



TRANSITIONS IN PLATEAU BORDERS AND

TOPOLOGICAL INSTABILITIES OF DEFORMED BUBBLE CLUSTERS

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TRANSITIONS IN PLATEAU BORDERS

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LIQUID FOAMS

Liquid Foams = clusters of bubbles

Bubbles - gas inside ; surrounded by liquid films

Edges meet at vertices

Liquid in Plateau borders = PBs



foams: dry i liquid films of negligible thickness meet in PB wet i liquid films of finite thickness meet in PB

2D foams - PBs run perpendicular to the plane of the foam PBs are polygonal regions with curved edges PBs are bounded by liquid surfaces (=interfaces between the liquid and the gas) of tension γ_L

 γ = film tension and γ =2 γ_L (PB tangent to film)

Wet foam: internal PBs - surrounded by bubbles peripheral PBs - border outside gas n=3

Internal PBs:

Small liquid fraction - PBs are triangular n=3
High liquid fraction - PBs with n=3, 4 or 5 (we did not study n>5)



□ experiments with foam clusters –

change the liquid fraction

Calculations with the Surface Evolver

Experimental procedure

Monolayer technique (Smith 1952; Vaz and Fortes 1997):

- **Vessel with detergent solution**
- **Blow air bubbles with a graduated syringe on a plate**
- **The plate was inverted onto the liquid in the vessel**
- Bubbles make contact with the liquid pool; gap between liquid surface and the plate = height h of the bubbles
- The liquid surface height can be varied by adding or removing solution \Rightarrow change the height *h* of the bubbles

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experimental set-up

detergent solution

h = gap between liquid surface and the plate or height of the cluster

Bubble volumes, $V_G = 320$ or 400 mm³ 3, 4 and 5 bubble clusters of the same volume Change the liquid fraction Small h: h \approx a (a =the capillary length= 1.85 mm) $a = (\gamma_L / \rho_L g)^{1/2}$ \Rightarrow clusters with an internal PB surrounded by n bubbles + n peripheral 3-PBs Increase h; observe the evolution of PBs and measured with image analysis software : - A₁ (total area of the internal PBs) - A_G (area of one bubble)

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Experimental results

Evolution of PB with n-fold symmetry surrounded by n identical bubbles:

Ex: n=3







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Experimental results

 \Box n= 3 – PB remains symmetrical at any h; for either positive or negative p_L - p_0

n=4 – increasing h —

4 =square PB (4-fold axis of symmetry) changes p_L-p₀=0 into 4' a rectangular PB (two mirror symmetry planes) two triangular (3+3) PBs of the same size (two mirror symmetry planes)

reversible sequence

□n=5 small h ⇒ configuration 5 - 5 fold symmetric PBs larger h ⇒ p_L - p_0 <0 ⇒ configurations of lower symmetry configuration 5' - mirror plane of symmetry configuration 5'' - mirror plane of symmetry configuration 5''' - no symmetry dissociation of PB configuration (4+3) mirror plane of symmetry configuration (4+3)' no symmetry

configuration (3+3+3) mirror plane of symmetry

Sequence 1: $5 \rightarrow 5' \rightarrow (4+3) \rightarrow (4+3)' \rightarrow (3+3+3)$ Sequence 2: $5 \rightarrow 5' \rightarrow 5'' \rightarrow 5'' \rightarrow (3+3+3)$



Surface Evolver calculations

Ken Brakke (1992)

equilibrium configurations

surface energy

excess pressure

- 1. Network: coordinates of vertices; define edges and cells
- 2. Perturbations in vertices = shuffling (jiggle command)
- 3. Apply T1 (neighbor switching) in short side edges (topological changes)
- 4. Relaxation



Analytical solutions for n-fold symmetrical clusters without peripheral PBs are easy to obