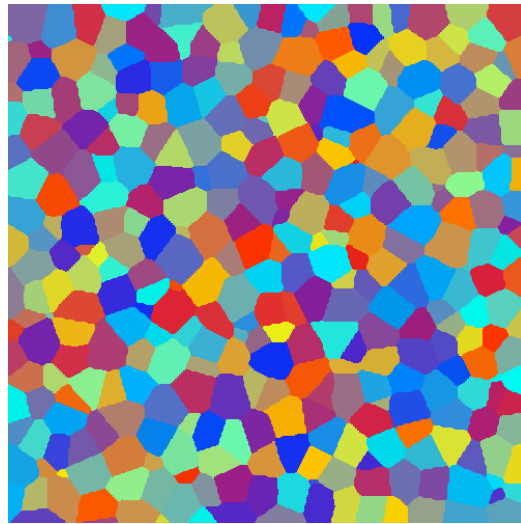
spatial distributions

- grains in matrix
- grains in crystalline aggregate

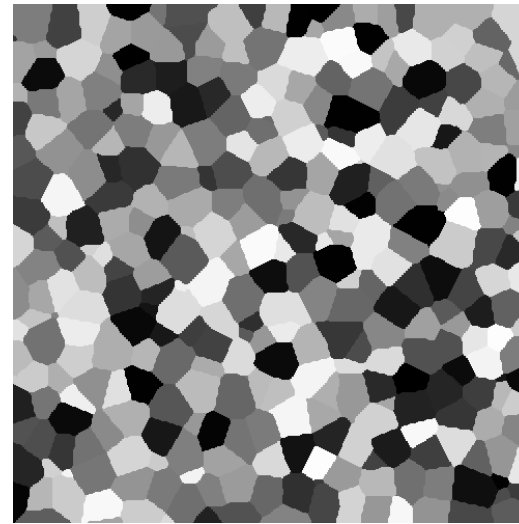
making random Voronois



5050anti5_z5.tif



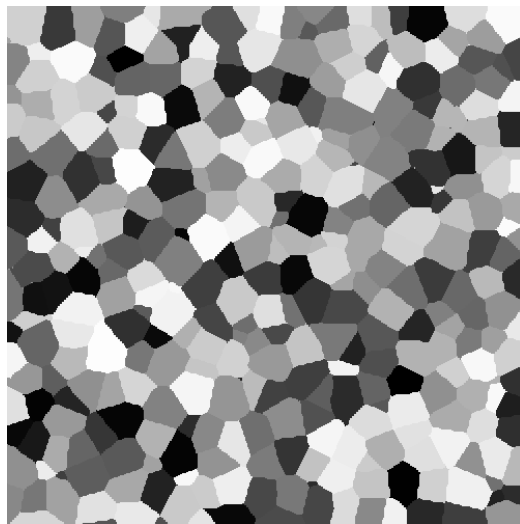
5050anti5Prep



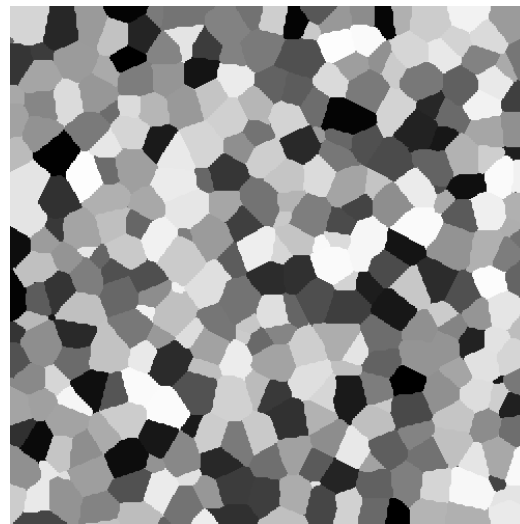
5050anti5Prep (R)



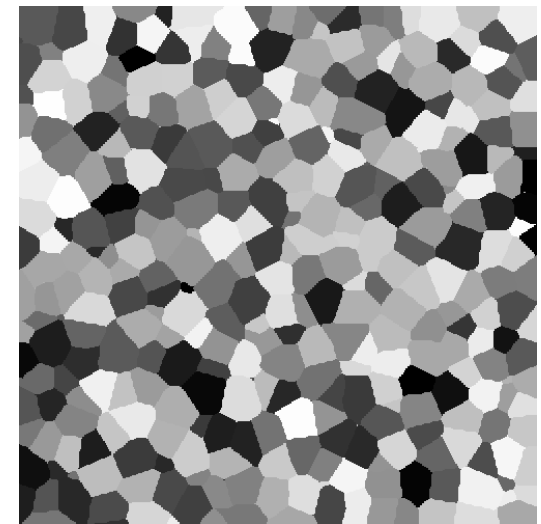
re-apply
random
Palette



5050anti5Prep R1



5050anti5Prep R2

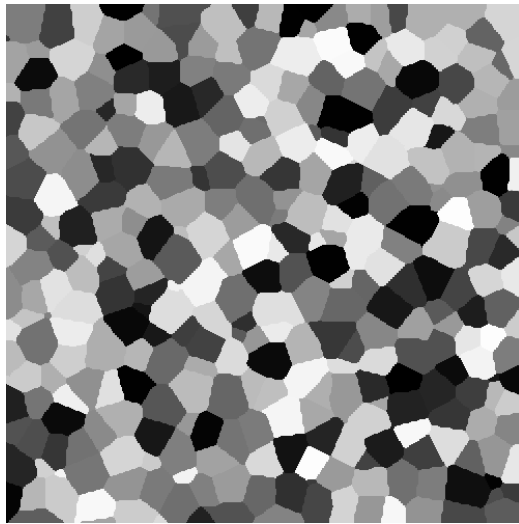


5050anti5Prep R3

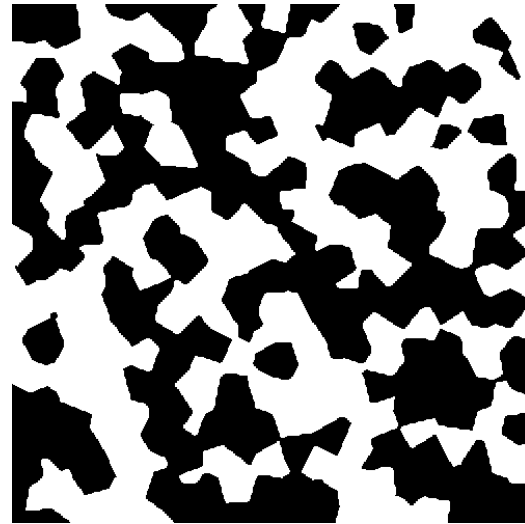
etc.

used random Palette on 5050anti5Prep R = processed R of 5050anti5_z5.tif

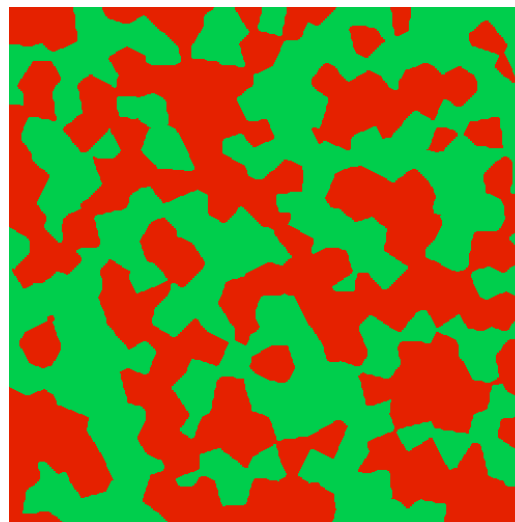
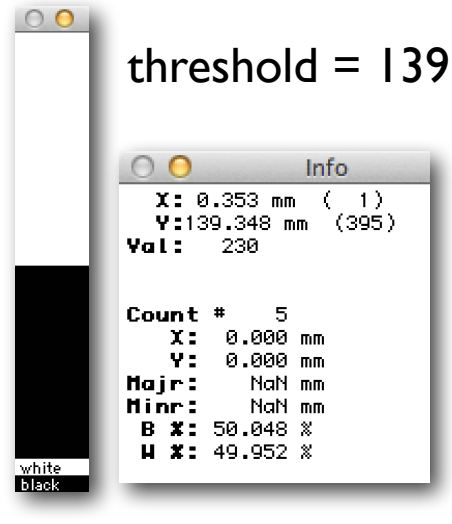
making random Voronois (continued)



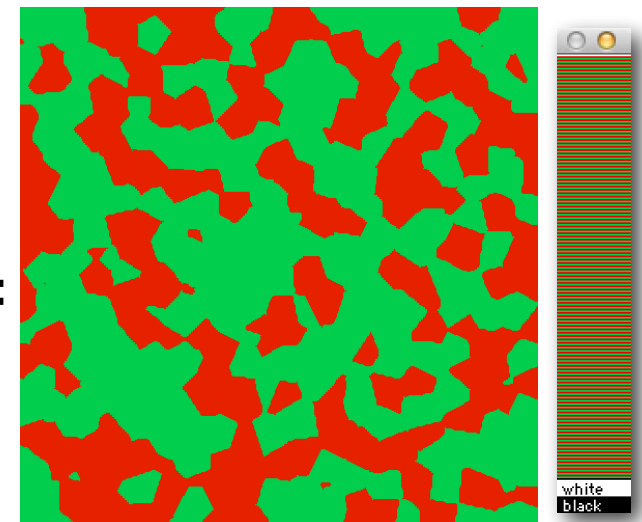
5050anti5Prep (R)



5050anti5Prep (R)



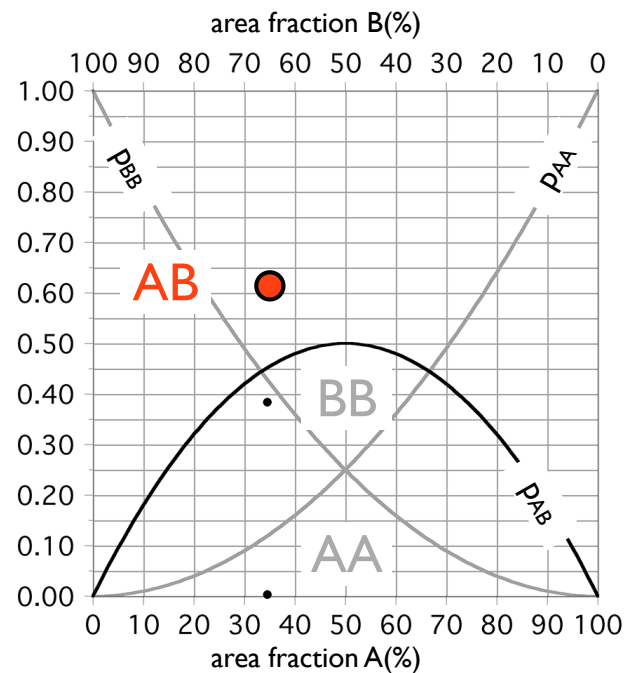
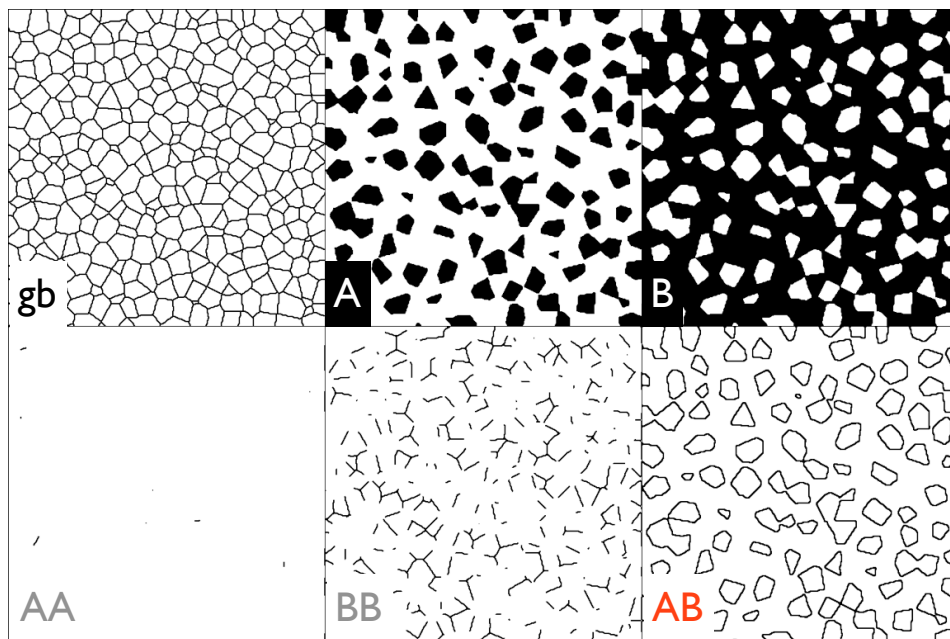
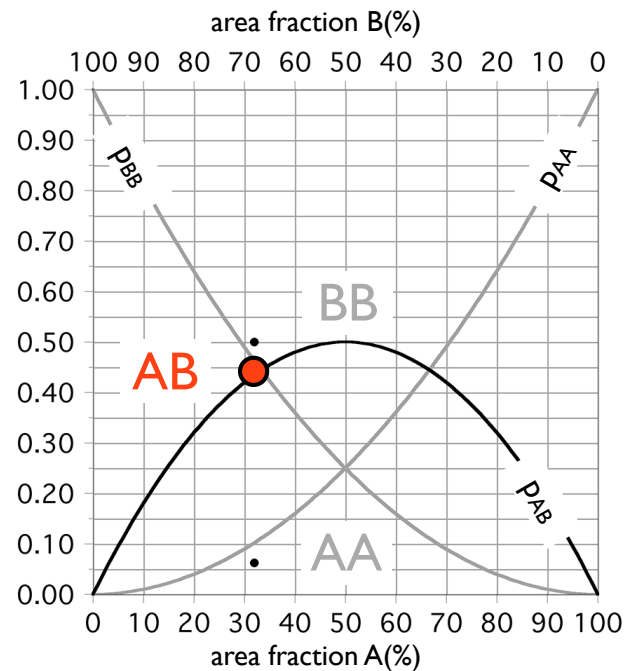
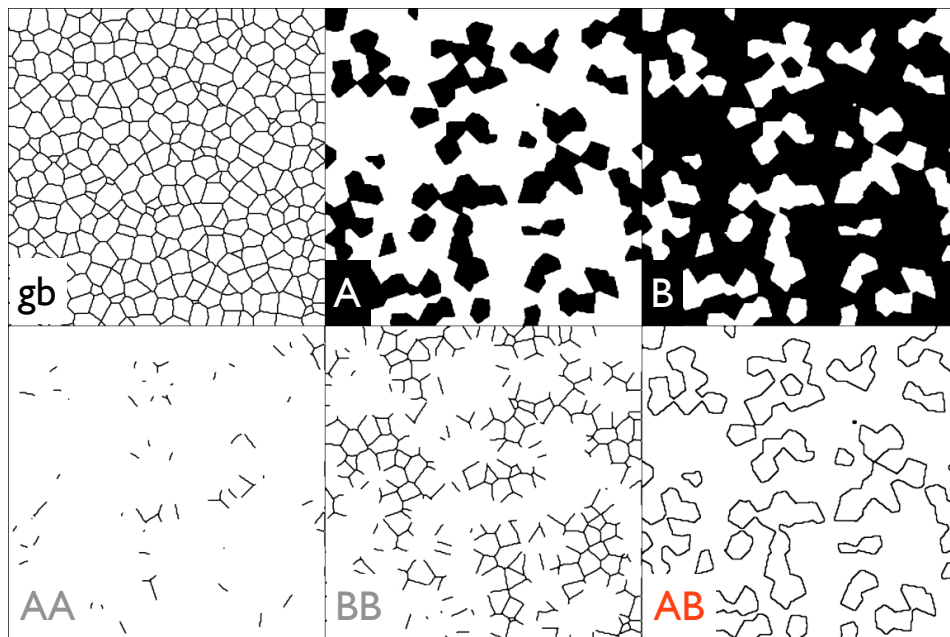
R 50.0% G 50.0%



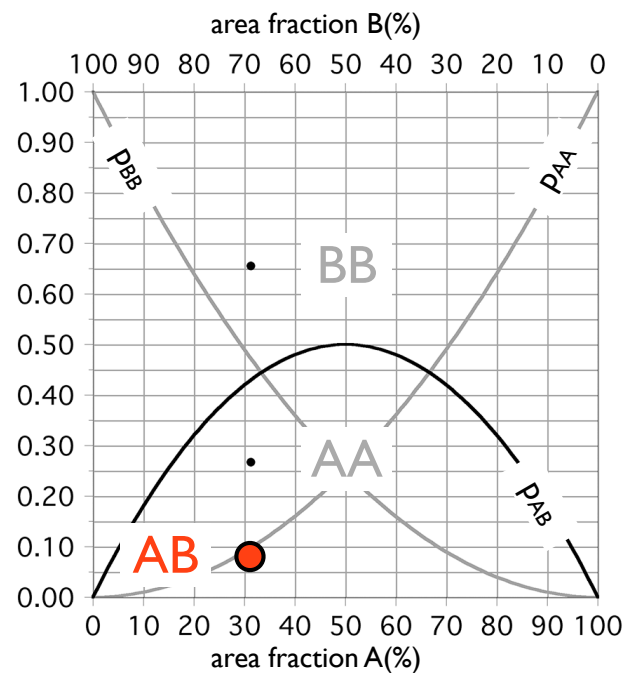
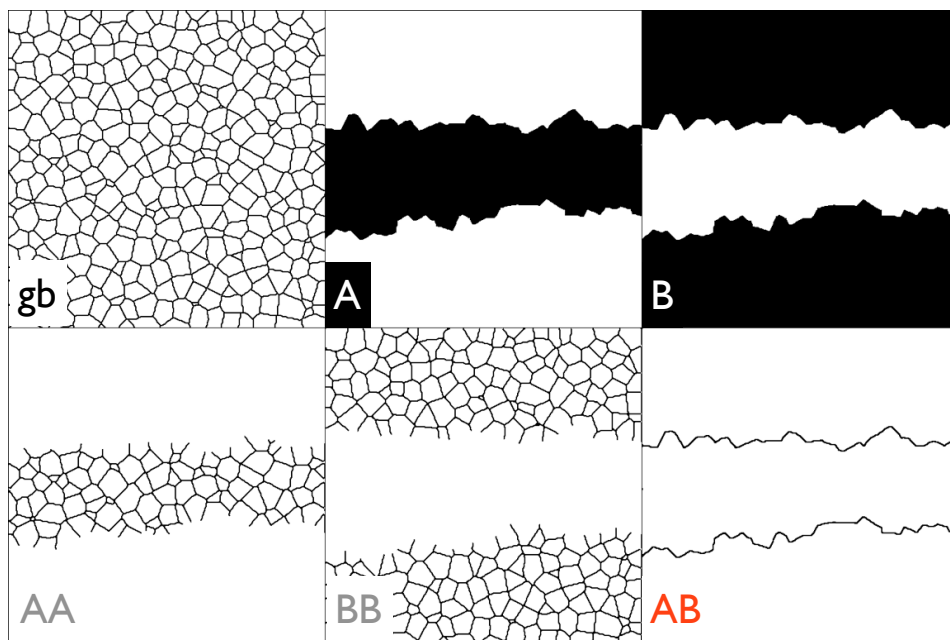
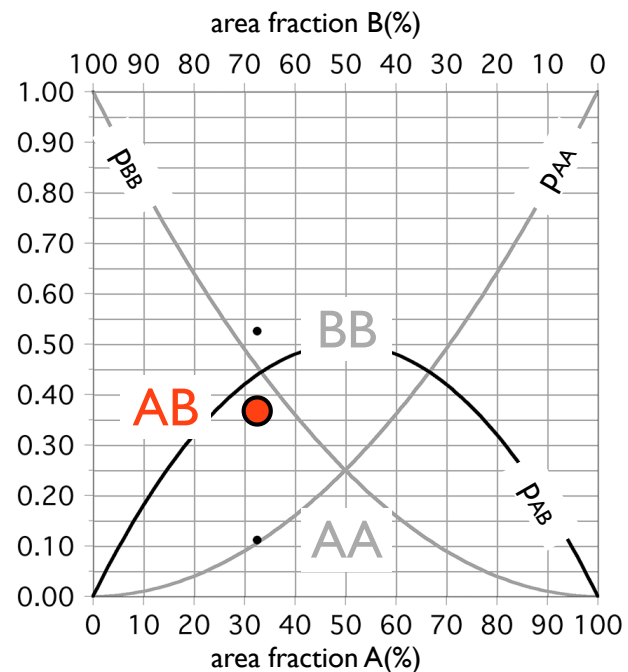
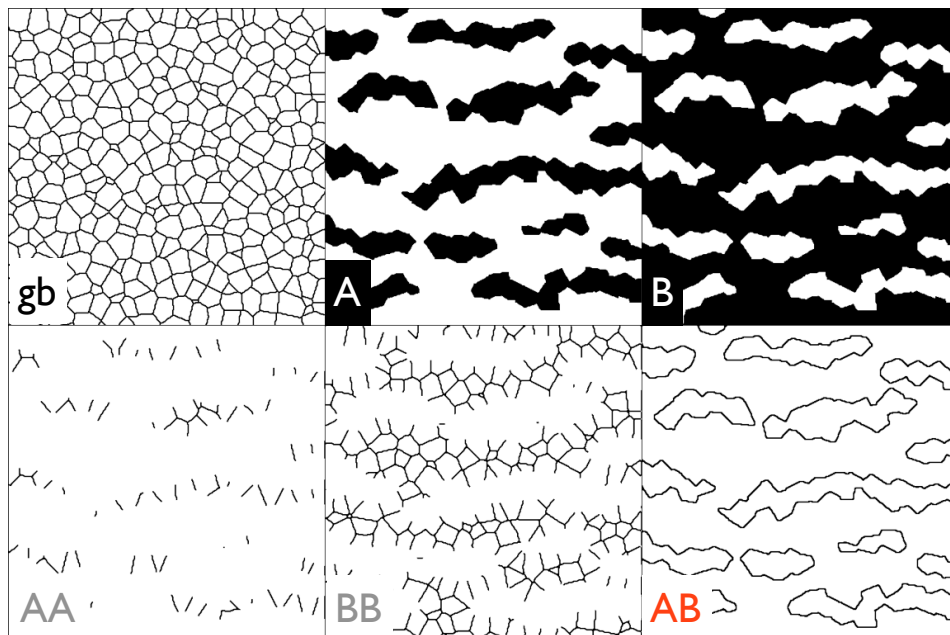
R 46.8% G 52.2%

threshold 5050anti5Prep instead of using macro [6], \Rightarrow control percentage

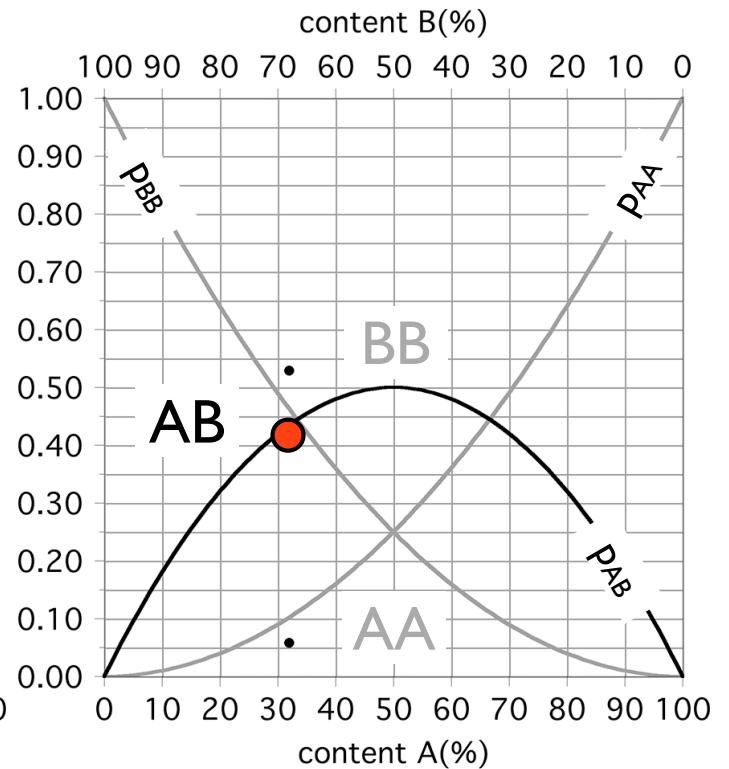
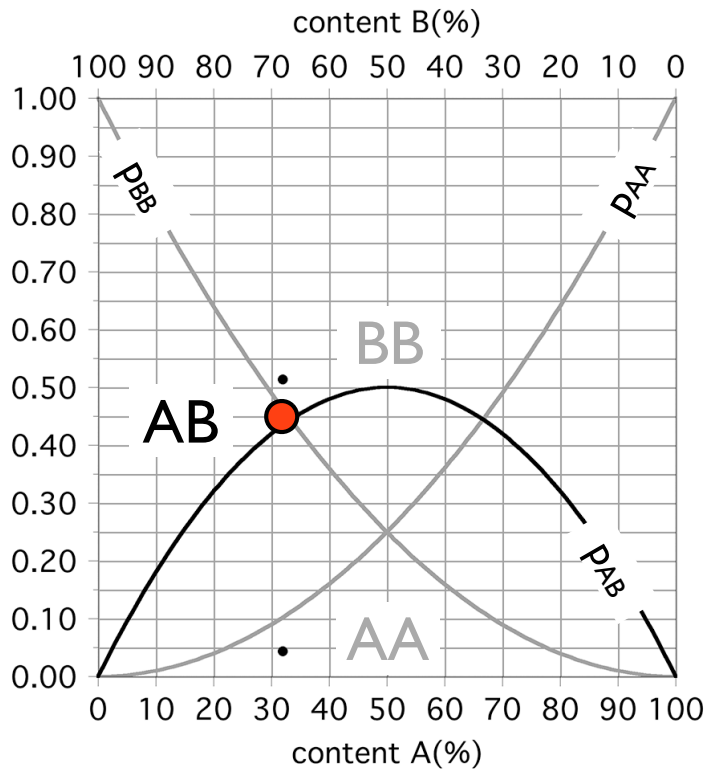
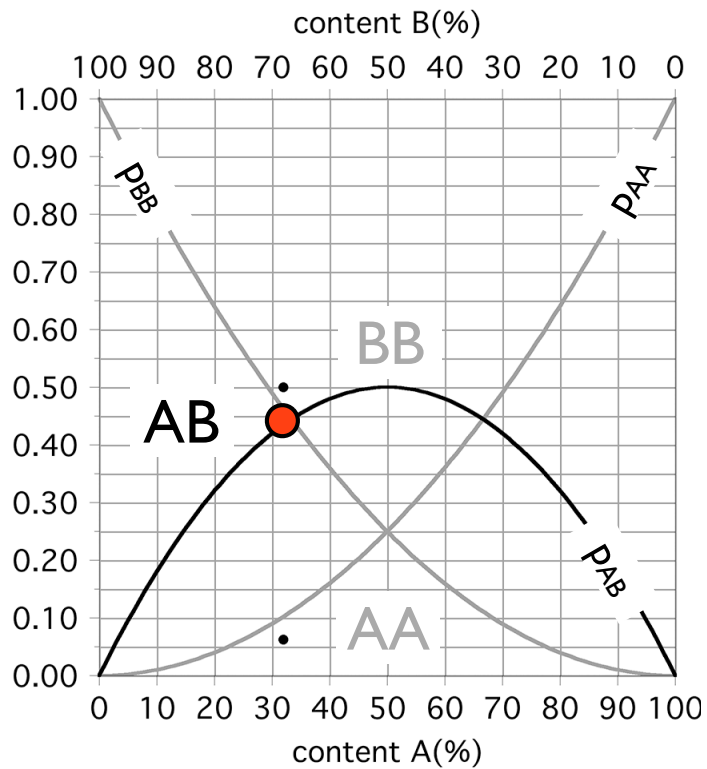
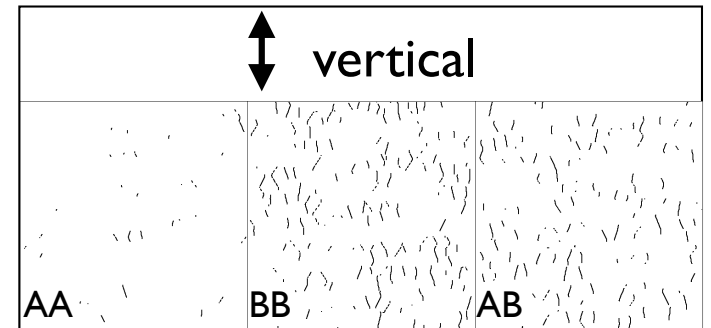
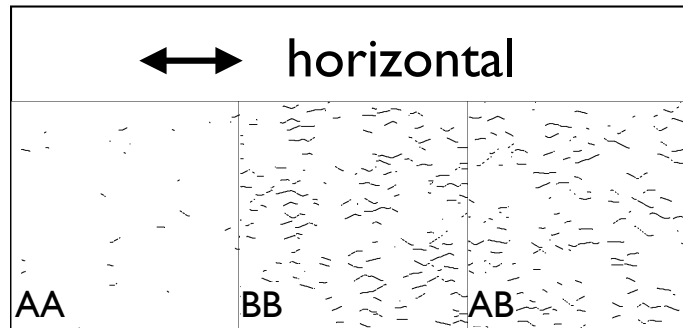
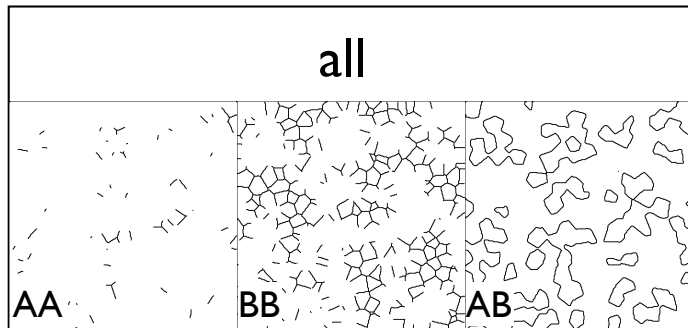
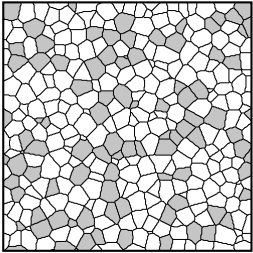
distributions and contact probabilities



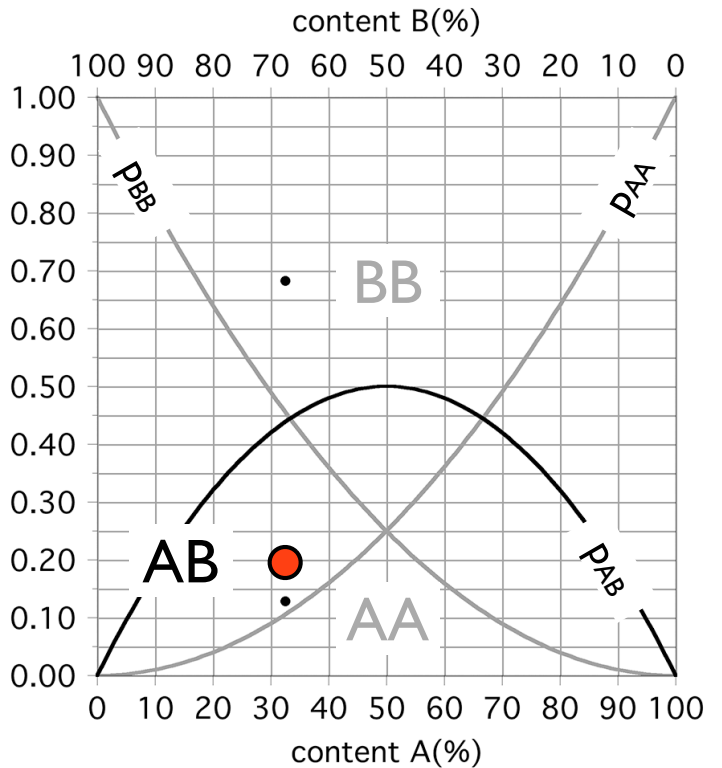
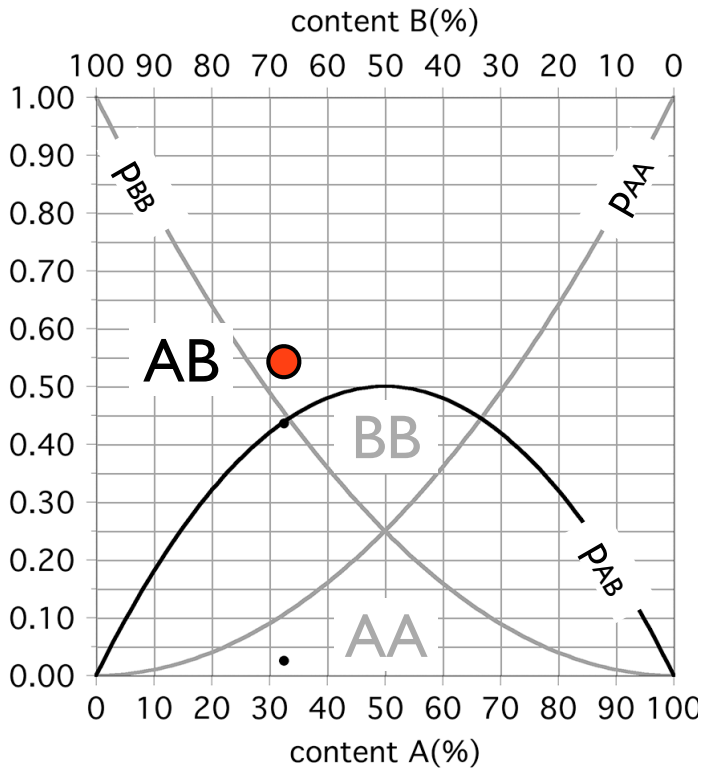
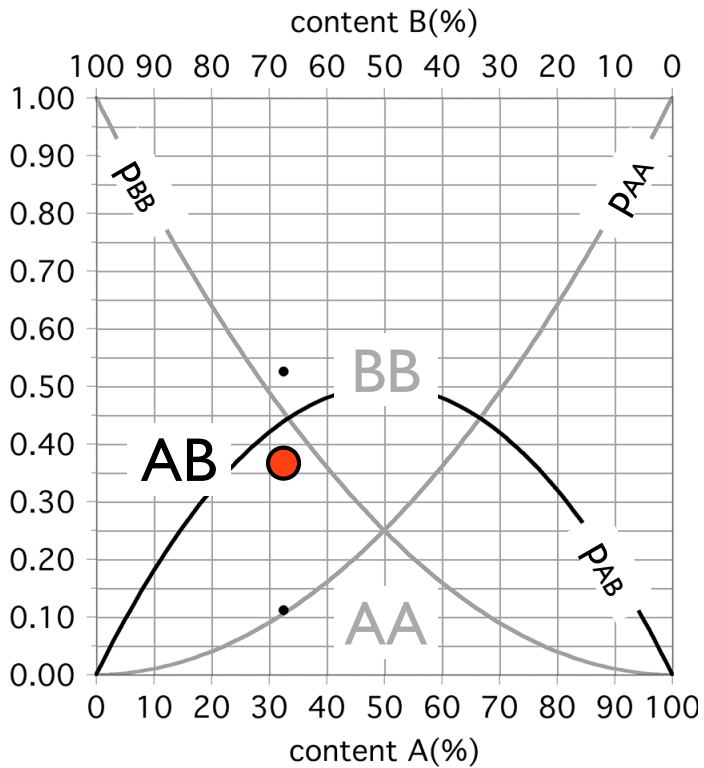
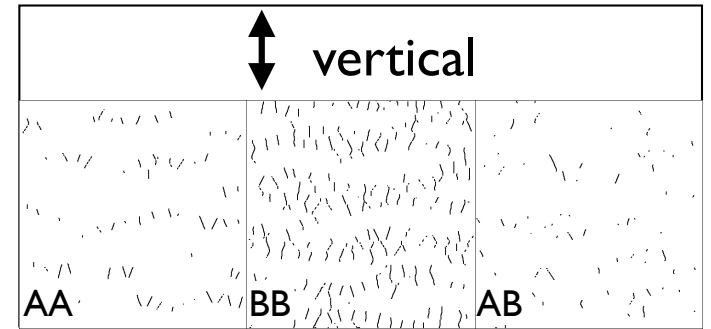
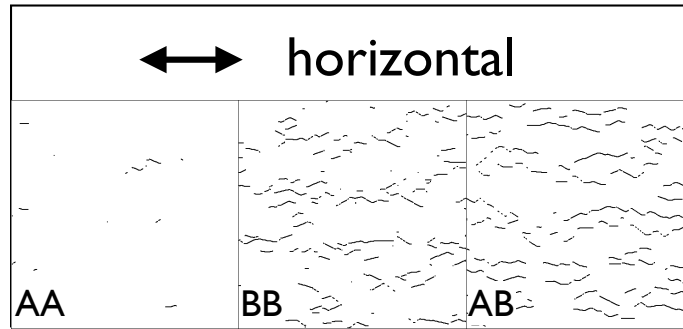
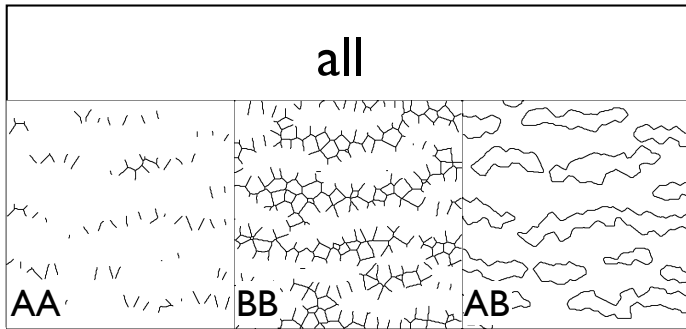
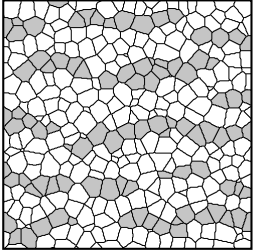
distributions and contact probabilities



isotropic random distribution

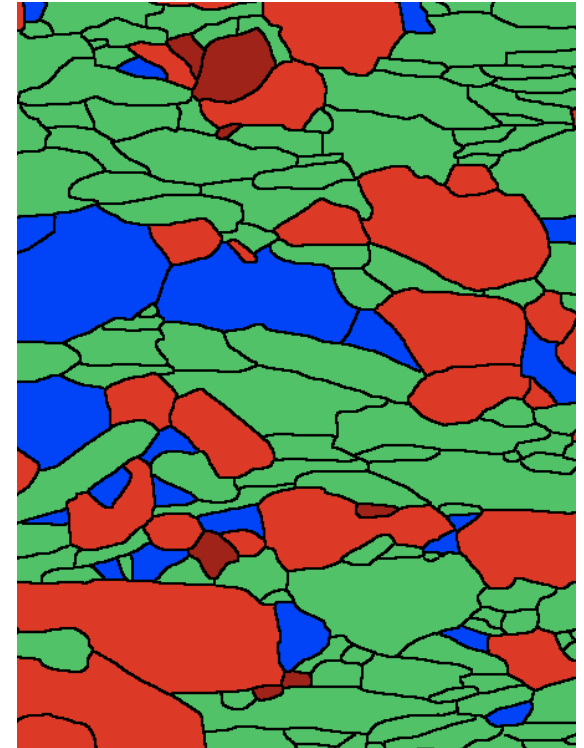
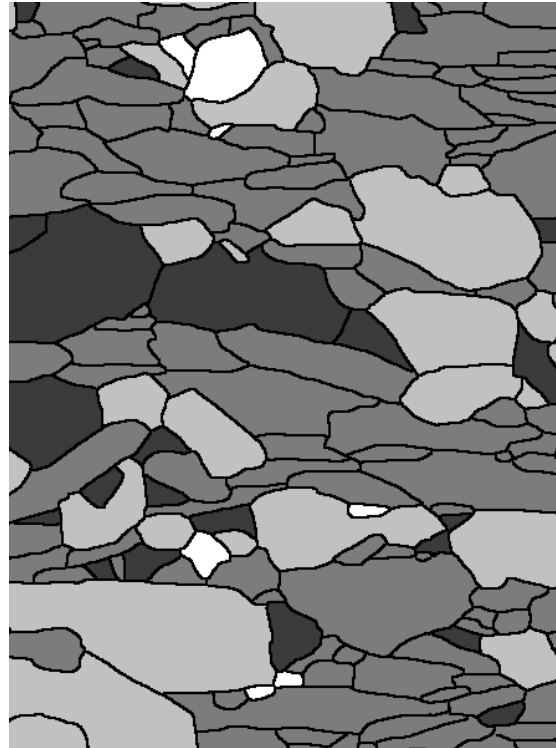
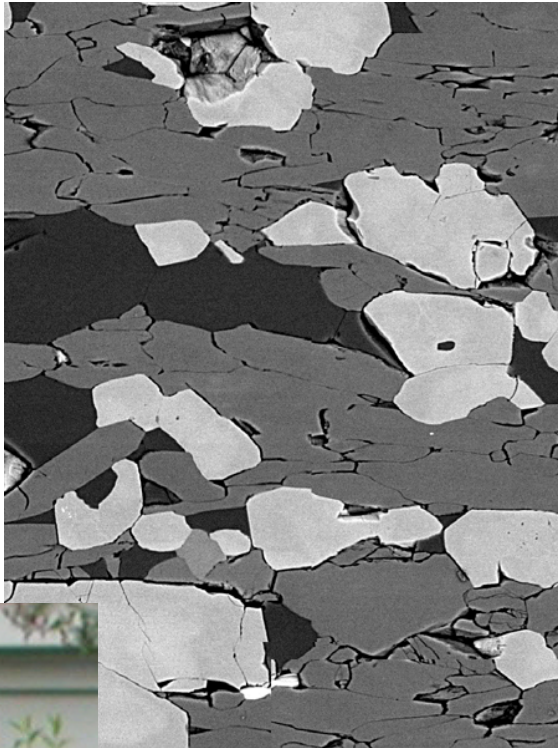


anisotropic random distribution











practical application

Meluzina eclogite

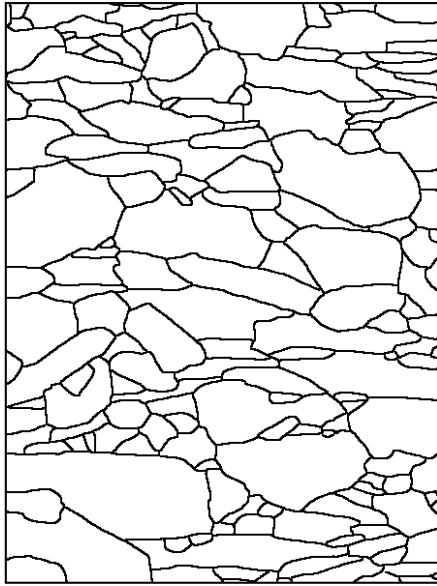


James Mackenzie

-  unidentified
-  garnet
-  omphacite
-  quartz

-  unidentified
-  garnet
-  omphacite
-  quartz

deriving phase and grain boundaries



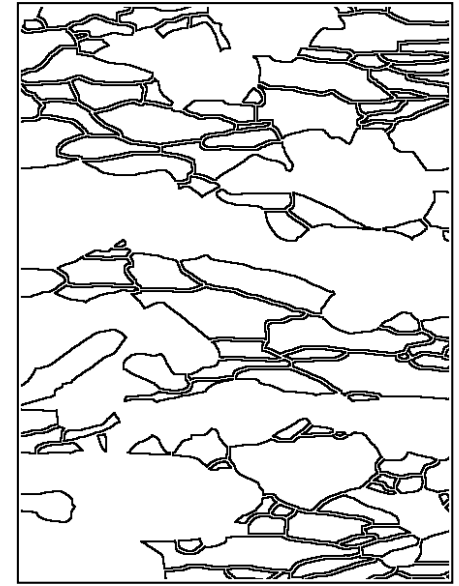
grain boundaries



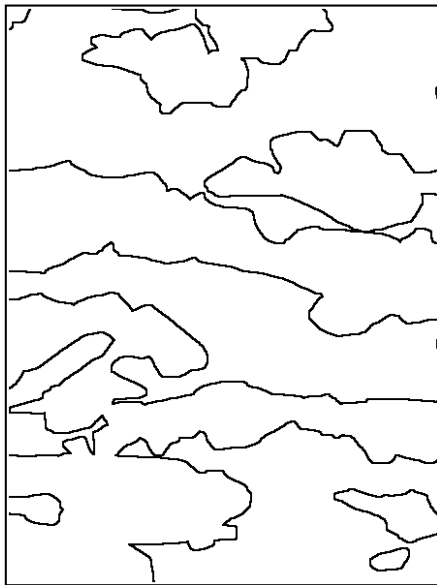
A grains



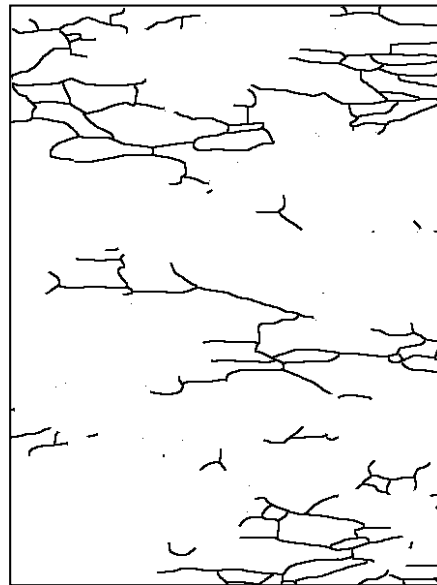
A phase



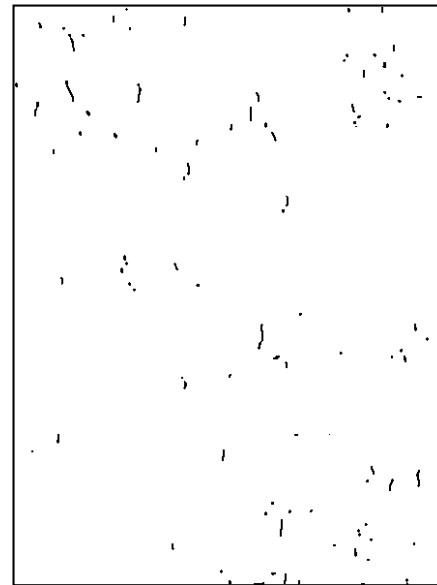
A grain outlines



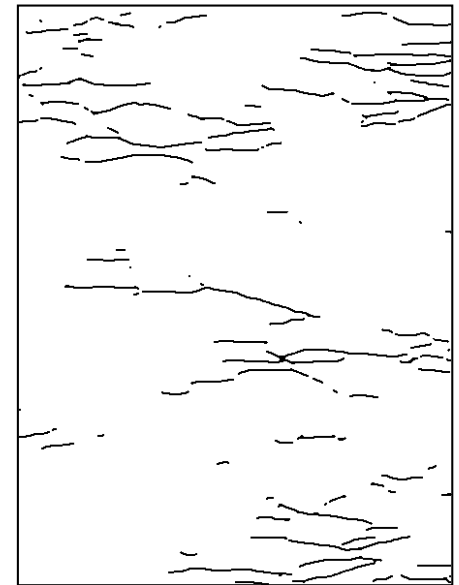
A phase outline



A grain boundaries

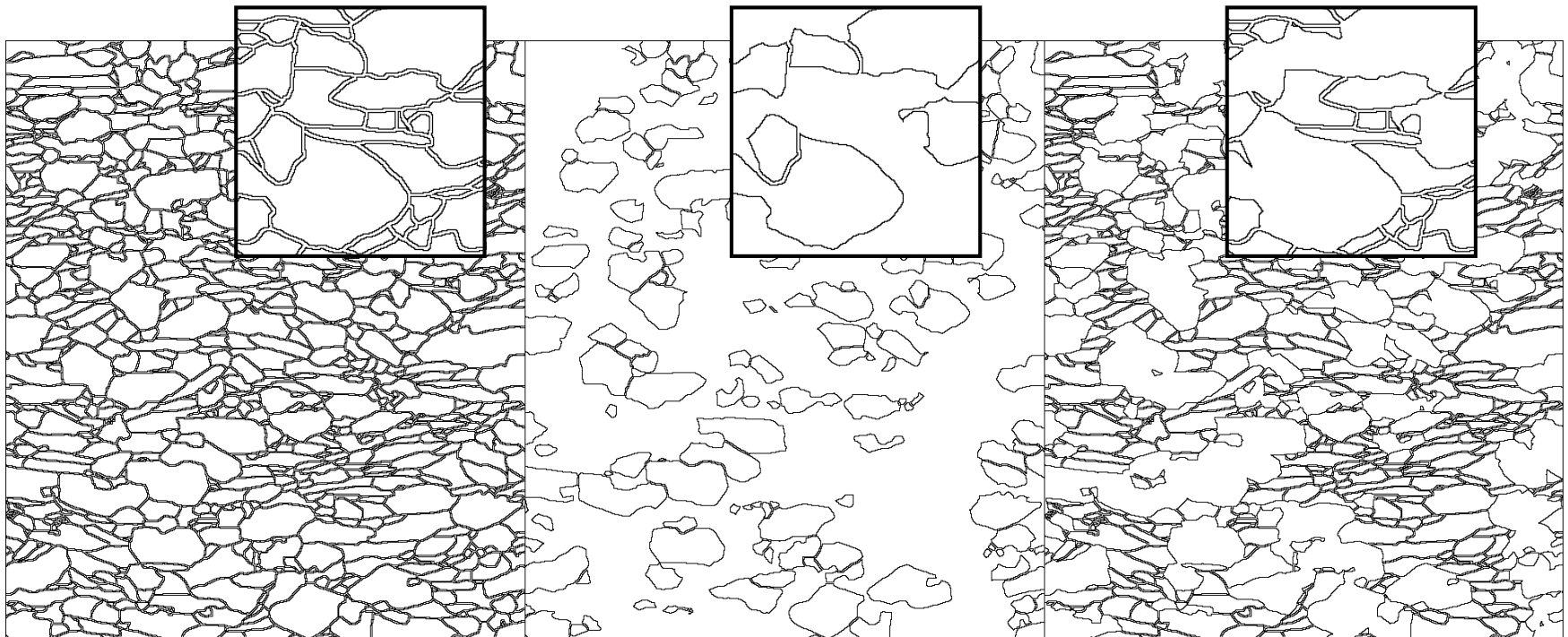
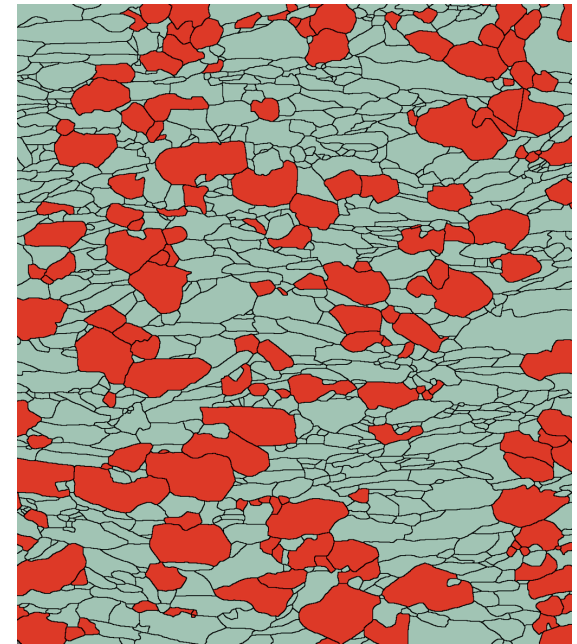
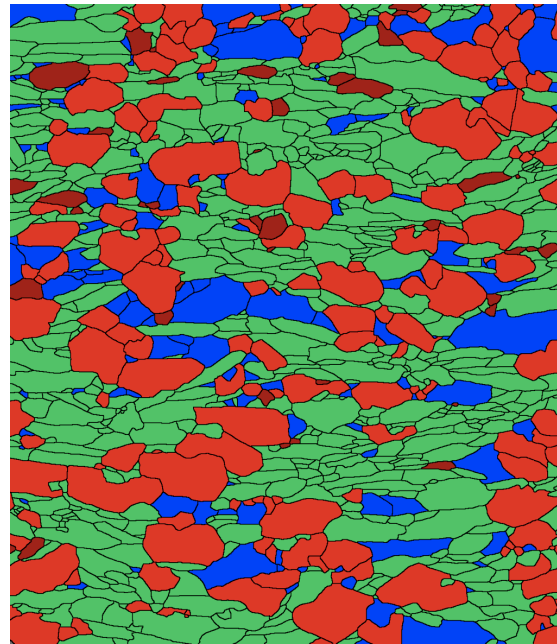
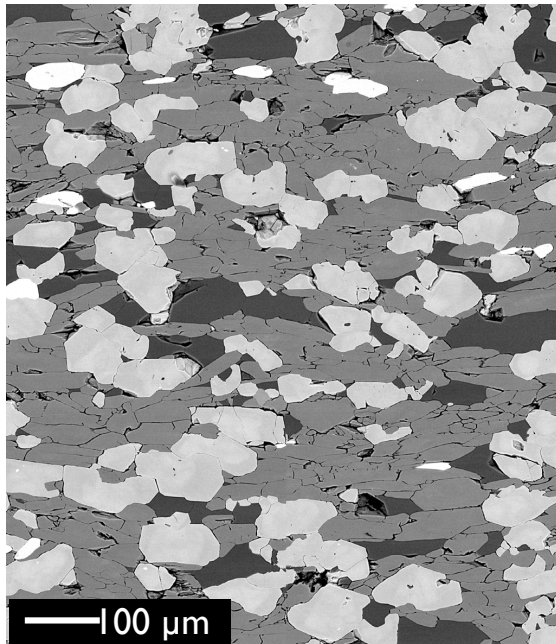


vertical A gb

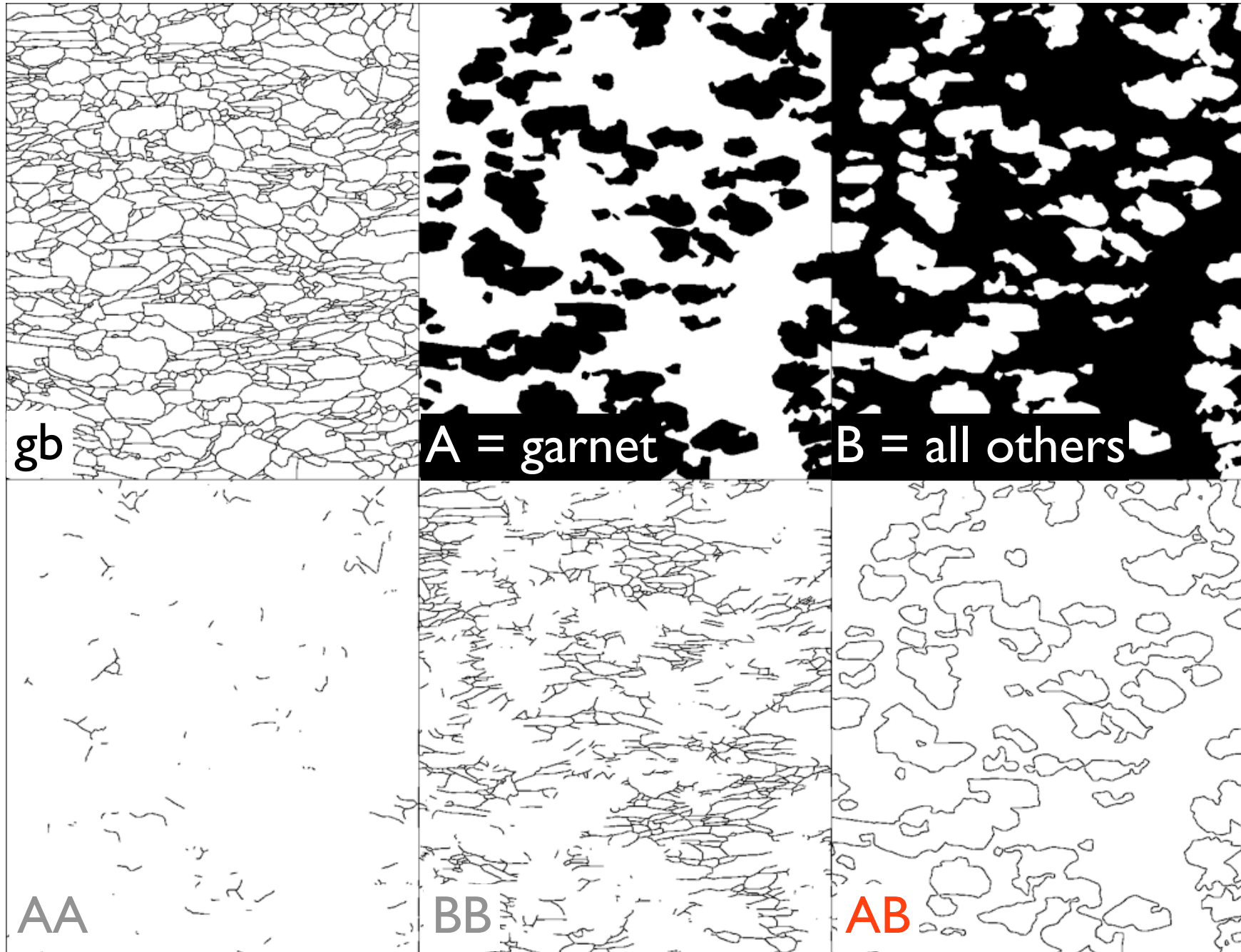


horizontal A gb

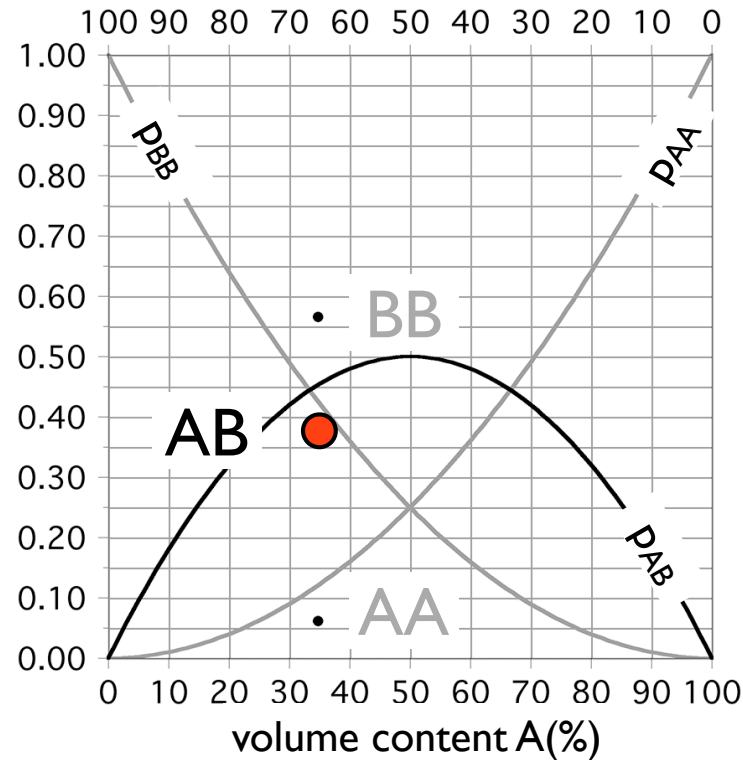
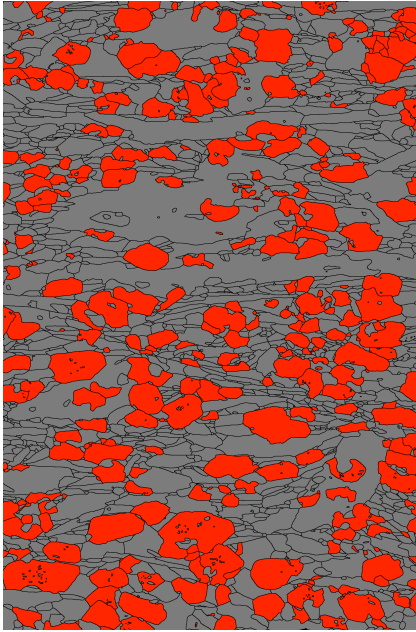
distribution of garnet in eclogite



garnet in eclogite

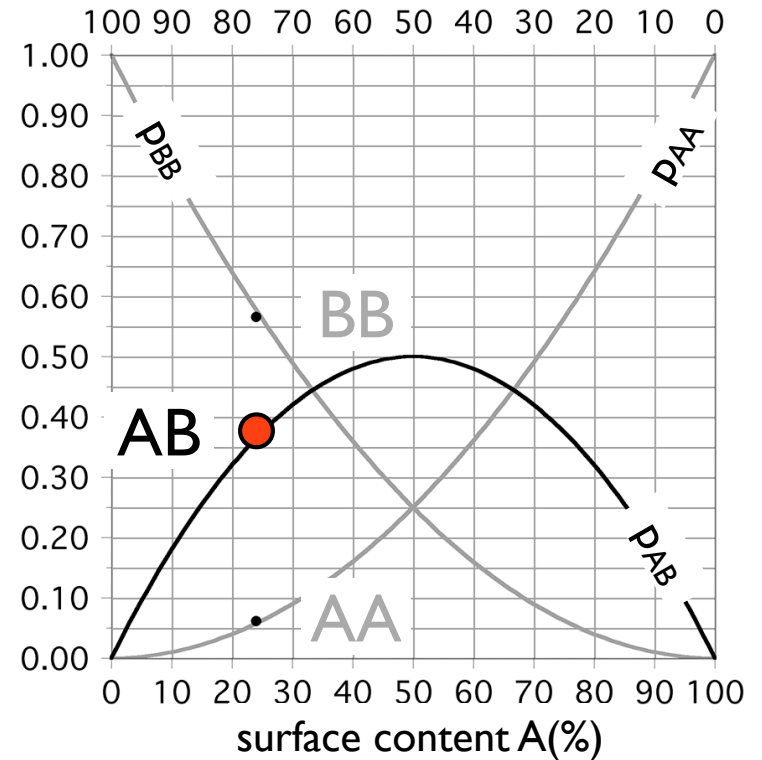


garnet in eclogite



in terms of
volume proportions

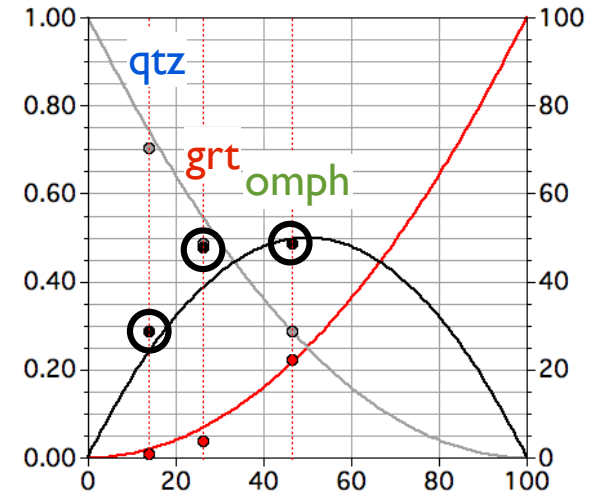
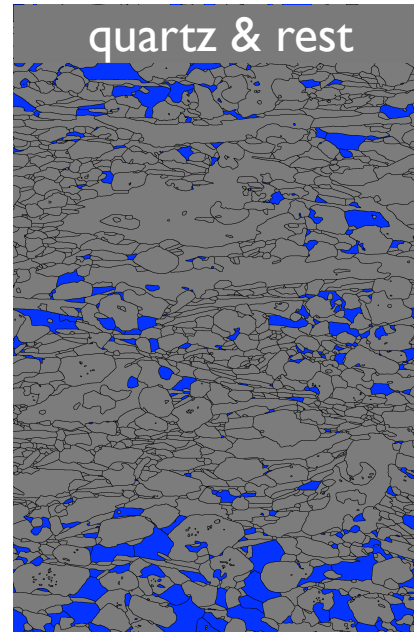
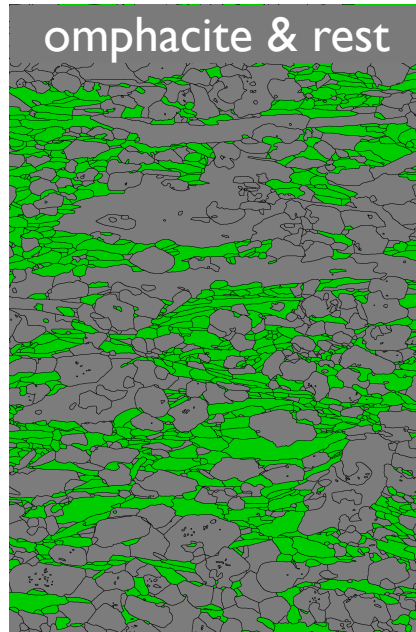
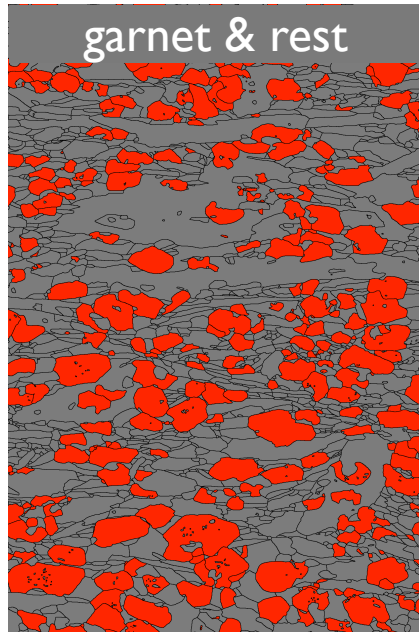
AB ~ random - clustered
AA anticlustered
BB clustered



in terms of
surface proportions

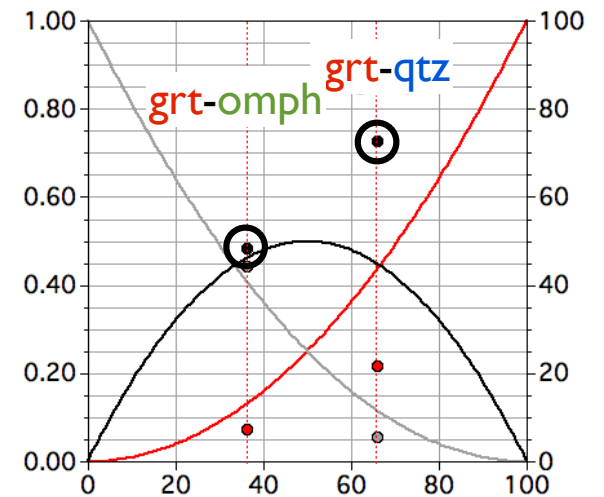
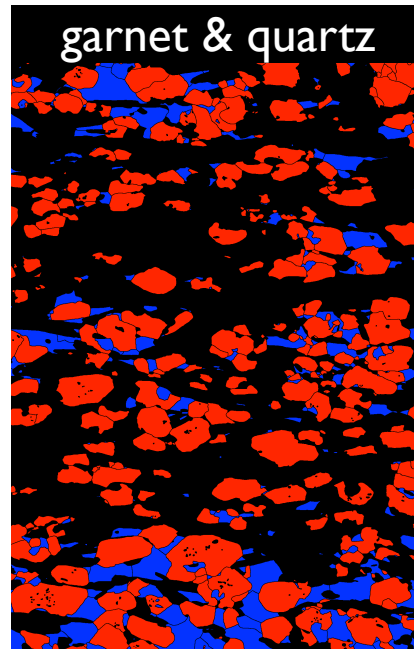
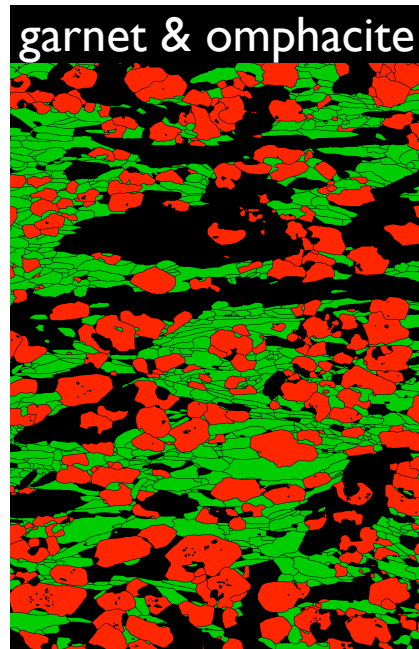
AB random
AA random
BB random

random - clustered - ordered ?



grt, omph, qtz = random
phase contacts on random curve

500 μm

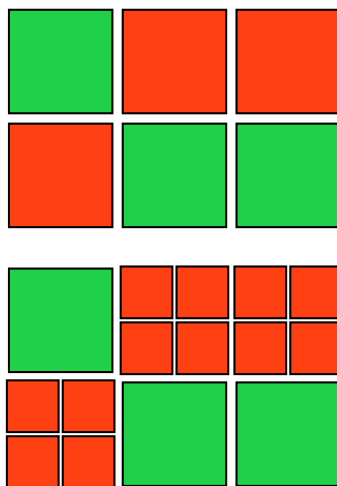
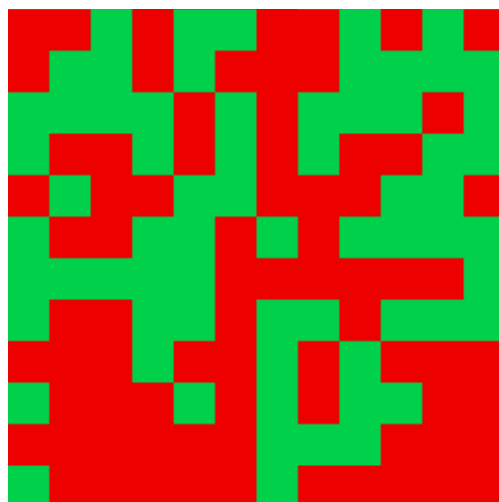


grt & omph = random
grt & qtz = ordered

2

influence of grain size

different grain sizes



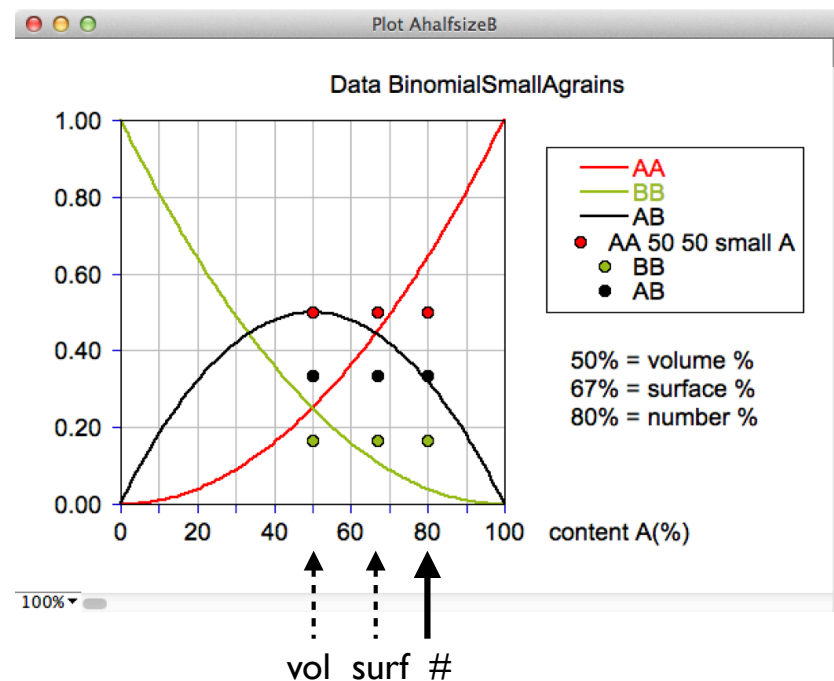
* each grain A divided in 4

volB = volB
 volA = volA
 volA* = volA
 nB = 50
 nA = nB
 nA* = 4 · nA
 surfB = surfB
 surfA = surfB
 surfA* = 2 · surfA

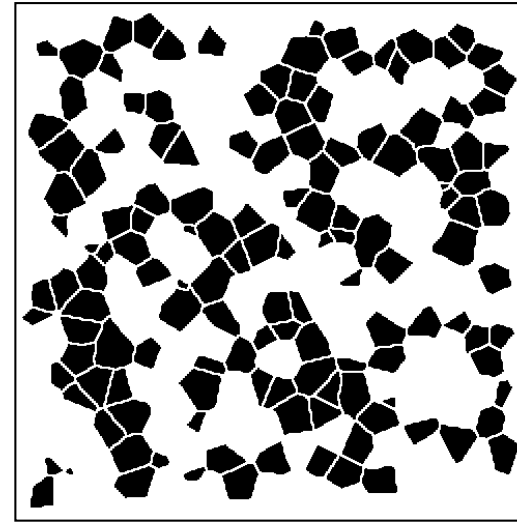
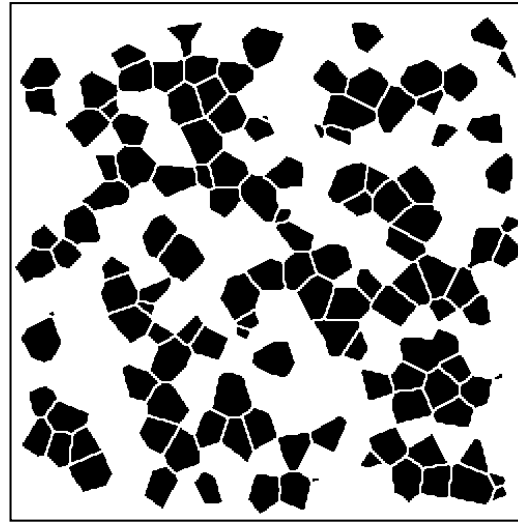
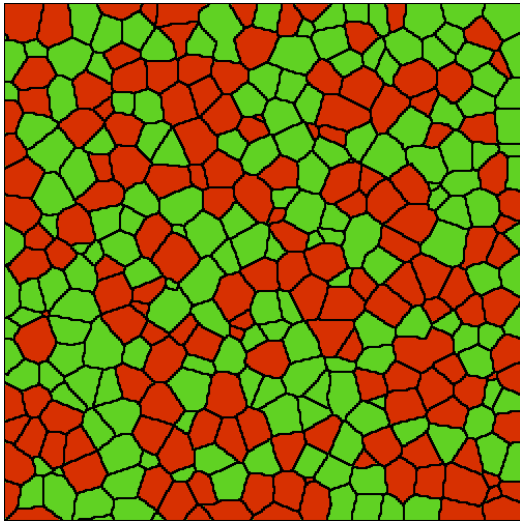
	A	50 A grains	A*	200 A grains
	AB	100	50	33.3
	BB	50	25	16.7
	AA	50	25	
	AA*	150		50.0

note: for A = 1/2 size of B
 number% best fit:
 phase = random
 BB = ordered
 AA* = clustered

	A	A = B	A smaller	%
0	vol A	50.000	50.000	
1	vol B	50.000	50.000	
2	total volume	100.00	100.00	
3				
4	no A grains	50.000	200.00	80.000
5	no B grains	50.000	50.000	20.000
6	total no grains	100.00	250.00	100.00
7				
8	surface A	50.000	100.00	66.670
9	surface B	50.000	50.000	33.330
10	total surface	100.00	125.00	100.00
11				
12	no contact AA	50.000	150.00	50.000
13	no contact BB	50.000	50.000	16.670
14	no phase AB	100.00	100.00	33.330
15				



2D grain size of Voronois



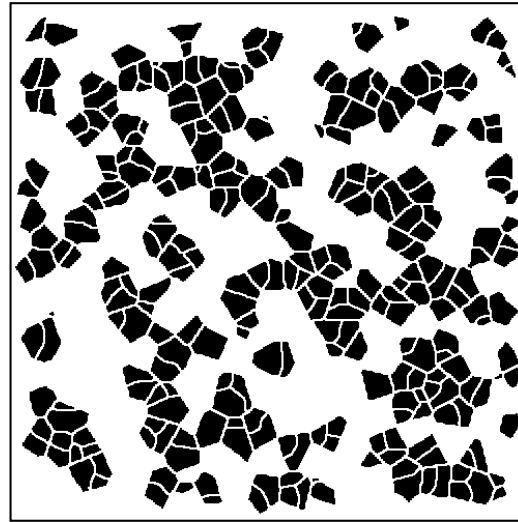
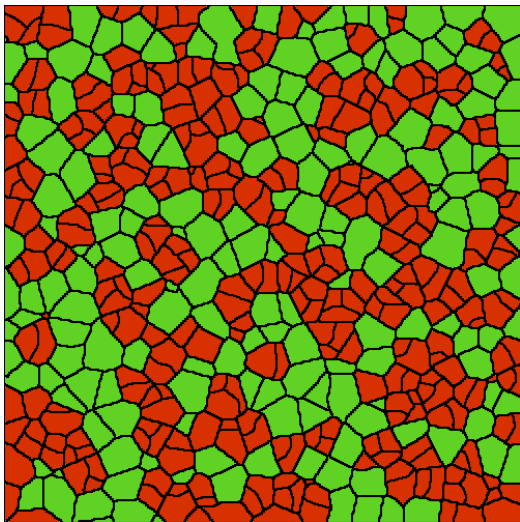
size $A \approx B$
no. $A \approx B$
no. $A_{small} \approx 2x$

A (129 grains)

mean=30.3, RMS=31.4

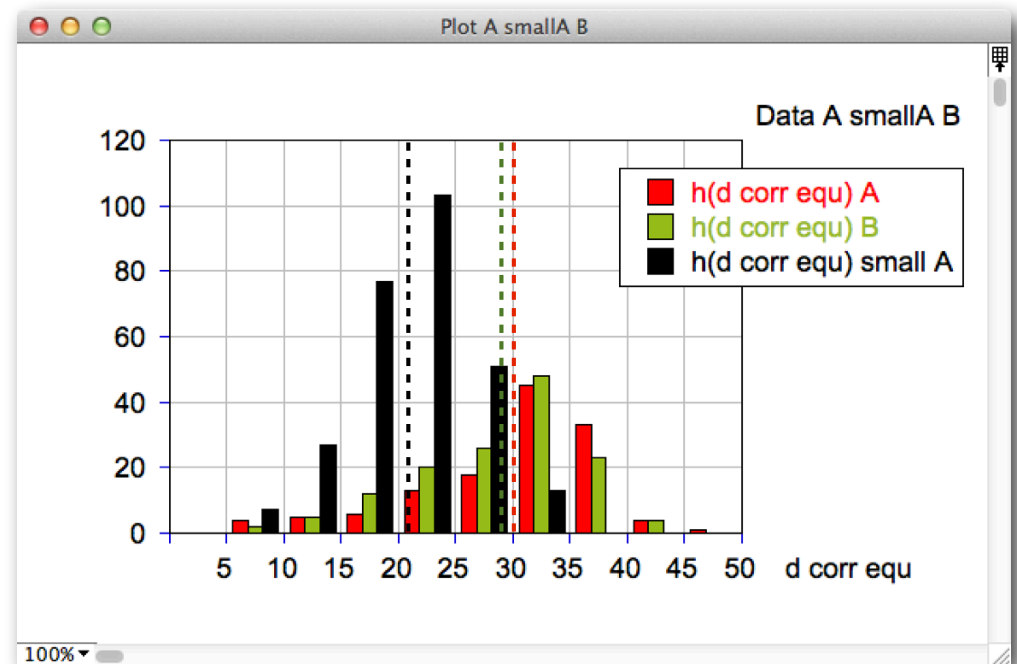
B (140 grains)

mean=28.9, RMS=29.9

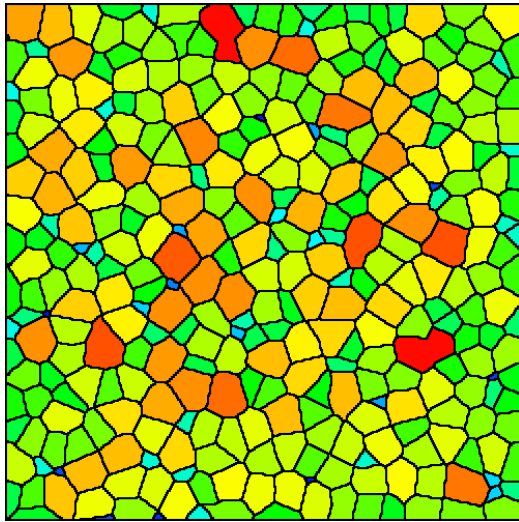


Asmall (278 grains)

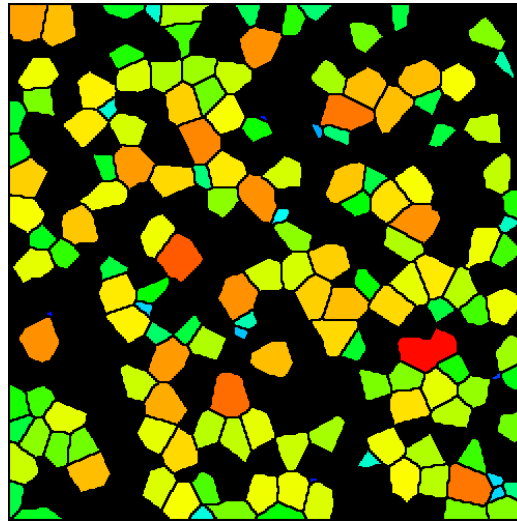
mean=21.0, RMS=21.7



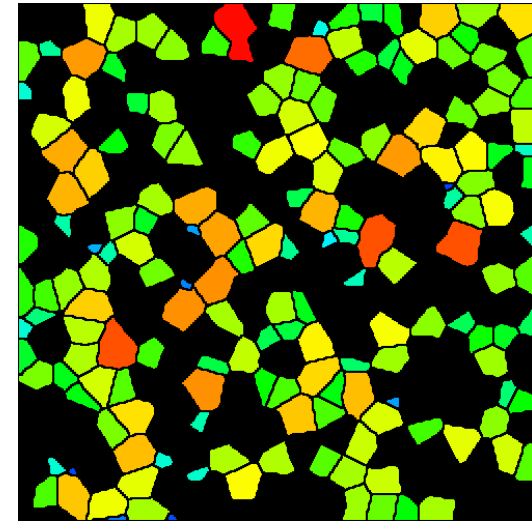
grain size $A <$ grain size B



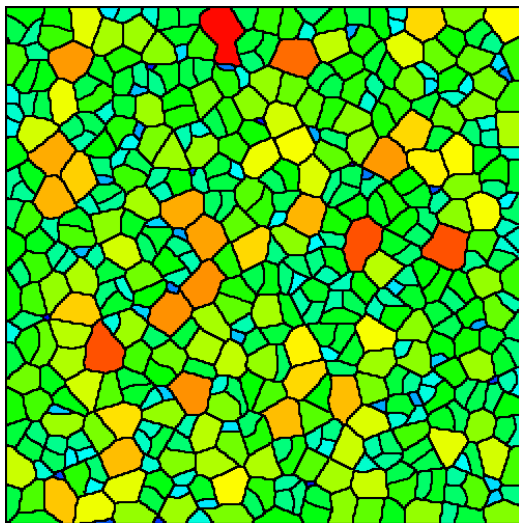
AB all



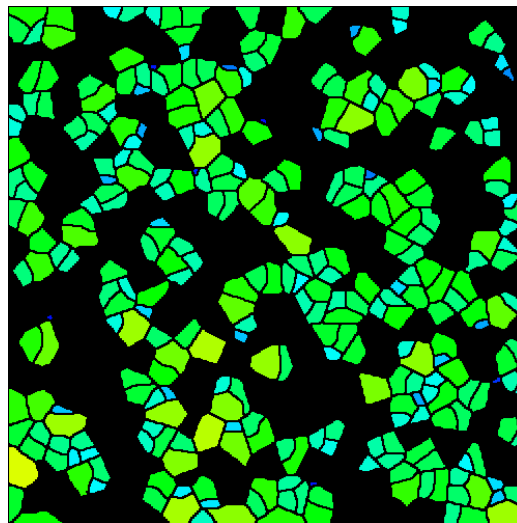
A



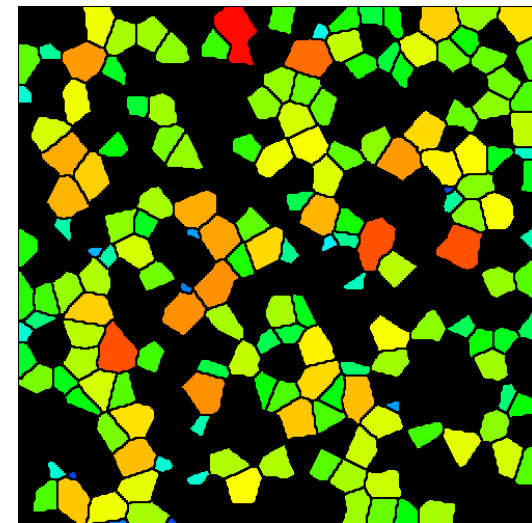
B



AsmallB all



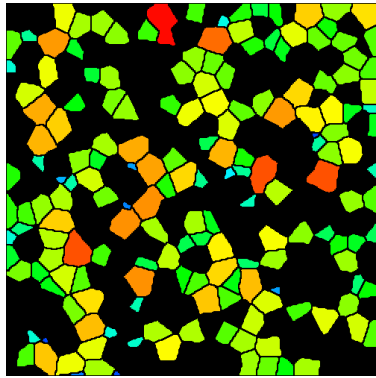
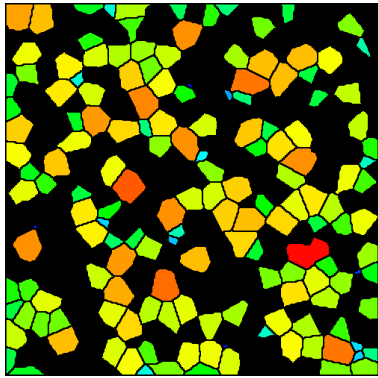
Asmall



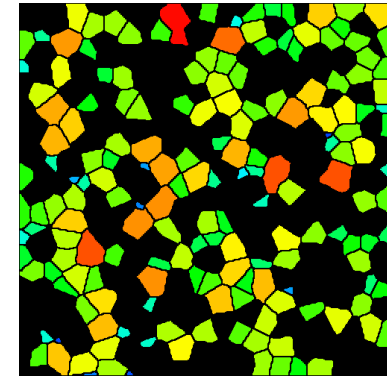
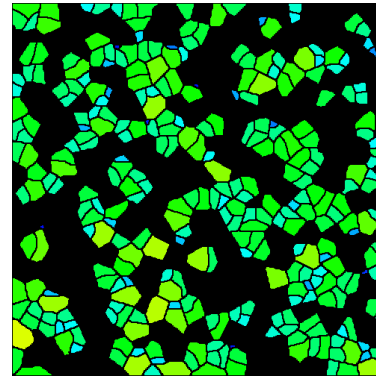
B unchanged

small phase A obtained by subdividing grains, B phase remains unchanged

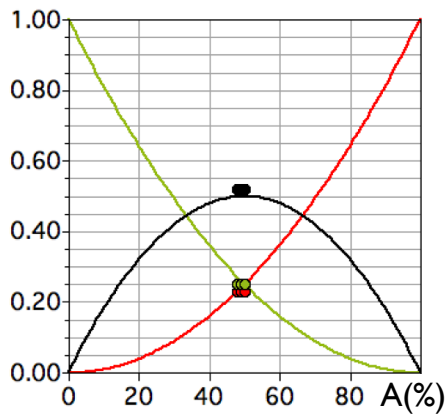
volume & - surface % - number %



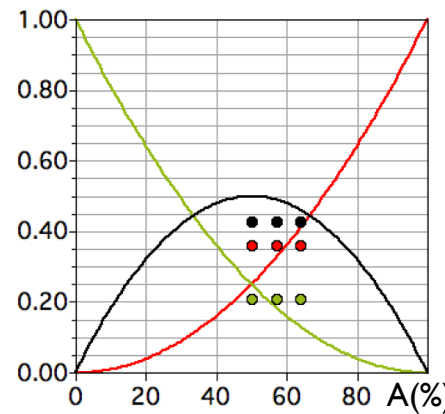
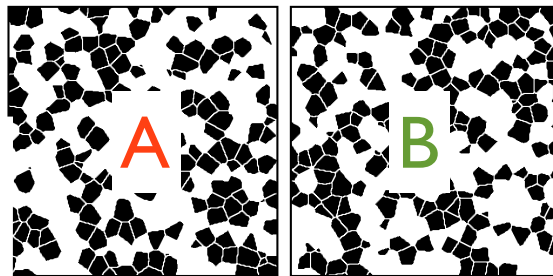
grainsize A = grainsize B



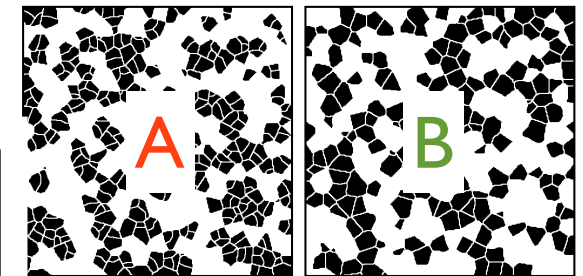
grainsize A < grainsize B
phase A 'recrystallized' in situ



vol% A = 50.0
surf% A = 48.9
no.% A = 47.8
no A:B = 160:175



vol% A = 50.0
surf% A = 57.3
no.% A = 64.2
no A:B = 314:175



3

variations

Menegon et al. 2013

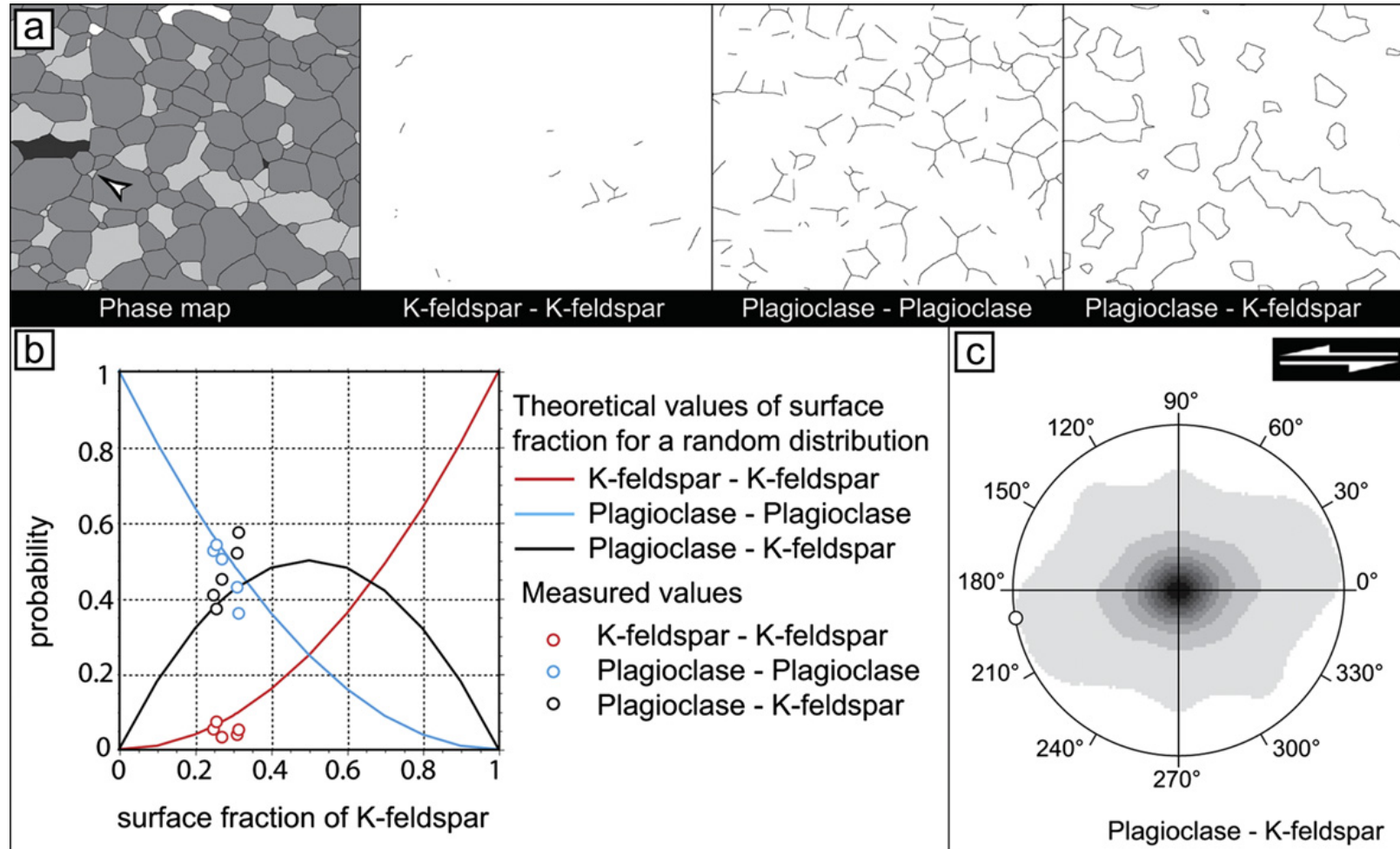


Fig. 7. (a) Example of phase map used in the analysis of spatial distribution of feldspars in the recrystallized matrix, and bitmaps of the grain- and phase boundaries derived from the map. The arrowhead indicates a K-feldspar grain at a quadruple junction between plagioclase grains. (b) The fraction of grain- and phase boundaries measured in five different phase maps are plotted as circles to evaluate the deviations from randomness. Curves show the theoretical fractions of grain- and phase boundaries as a function of the surface fraction of K-feldspar expected for a random distribution in a two-phase mixture (c) Autocorrelation function (ACF) calculated for the plagioclase – K-feldspar phase boundaries. The ACF is shaded at 10% multiples of the ACF max (located in the centre). For the sake of clarity, the ACF is scaled such that the 10% contour touches the superposed reference circle. The black circle indicates the highest correlation length at 12° (measured anticlockwise) from the trace of the mylonitic foliation (E–W).

The microstructure of recrystallized feldspars is characterized by the predominance of phase boundaries over grain boundaries

Kilian et al. 2011

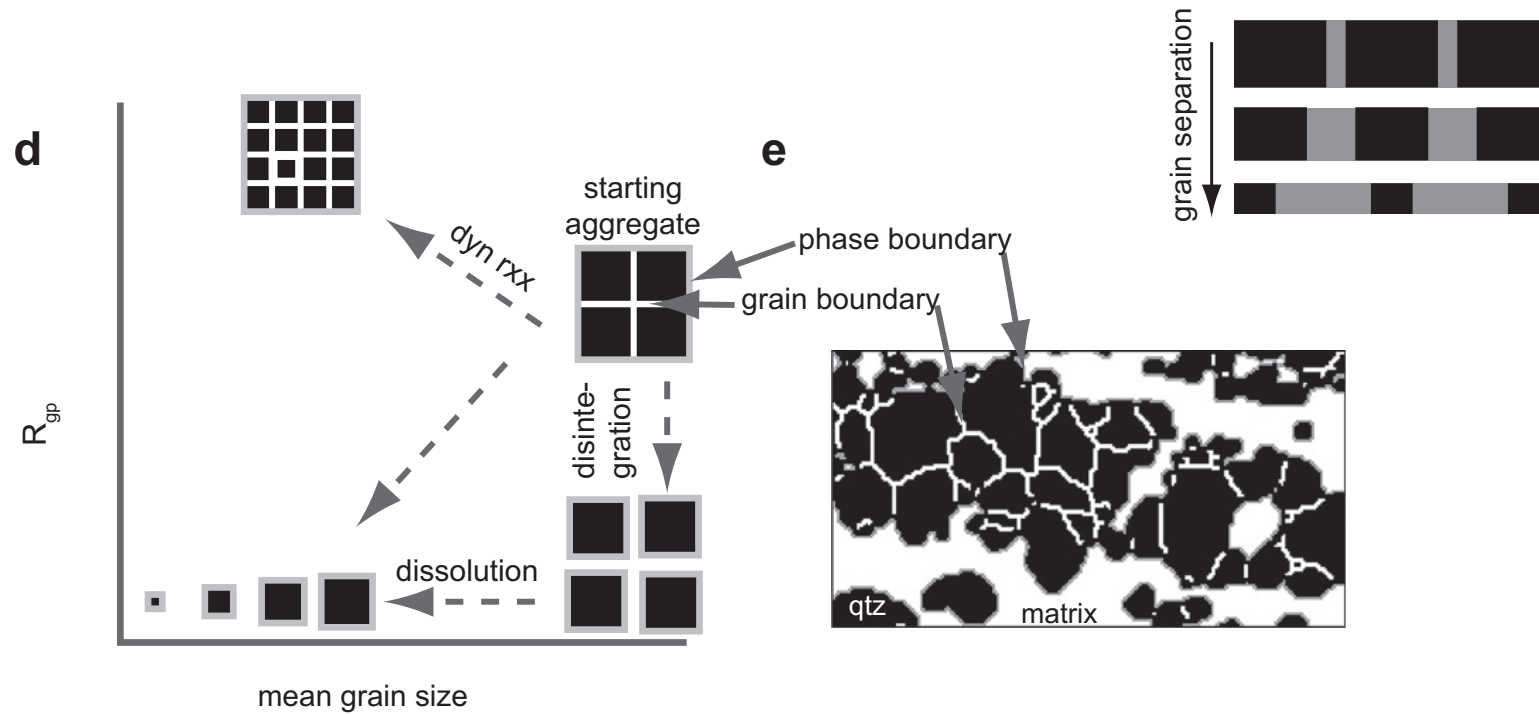


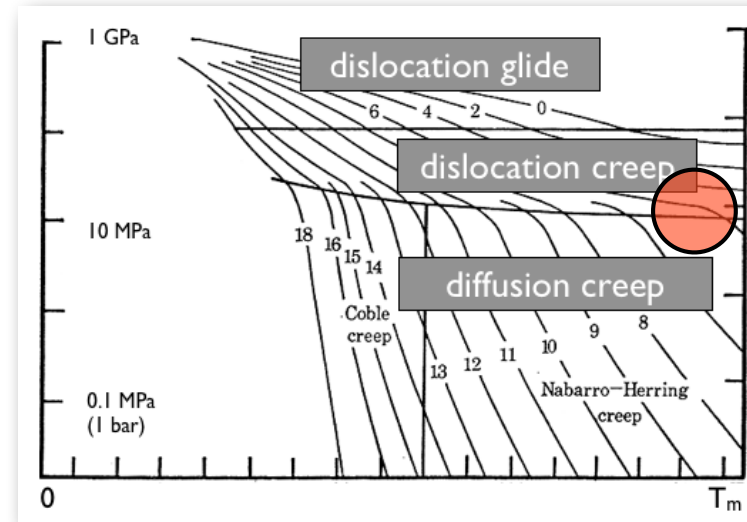
Fig. 15. Phase distribution and grain size evolution. (a) Phase map of quartz (black), plagioclase (dark gray), K-feldspar (light grey) and mica (white). The map shows a quartz aggregate in the transition between the mylonite and the ultramylonite with different stages of disintegration. (b) R_{gp} (ratio of grain boundary/phase boundary area) versus the mean grain size in 6 different, laterally disintegrating aggregates. Symbols correspond to individual, homogenous aggregates. (c) Relationship between the mean quartz grain size in a layer and the relative quartz volume fraction. The relative quartz volume fraction in a layer decreases with increasing disintegration, as grains are separated by K-feldspar. In layers with a higher grain separation the mean grain size is smaller. (d) Schematic model of R_{gp} – grain size evolution. Black squares are separated from another by grain boundaries (white) and from the matrix by phase boundaries (dark grey) (e) Example of an analyzed area. The evaluation of R_{gp} and the grain size would correspond to a single point in the diagram b).

practical application

dislocation creep vs. diffusion creep

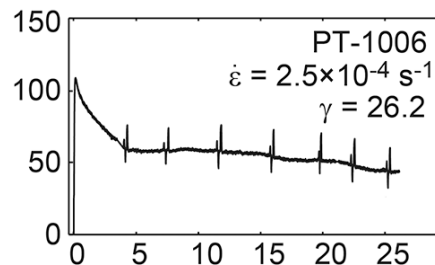


Gas medium High pressure Torsion apparatus (UMN)



Miki Tasaka Mark Zimmerman David Kohlstedt

70% iron-rich olivine
30% orthopyroxene
hotpressed @1200°C
d ~15 μm



$p_c = 300 \text{ MPa}$
 $T = 1200^\circ\text{C}$

in olivine:
MeO dissolves at maximum σ_1 .
Reaction ol \rightarrow opx

in orthopyroxene:
MeO diffuses to tension σ_3 .
Reaction opx \rightarrow ol

olivine
 opx
 MeO
 diffusion pass

r1: Me_2SiO_4 (ol) - MeO \rightarrow MeSiO_3 (opx)
 r2: MeSiO_3 (opx) + MeO \rightarrow Me_2SiO_4 (ol)

motivation

$$\dot{\epsilon} = A \cdot \Delta\sigma^n \cdot \exp(-Q/RT)$$

increasing strain

$$\dot{\epsilon} = A \cdot \Delta\sigma^n \cdot d^m \cdot \exp(-Q/RT)$$

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Research Article

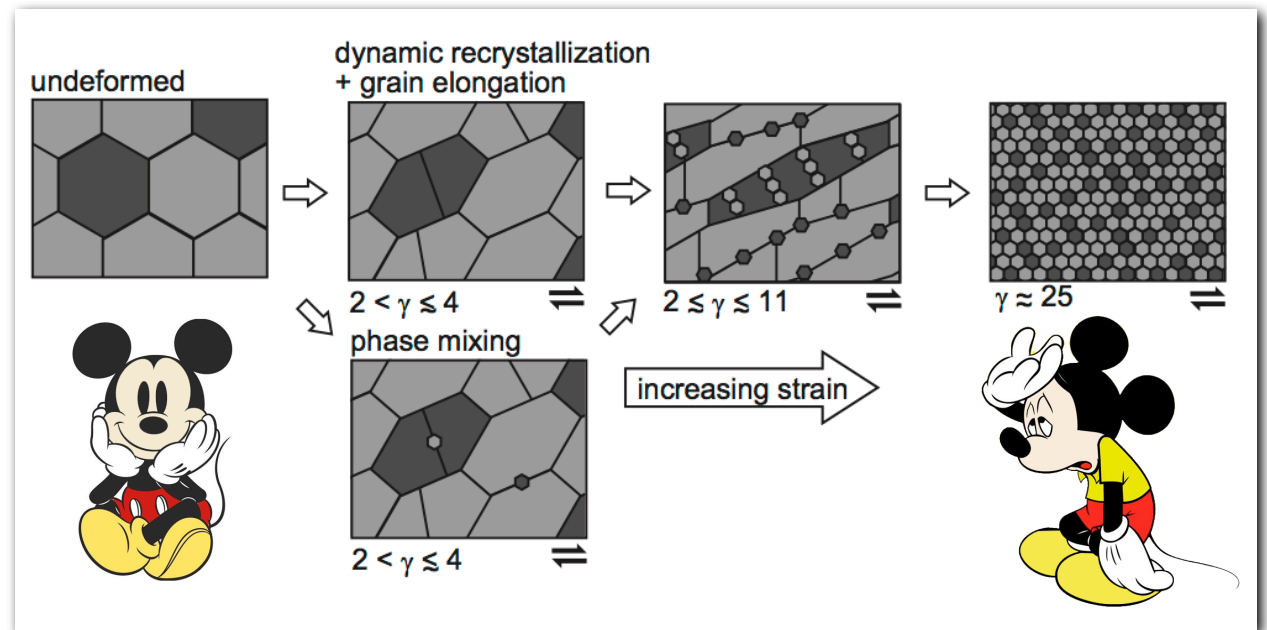
Rheological weakening of olivine + orthopyroxene aggregates due to phase mixing, Part 1: Mechanical behavior

Miki Tasaka , Mark E. Zimmerman, David L. Kohlstedt

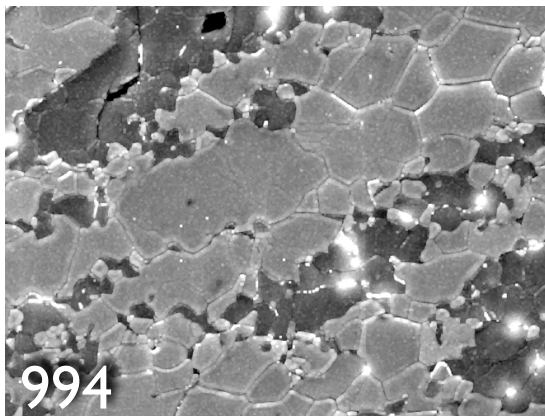
Accepted manuscript online: 8 September 2017 [Full publication history](#)

DOI: 10.1002/2017JB014333 [View/save citation](#)

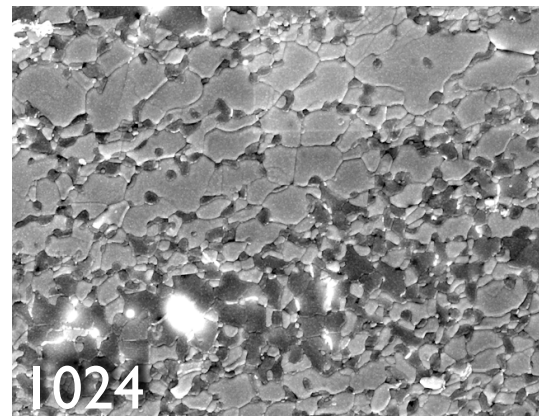
Tasaka et al. (JGR, 2017)



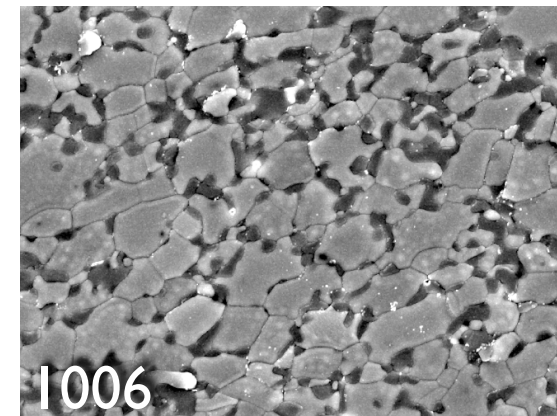
dislocation creep ?



phase mixing ?



diffusion creep ?



10 μ m

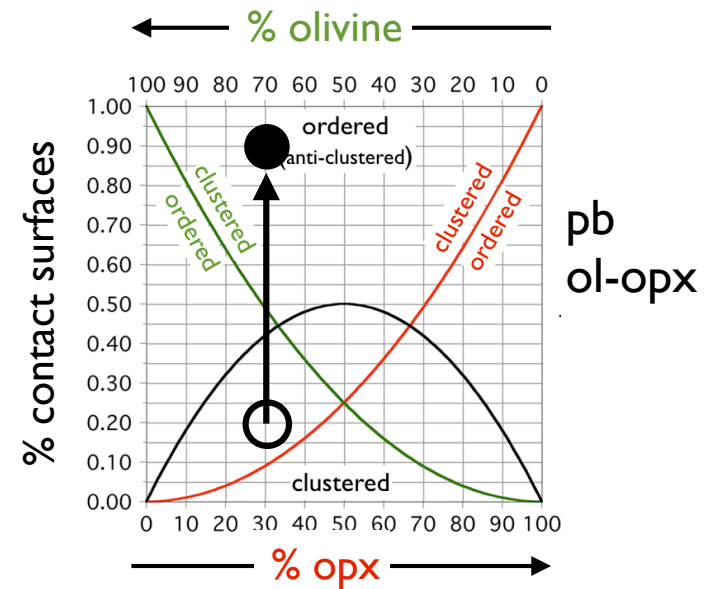
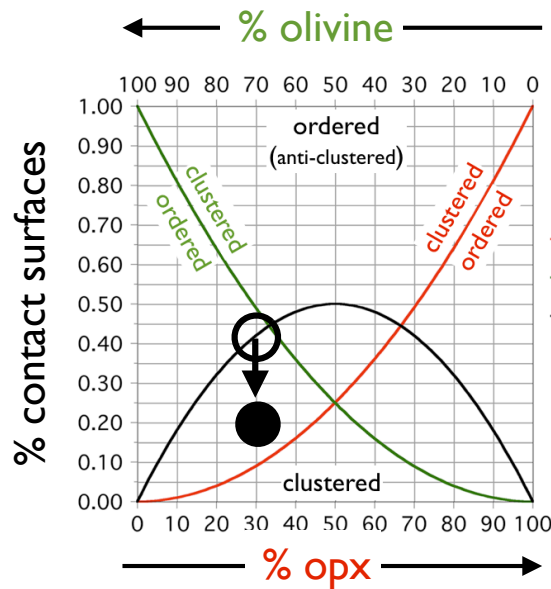
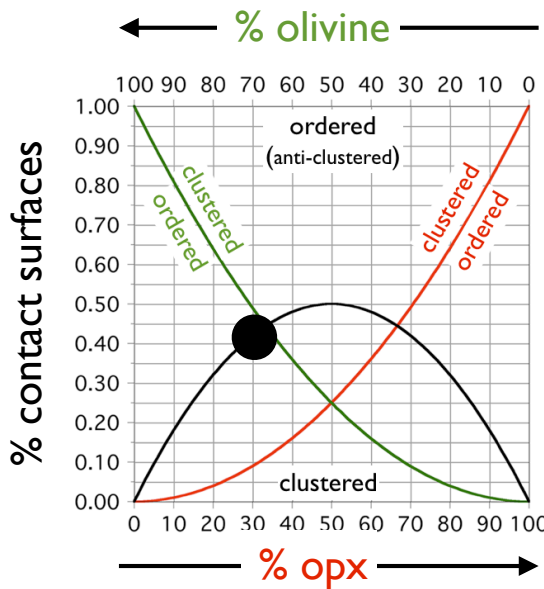
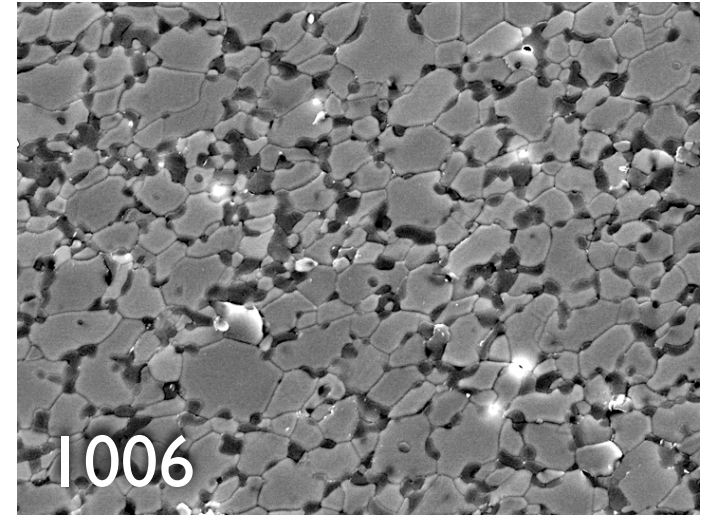
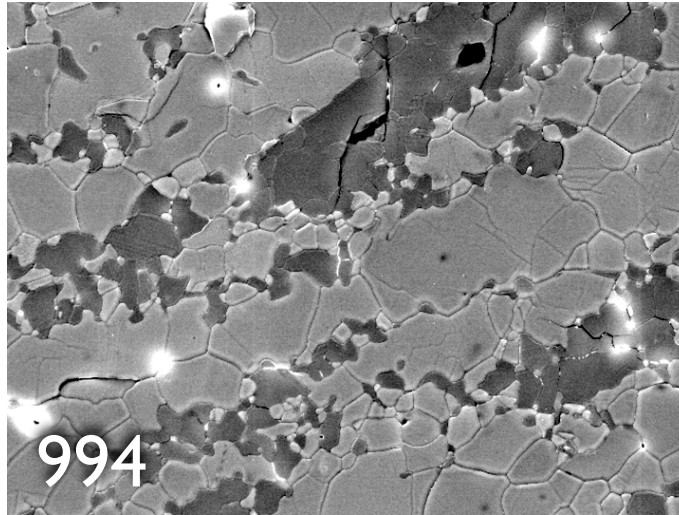
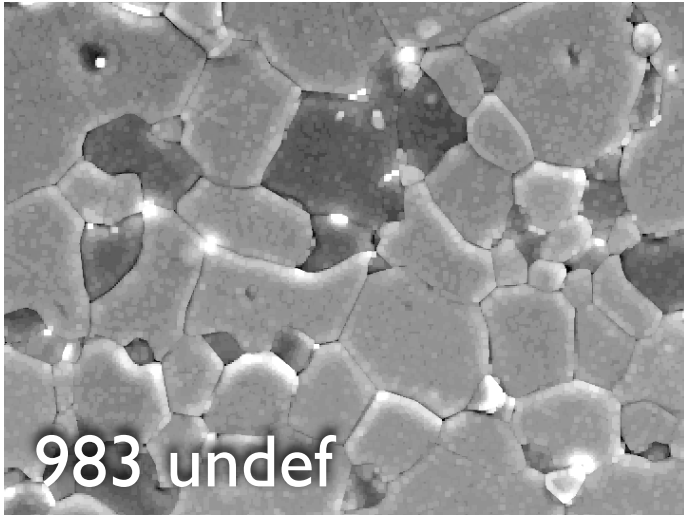
which spatial distributions do we expect ?

starting material

dislocation creep

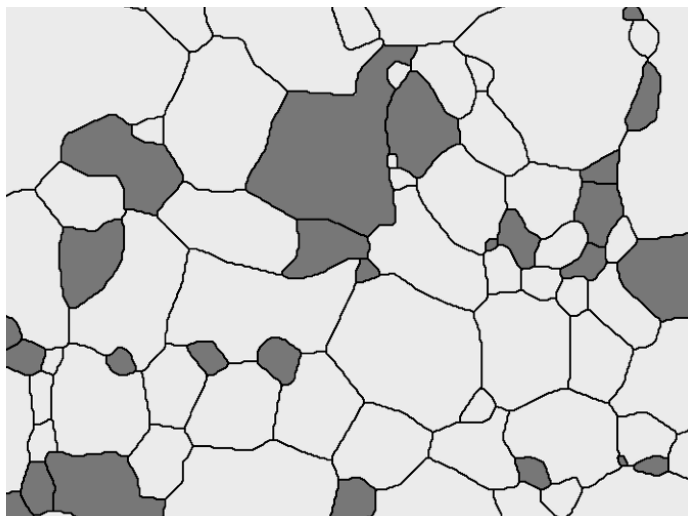


diffusion creep

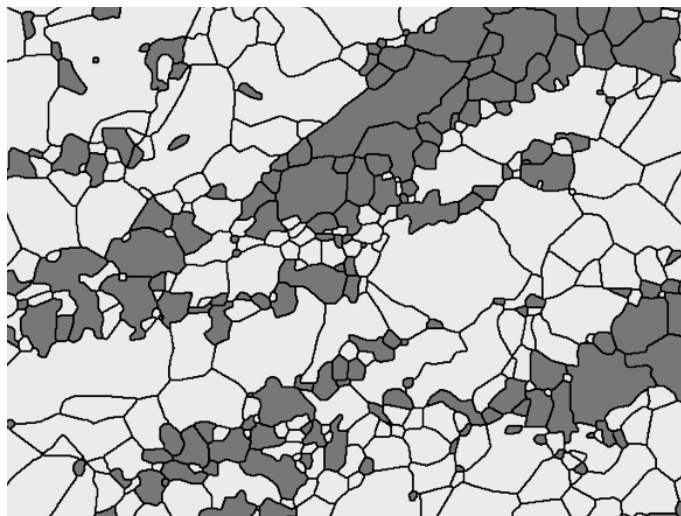


and which spatial distributions do we get ?

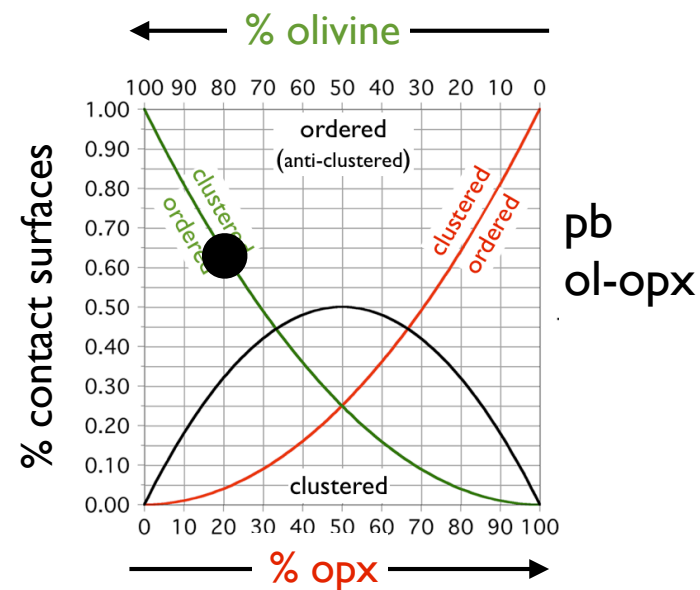
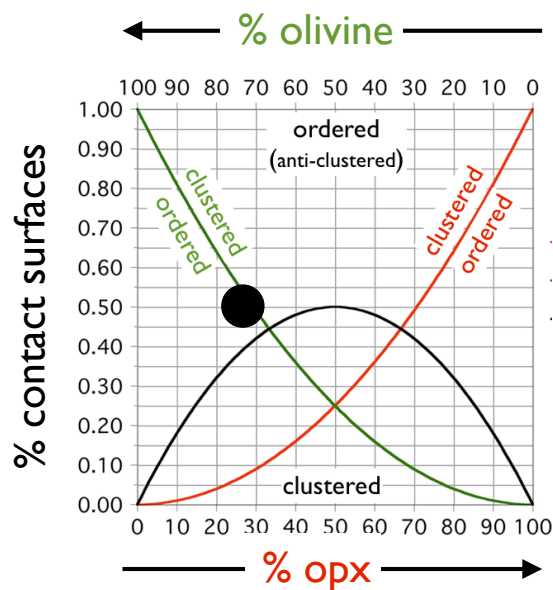
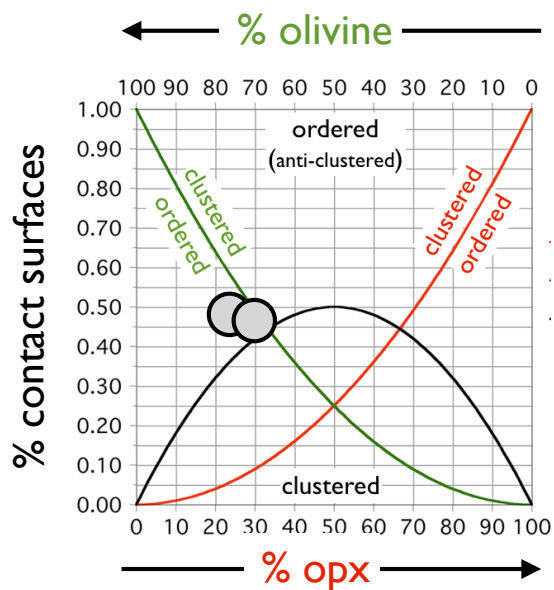
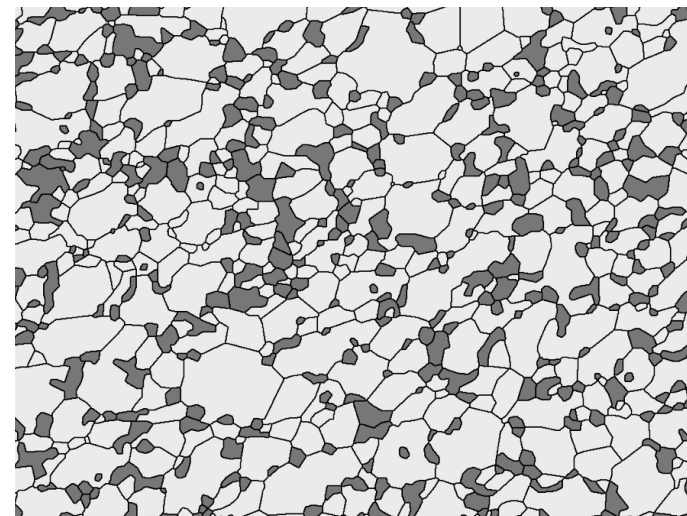
983 starting material



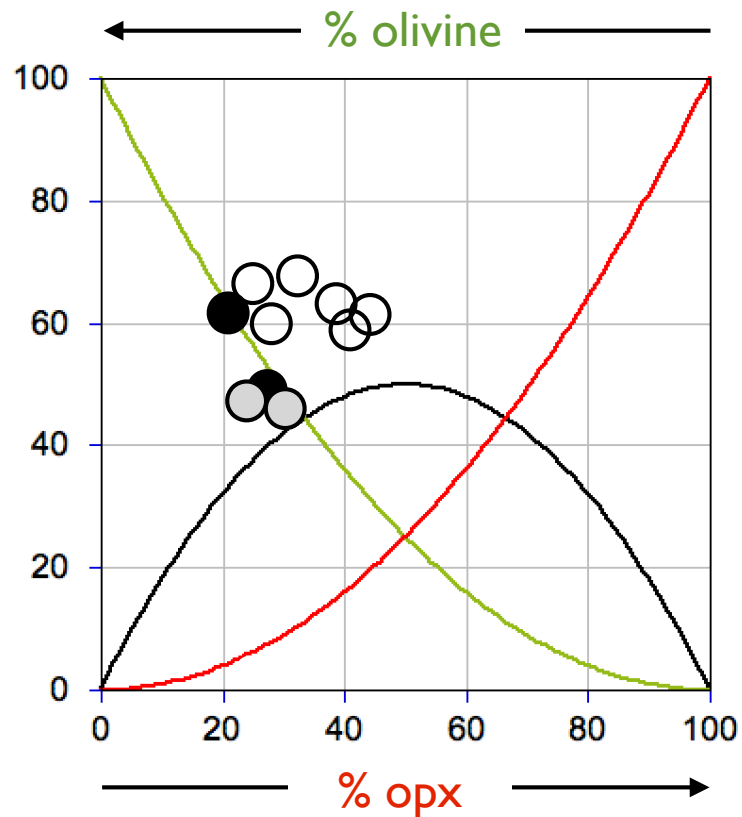
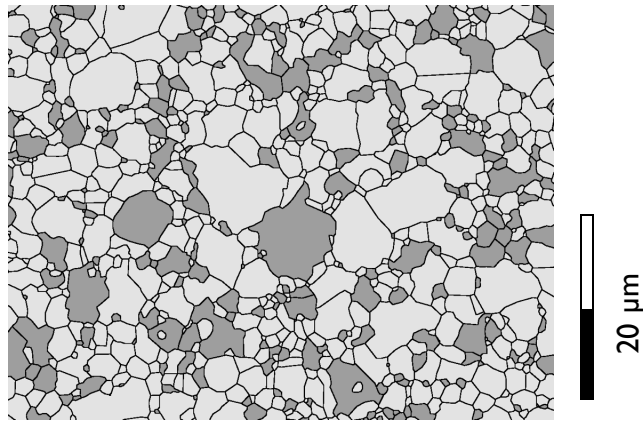
994 dislocation creep



1006 diffusion creep



even the starting material is ordered !!



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Research Article

Rheological weakening of olivine + orthopyroxene aggregates due to phase mixing, Part2: Microstructural development

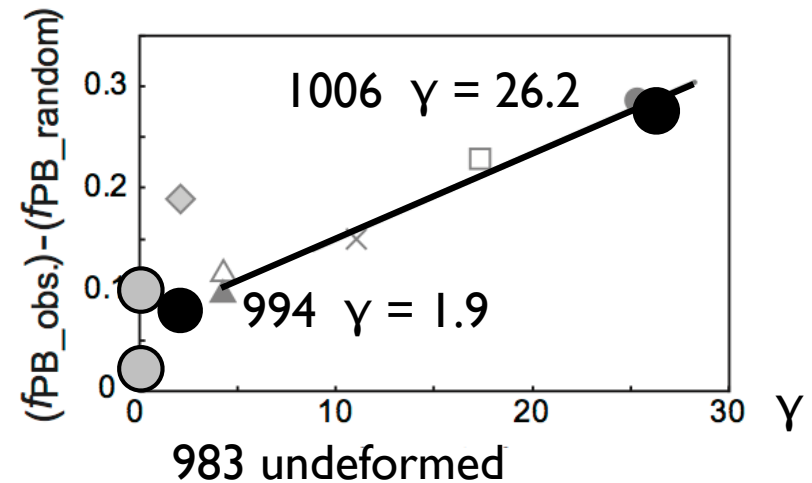
Miki Tasaka , Mark E. Zimmerman, David L. Kohlstedt, Holger Stünitz, Renée Heilbronner

Accepted manuscript online: 8 September 2017 [Full publication history](#)

DOI: 10.1002/2017JB014311 [View/save citation](#)

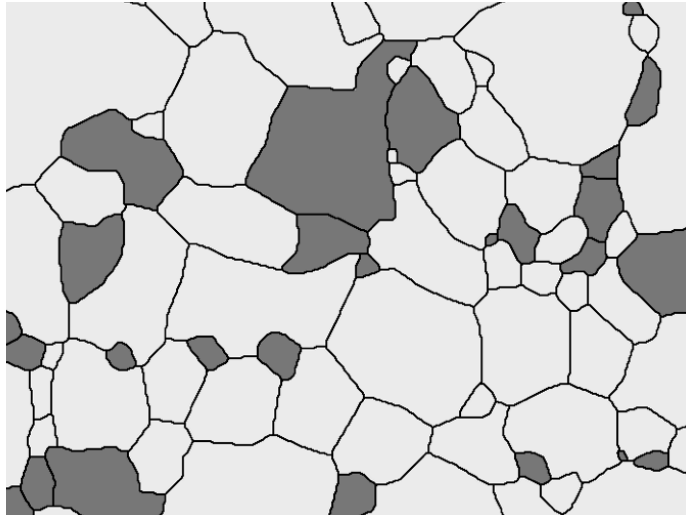
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Tasaka et al. (JGR, 2017)



which spatial distributions do we really get ?

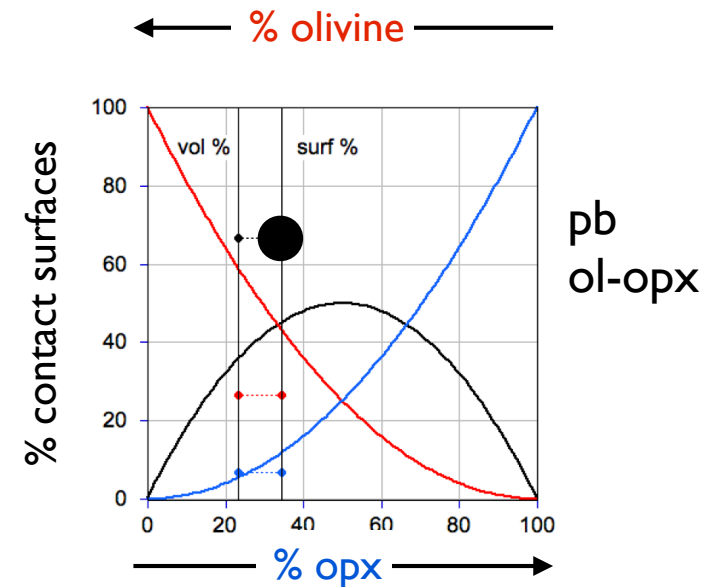
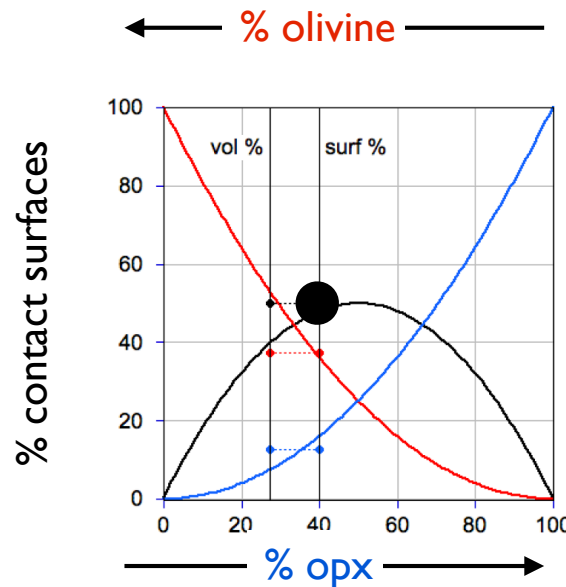
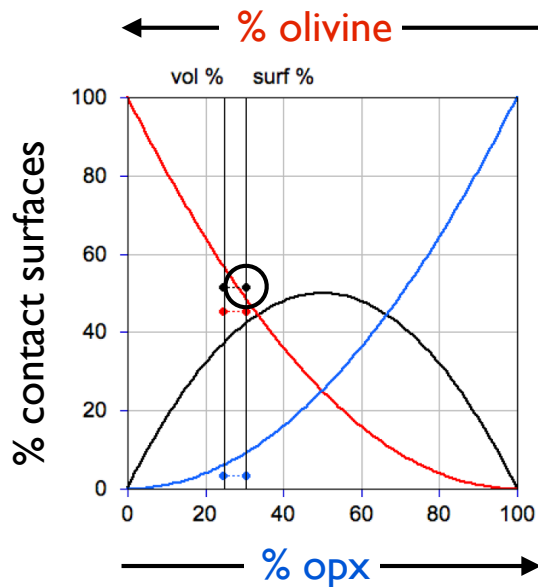
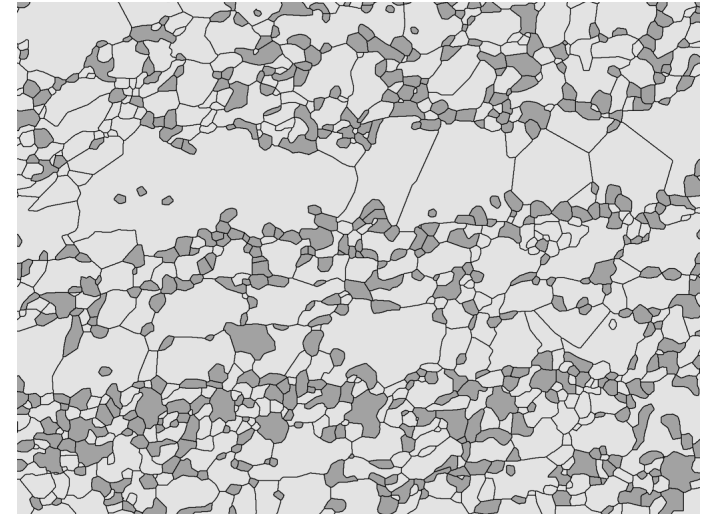
983 starting material



994 dislocation creep



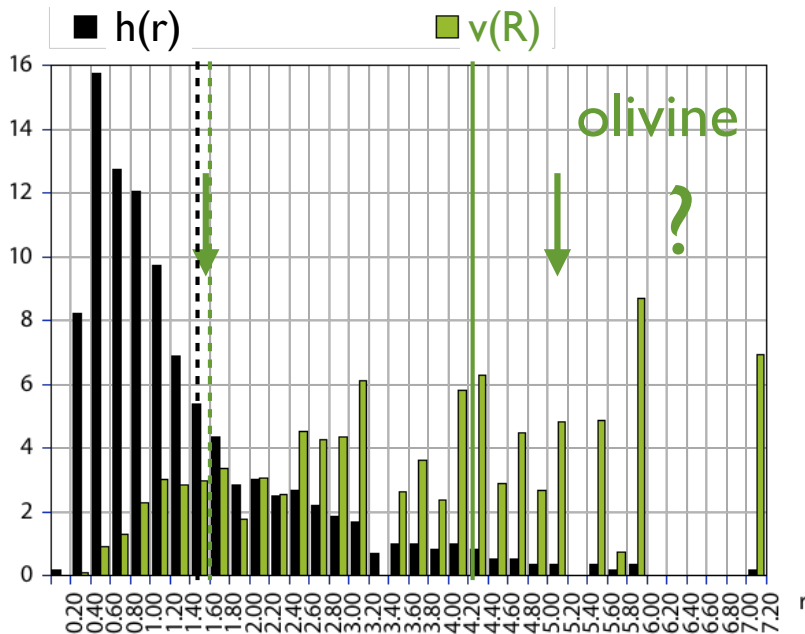
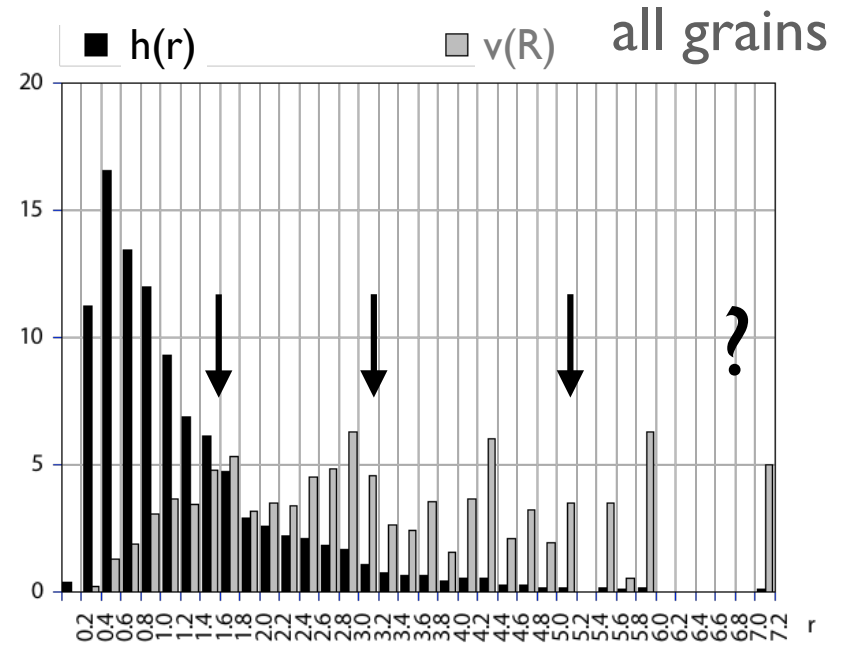
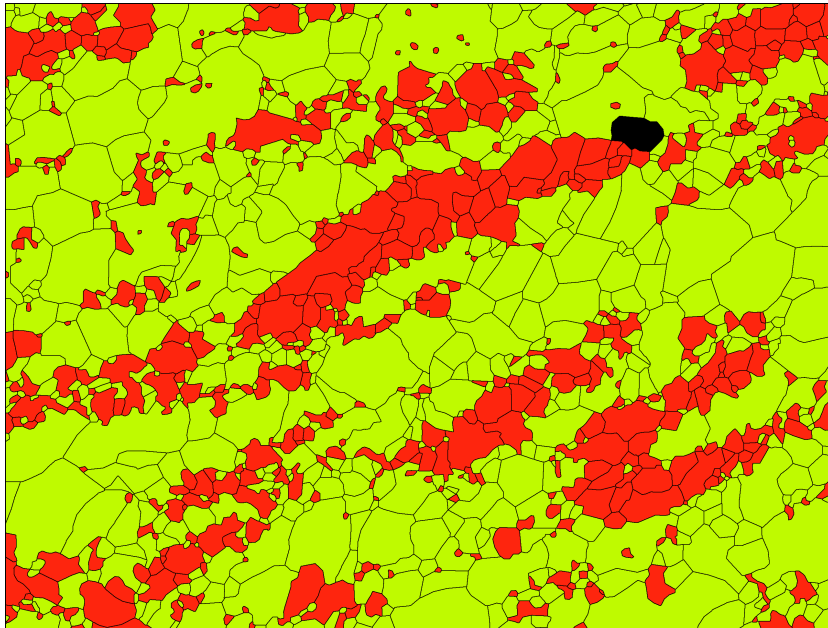
1024 diffusion creep



pb
ol-opx

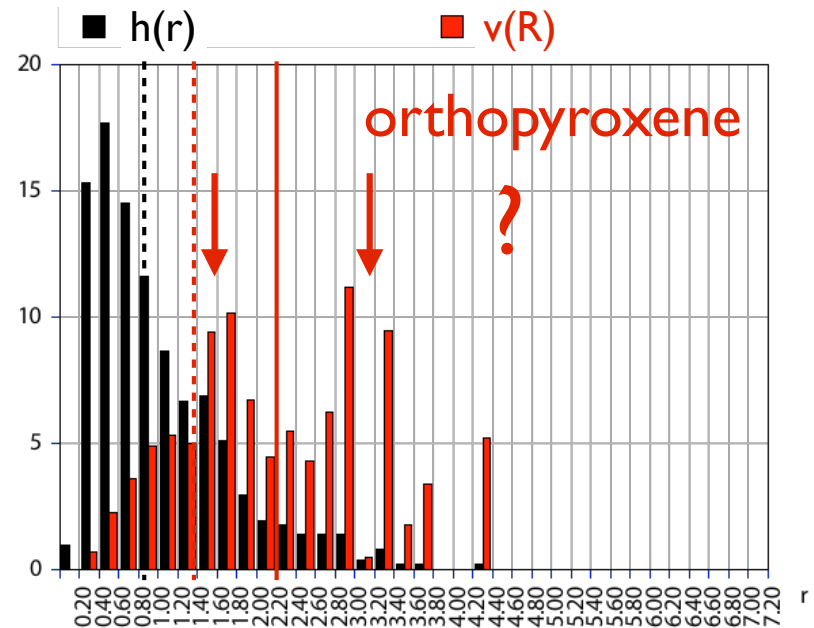
plotted against surface fraction !

also, grain size distributions ...



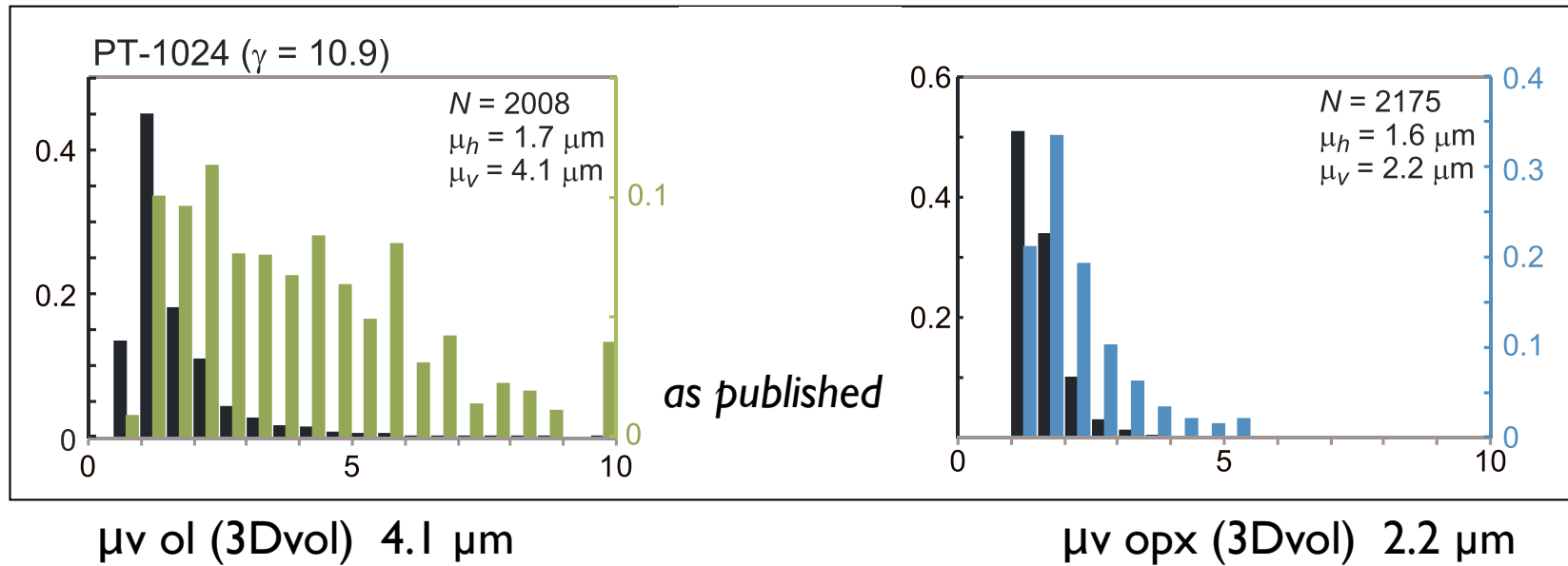
Miki's

- $\mu(d)$
- $\mu(D)$
- $\mu(V)$

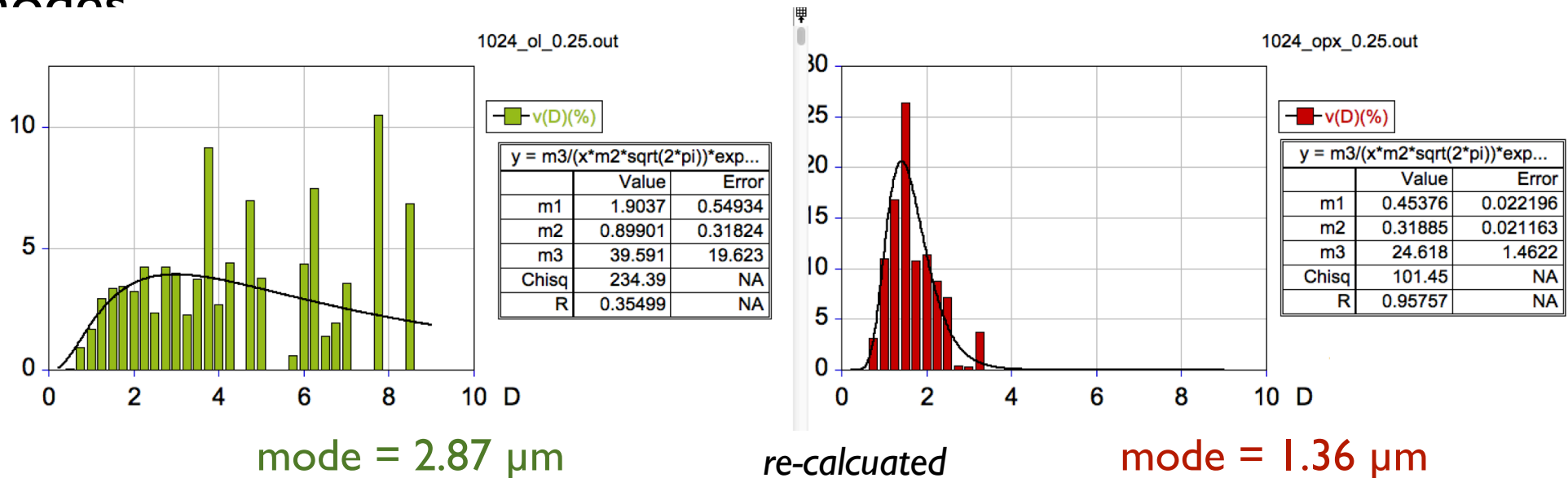


means... and modes

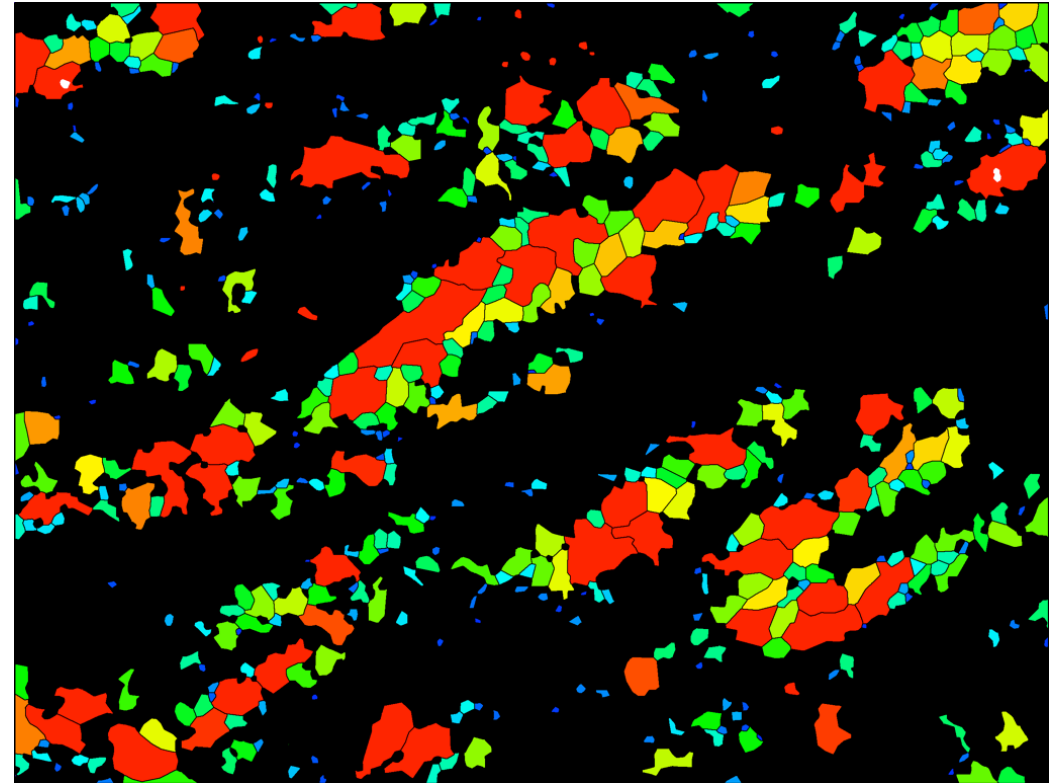
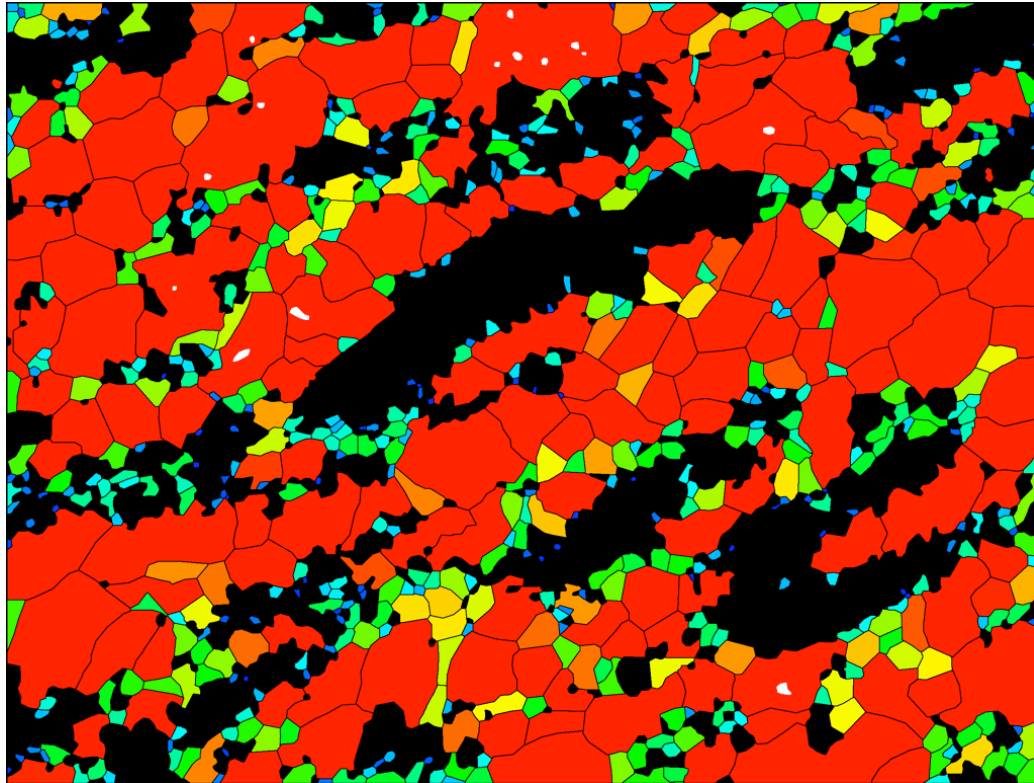
means



modes



grain size of ol and opx



olivine

ortho-pyroxene

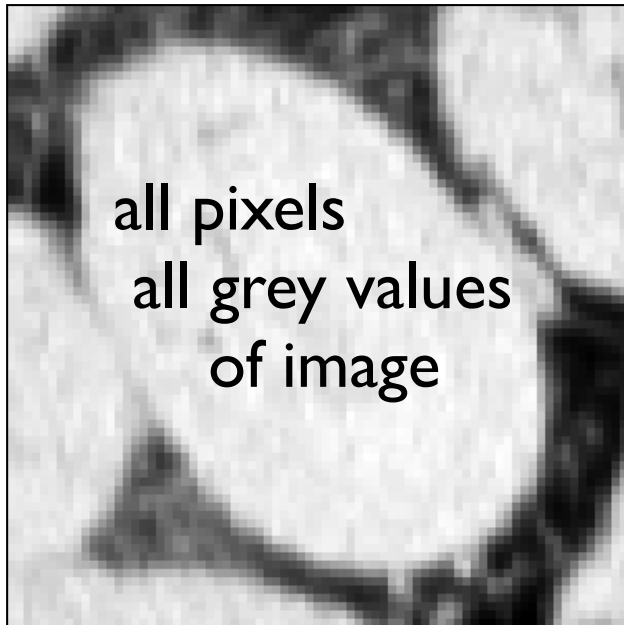


3

image statistics

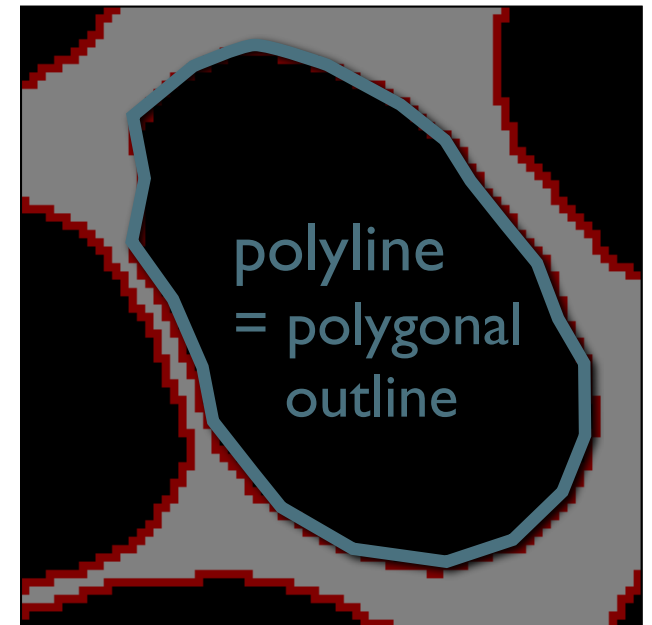
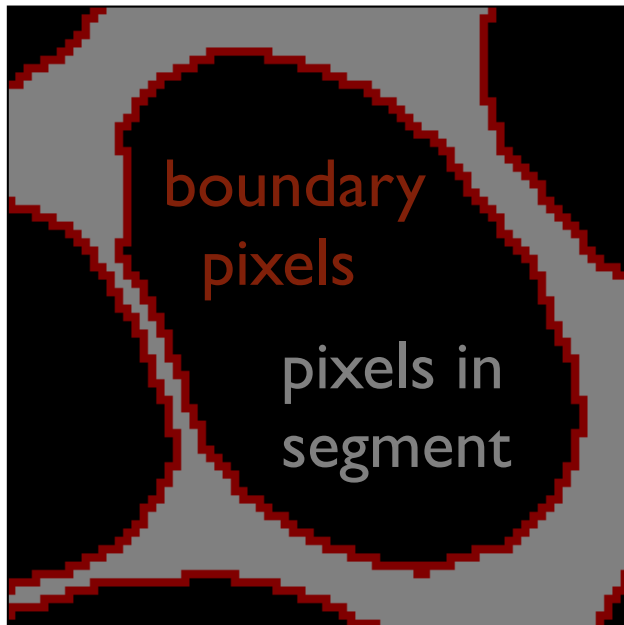
analysis of image 'as-is'

image



(no segmentation)

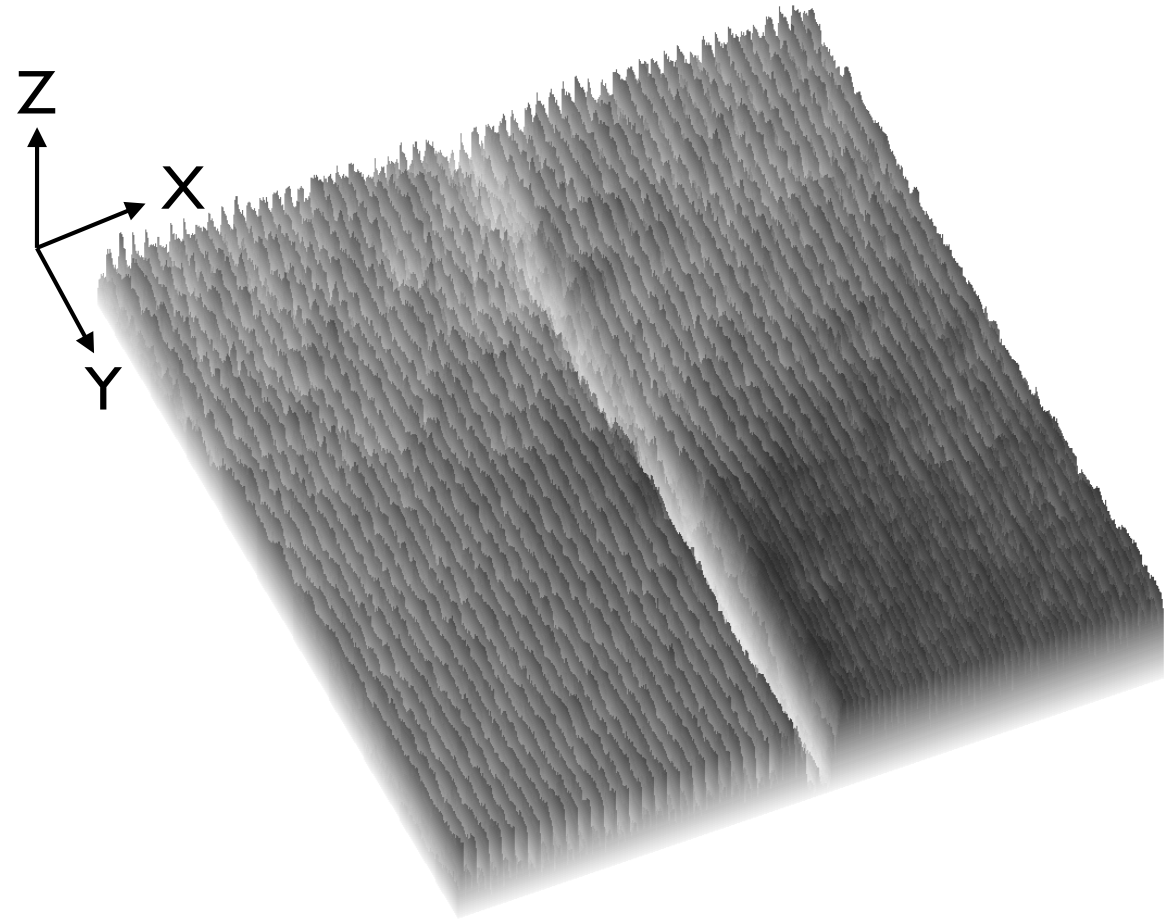
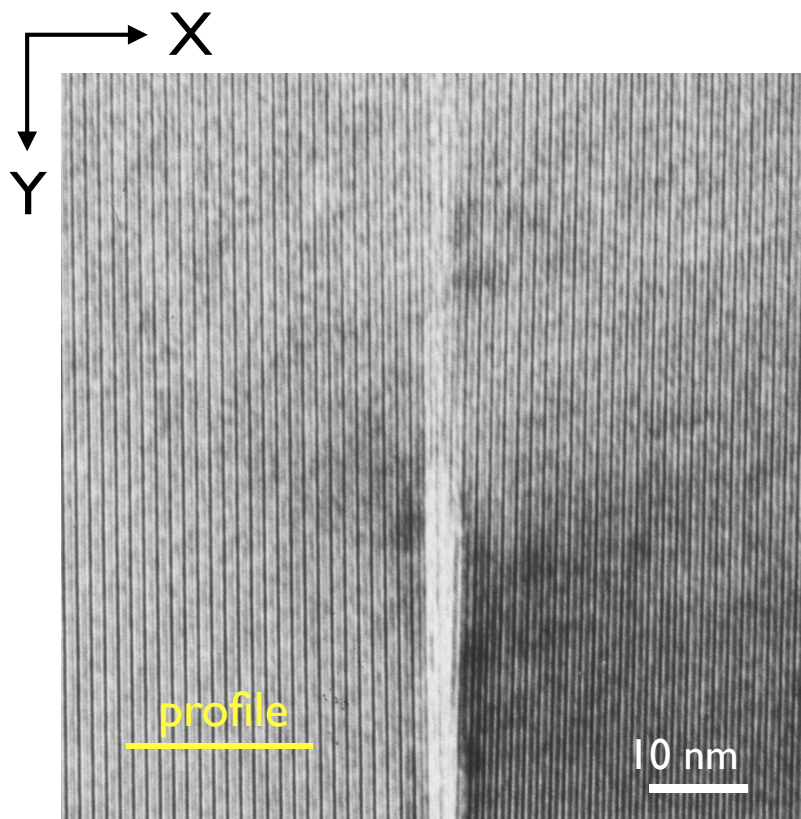
bitmap



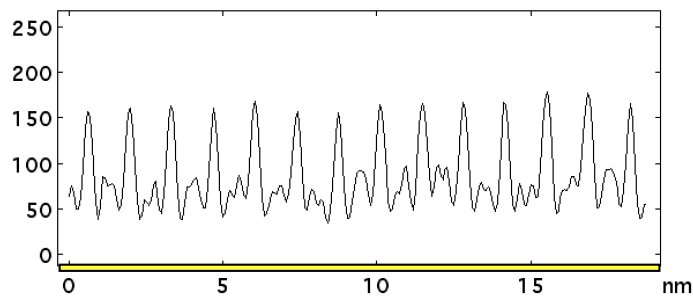
Fourier transforms

- 1D / 2D FFT
- diffraction patterns

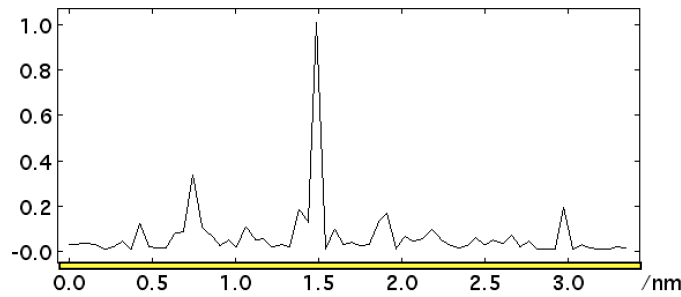
profiles \rightarrow 1D FFT \rightarrow ACF



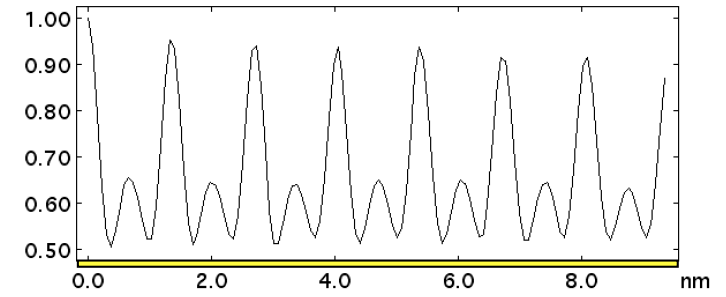
grey values (GV)



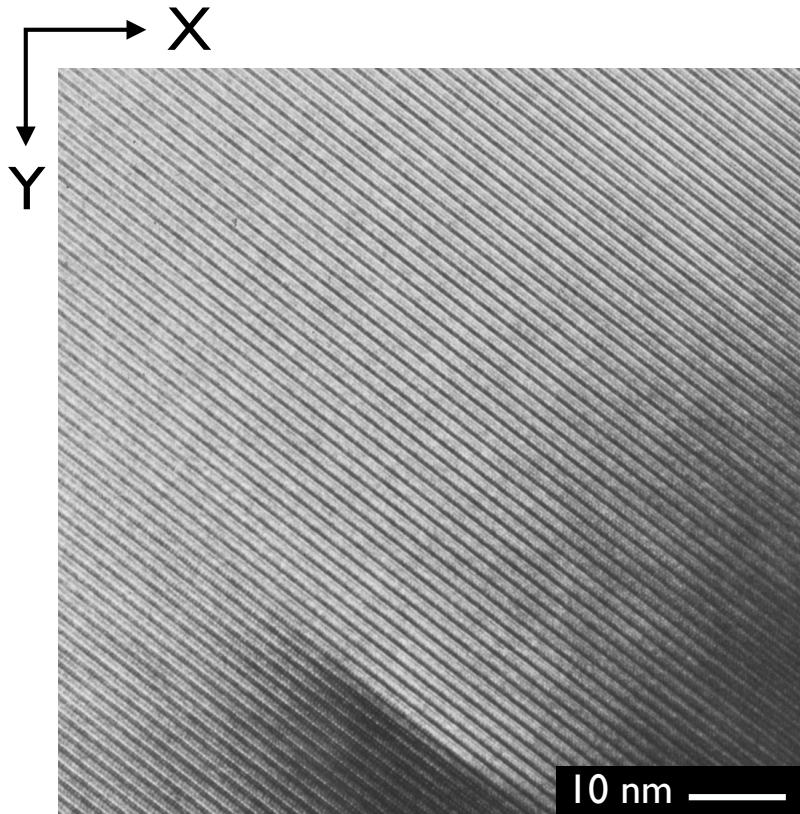
Fourier transform (FFT)



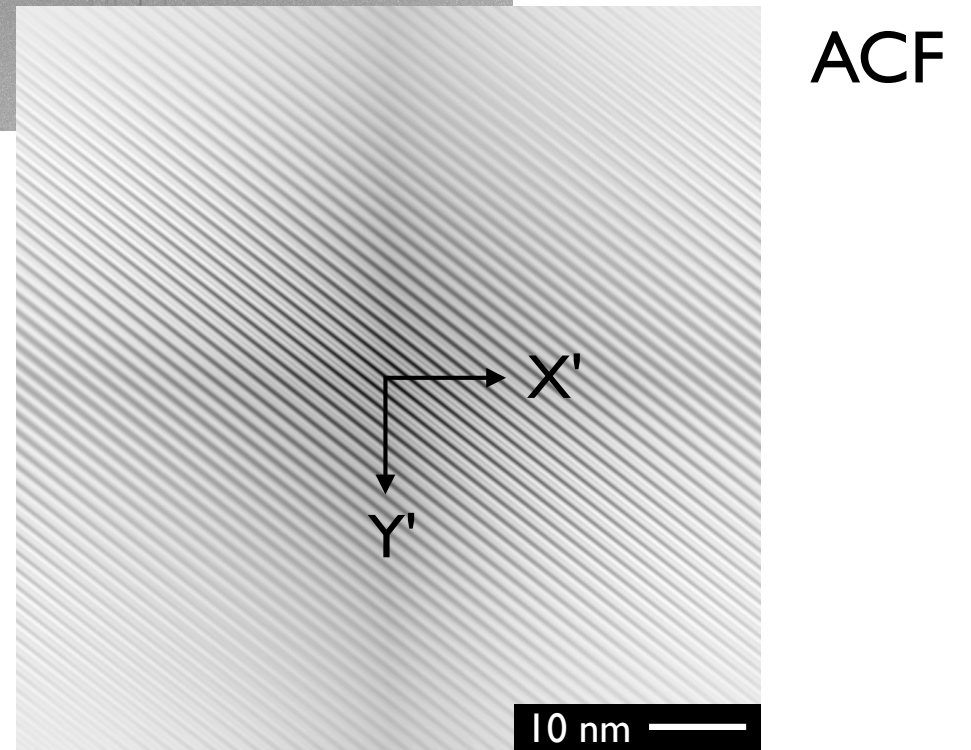
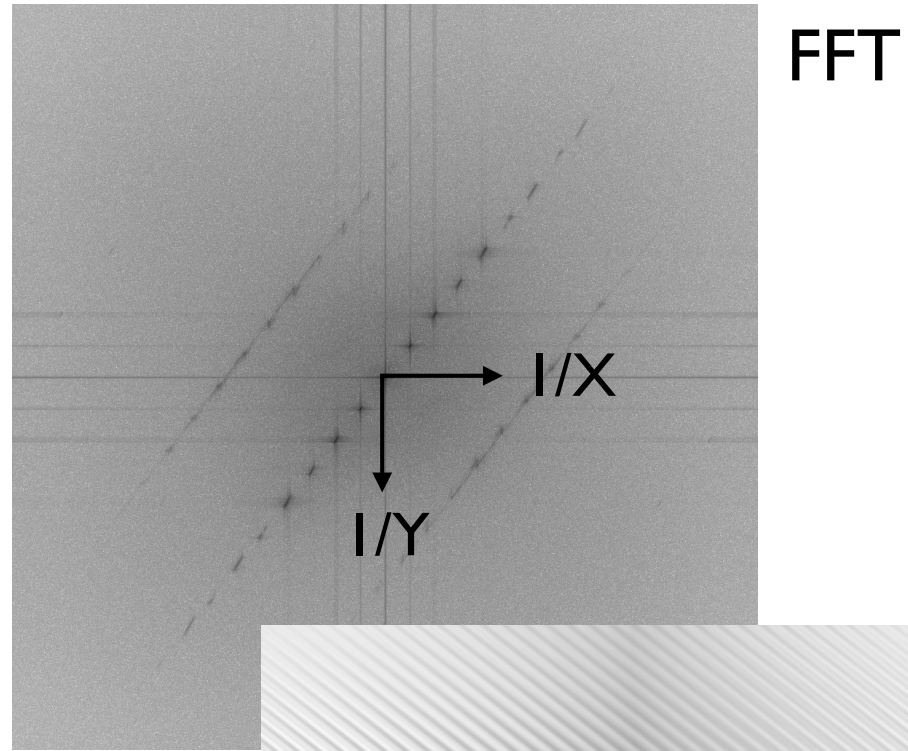
autocorrelation (ACF)



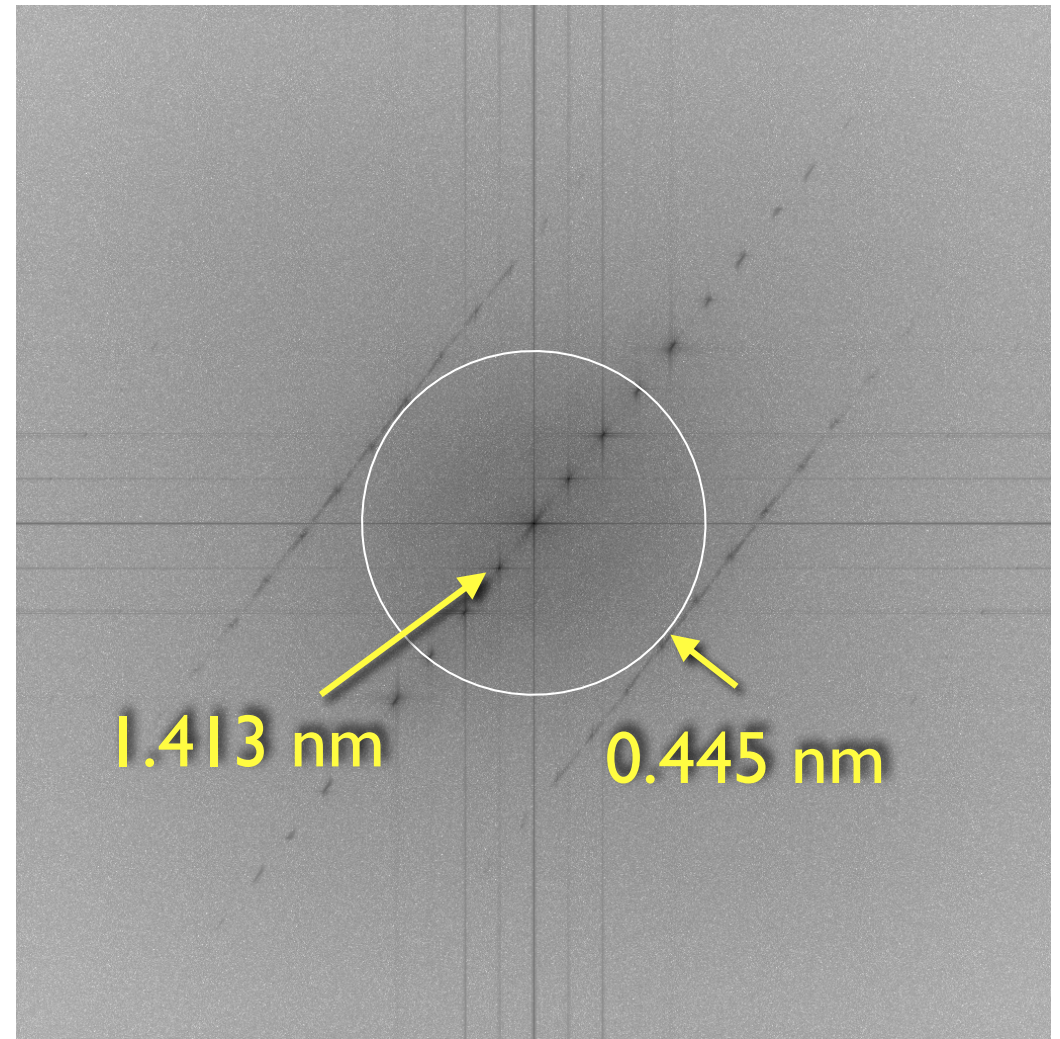
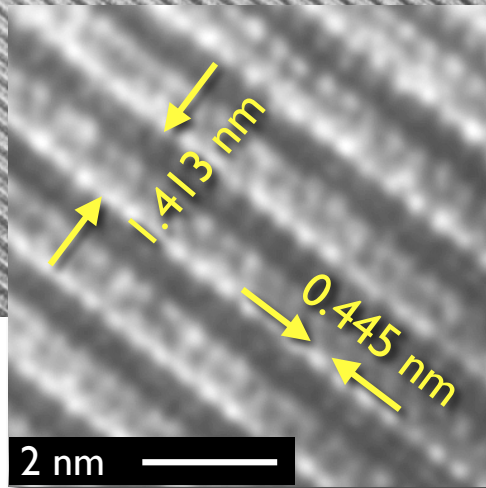
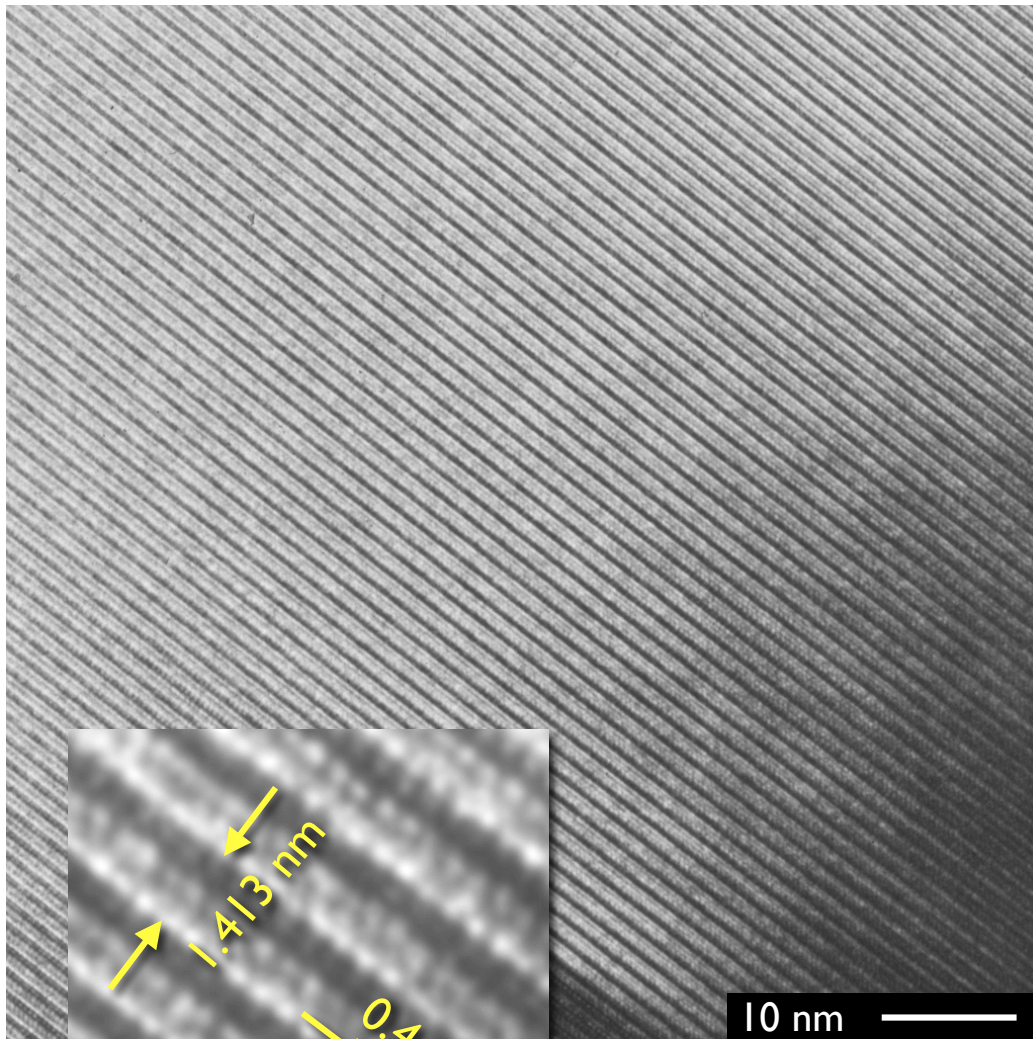
2D FFT and ACF



*high resolution TEM image of chlorite
courtesy Andreas Kronenberg*

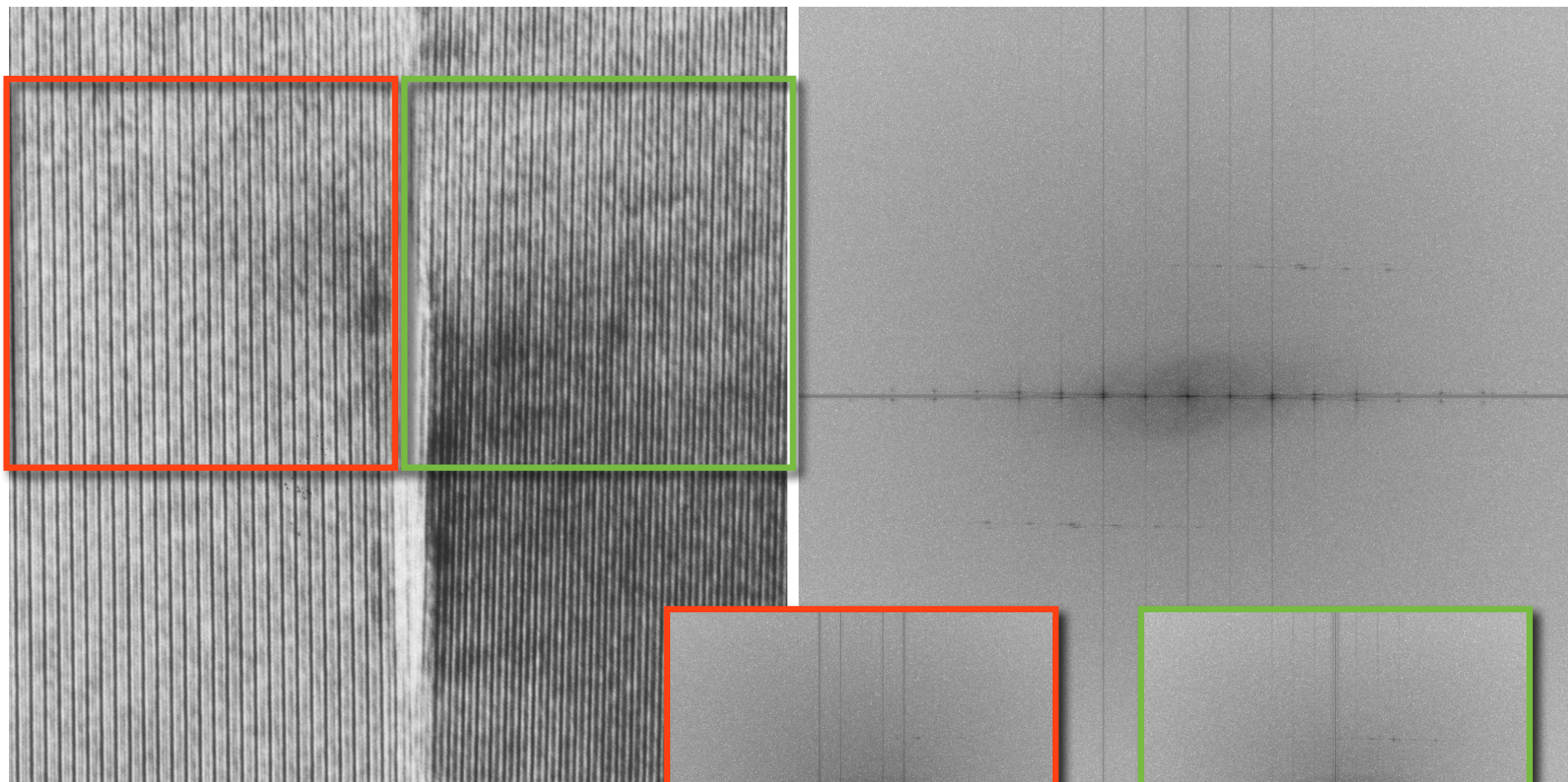


FFT - diffraction patterns



*high resolution TEM image of chlorite
courtesy Andreas Kronenberg*

FFT - low angle boundary



high resolution TEM image of chlorite
courtesy Andreas Kronenberg

Info			
X:	-83	Val:	169
Y:	3	Log:	5.03
Frq:	2.205 per nm		
Per:	0.45 nm		177.9°

Info			
X:	-84	Val:	173
Y:	-2	Log:	5.14
Frq:	2.230 per nm		
Per:	0.45 nm		181.4°

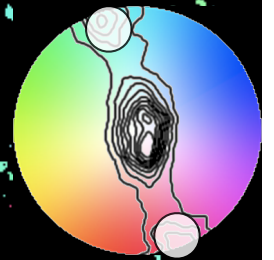
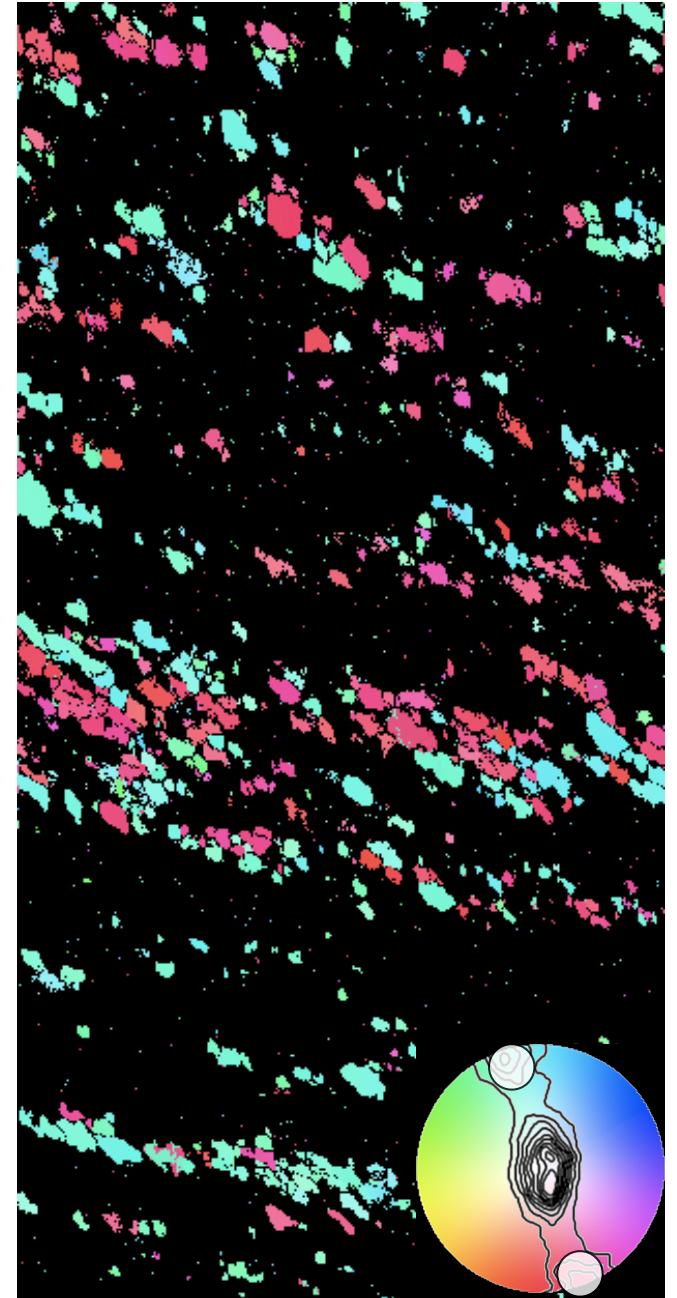
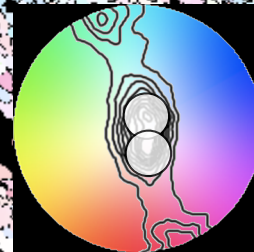
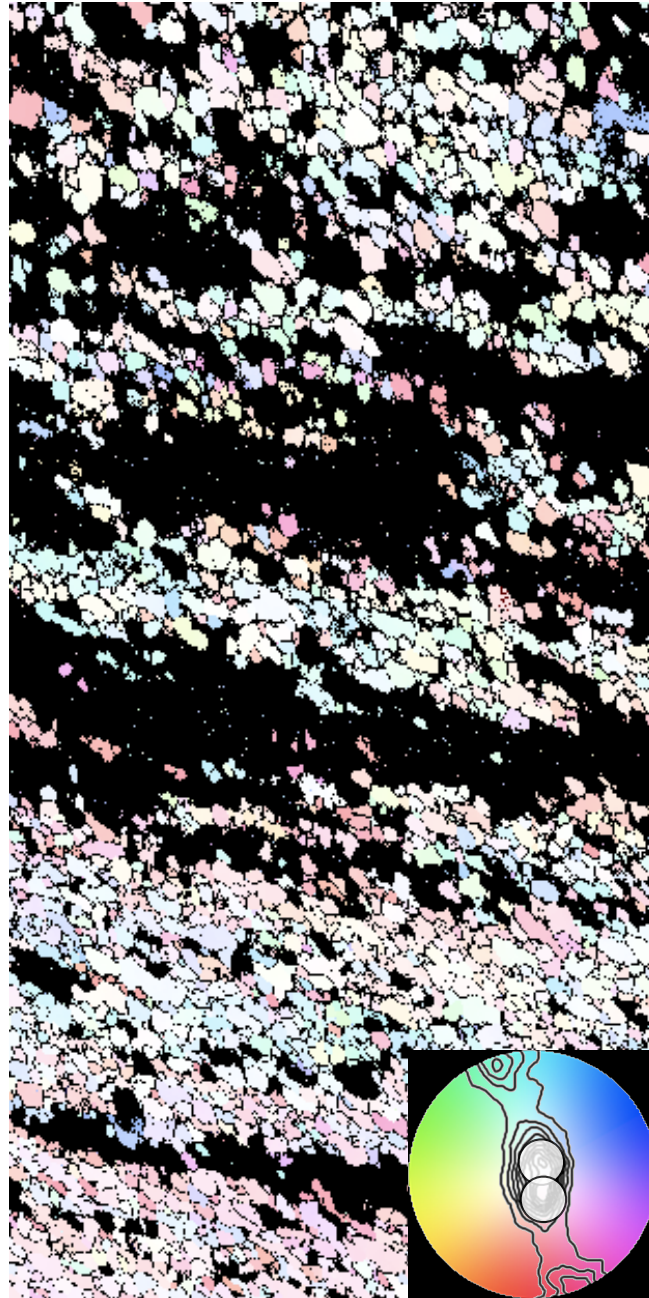
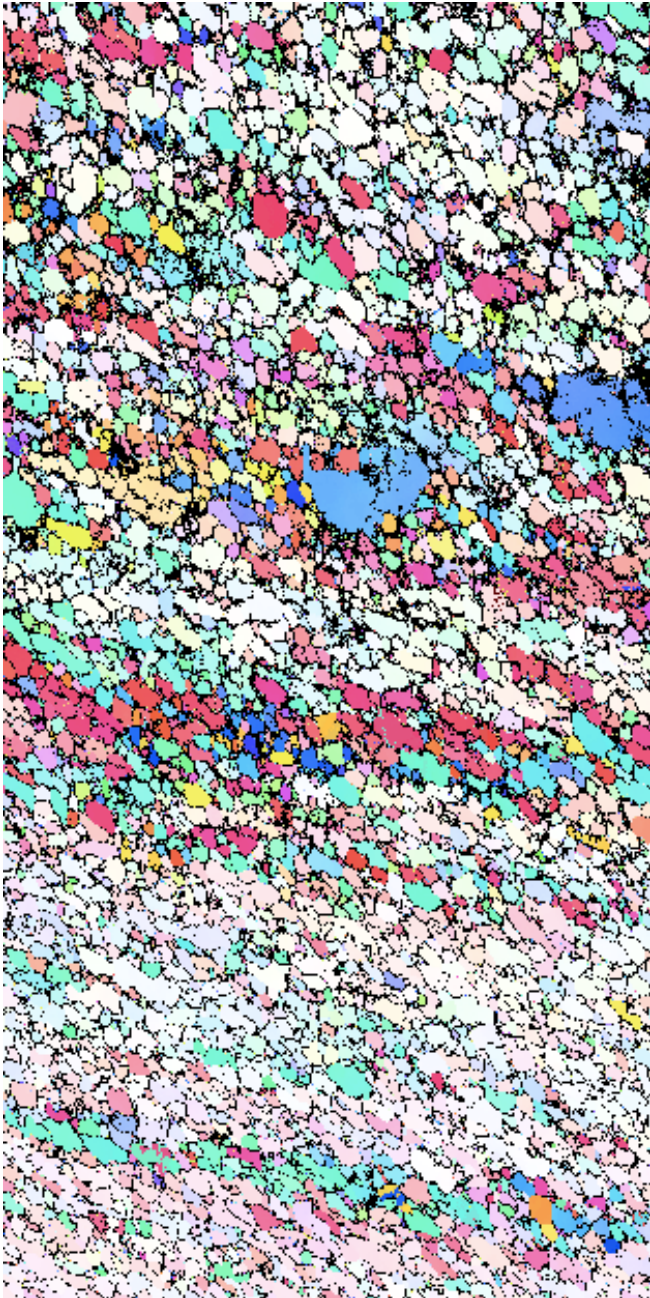
4

autocorrelation

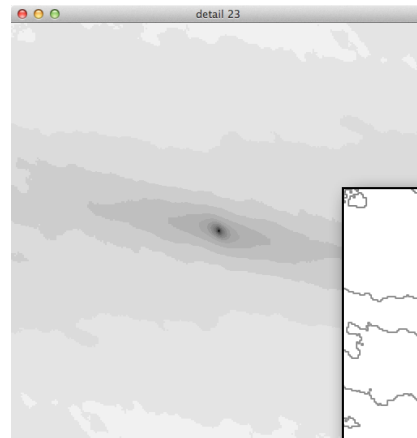
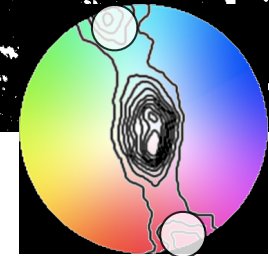
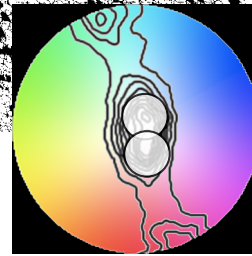
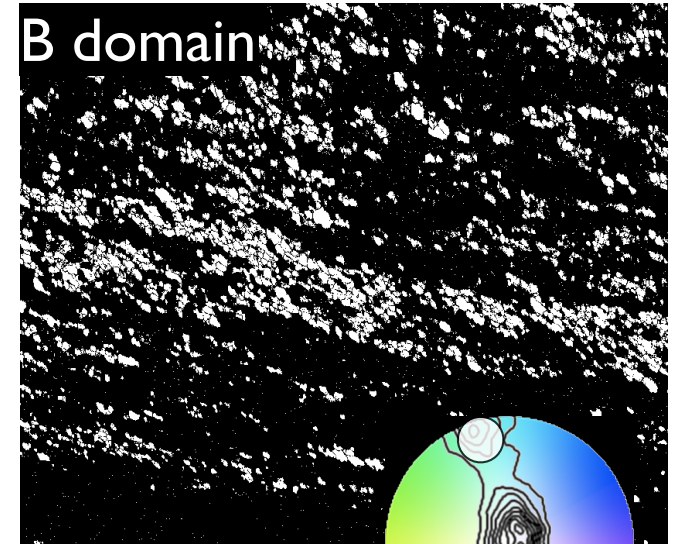
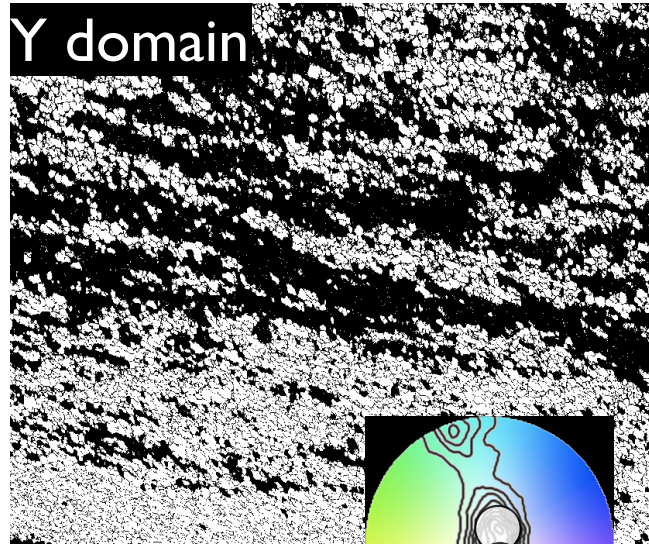
autocorrelation

- shape
- strain
- grain size
- spatial distribution

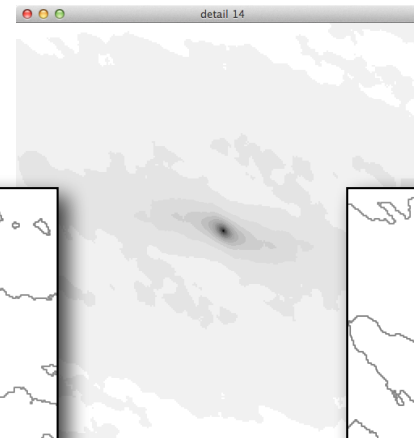
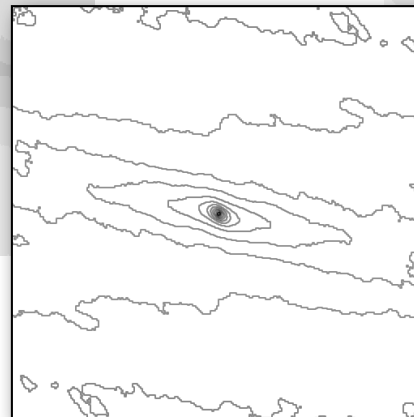
density slicing in orientation space



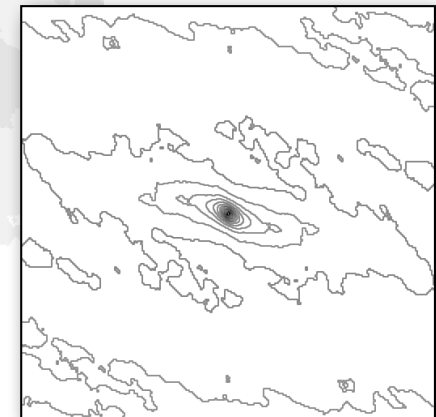
ACF - grain shape as $f(\text{CPO})$



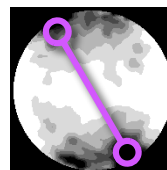
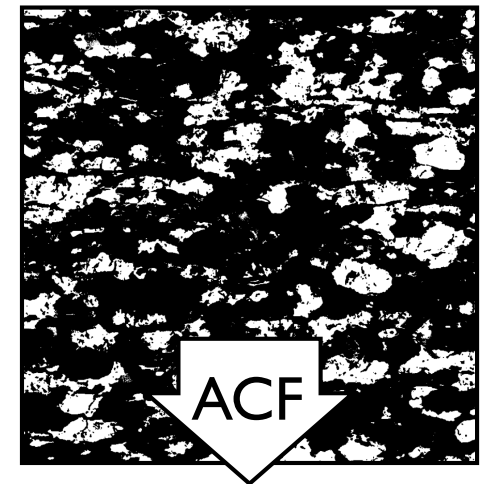
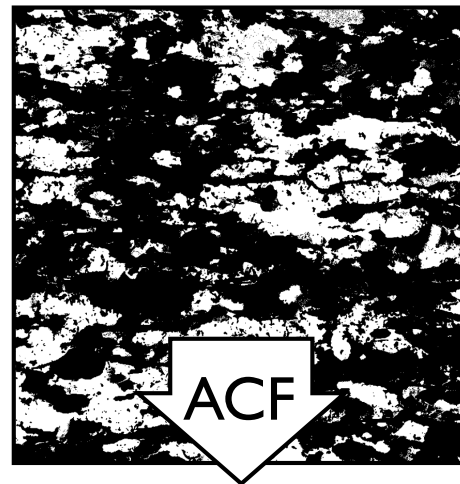
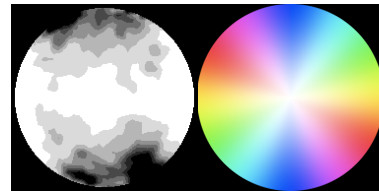
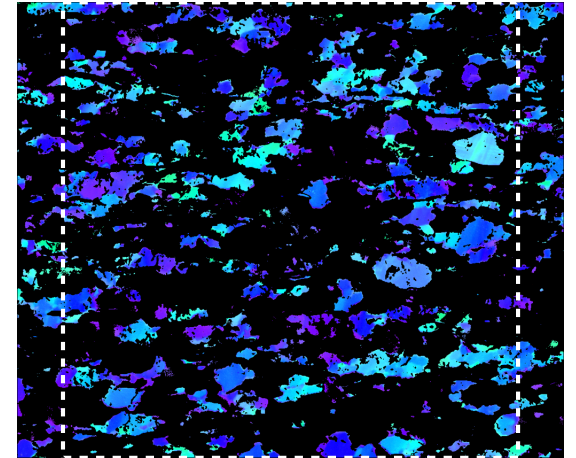
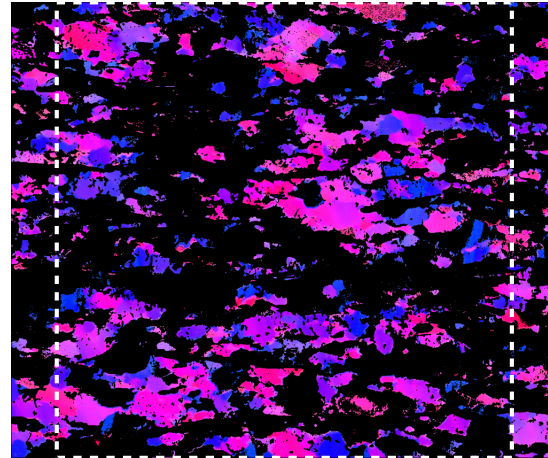
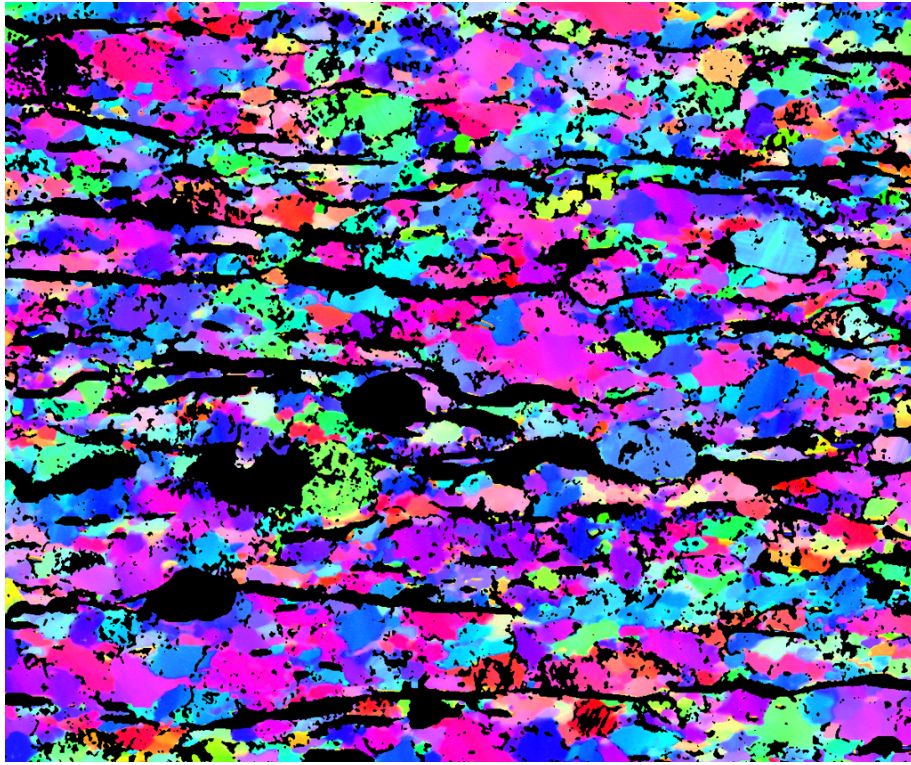
ACF of
Y domain



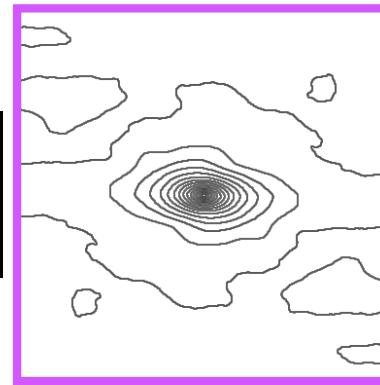
ACF of
B domain



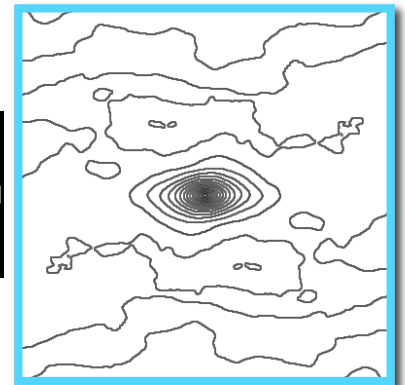
ACF - shape of texture domains



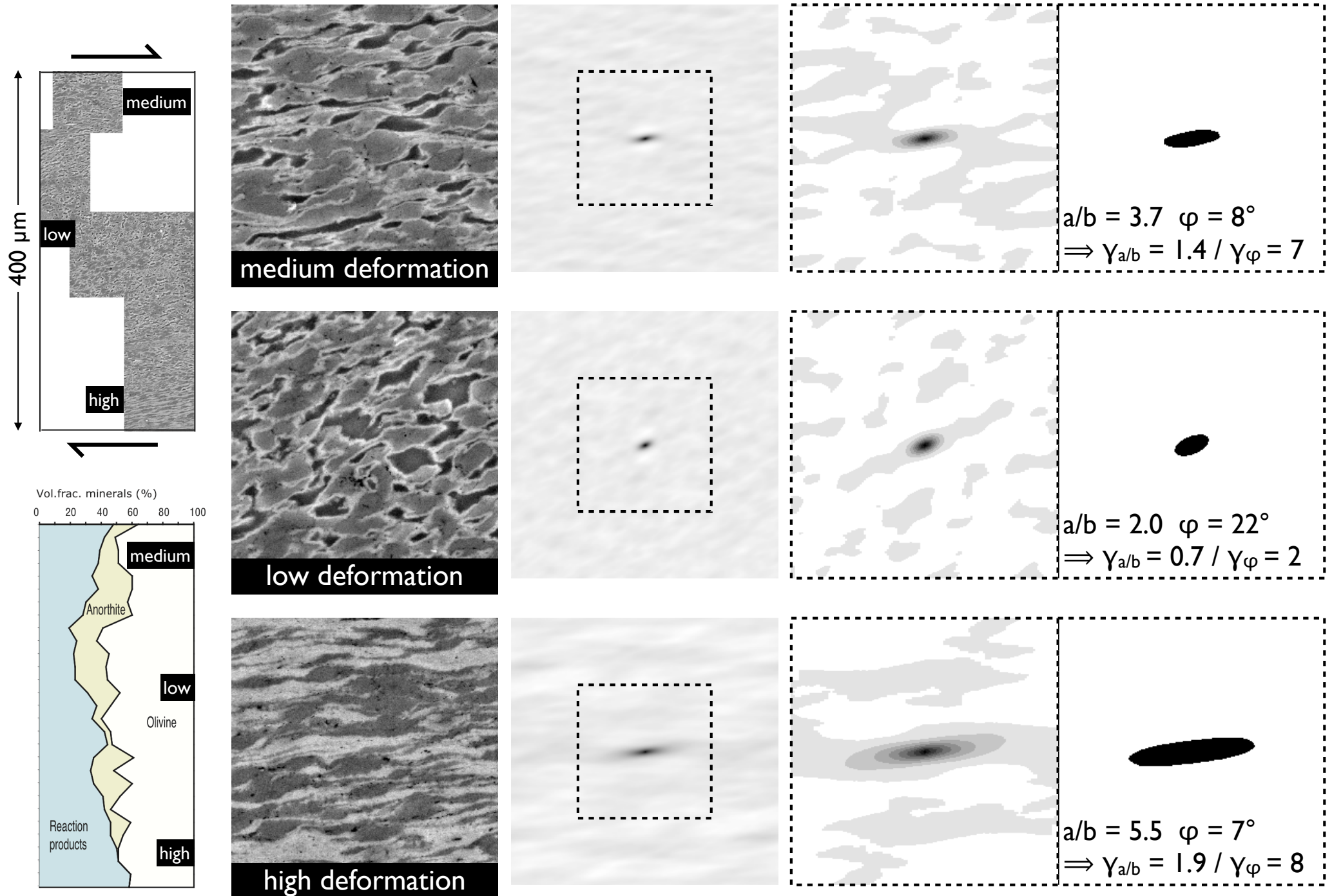
sinistral domain



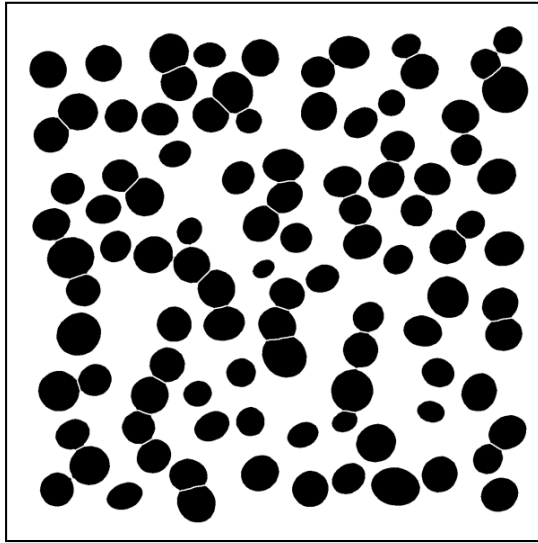
dextral domain



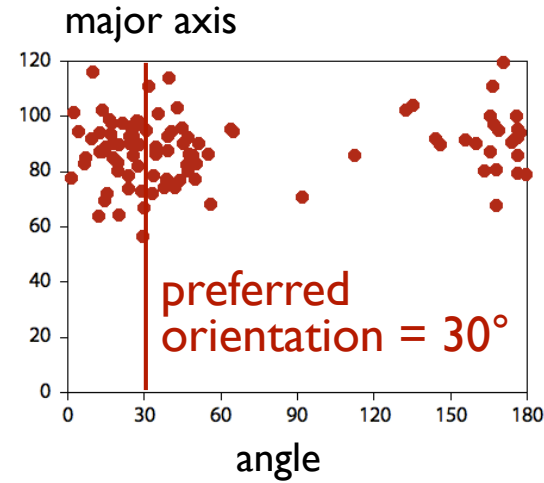
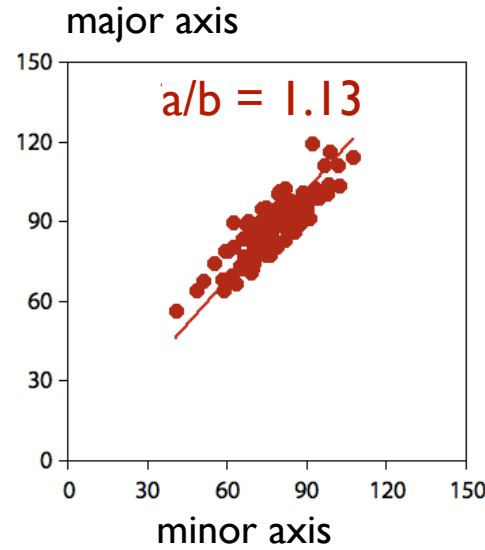
ACF → anisotropy/orientation → strain



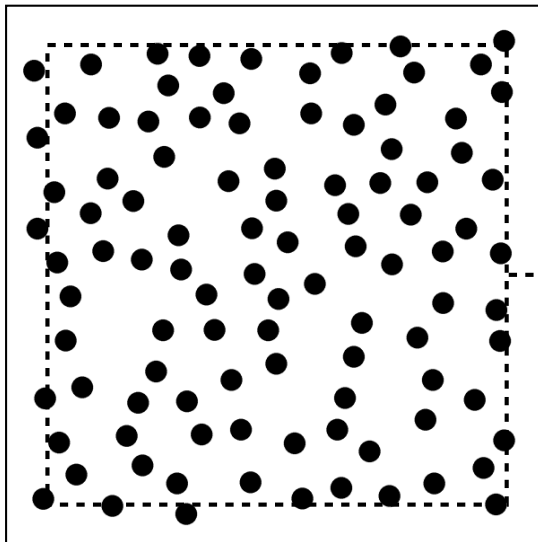
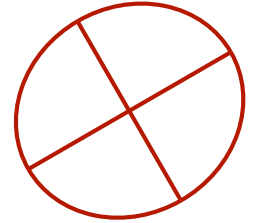
centerpoint distribution



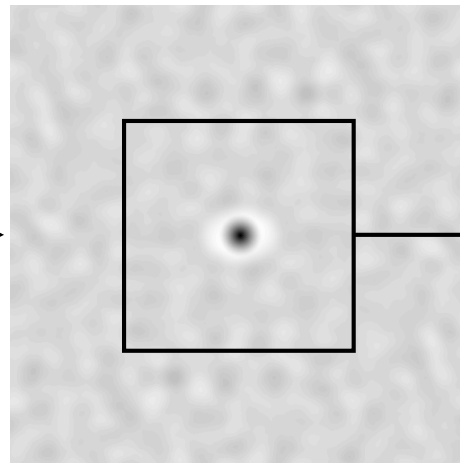
ooides bitmap



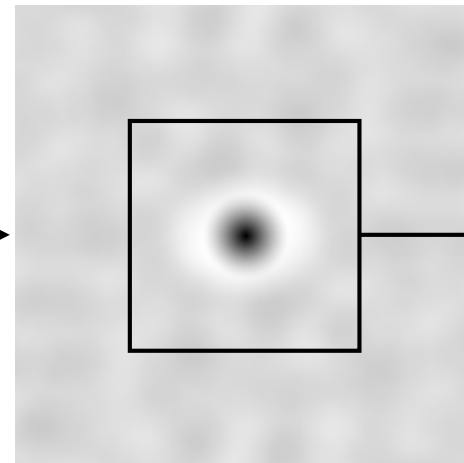
shape of particles



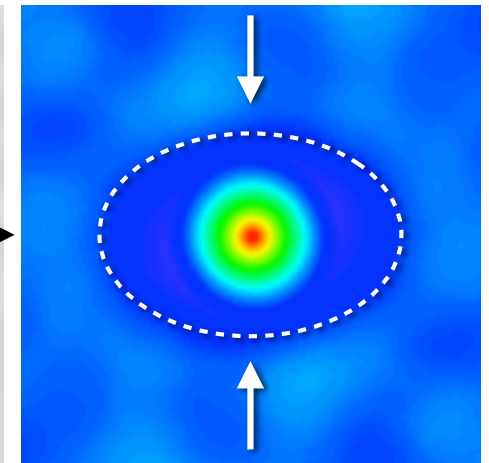
plot of centerpoints



ACF



matrix compaction



'halo' = anticorrelation of centerpoints
(→ Fry plot)

end
spatial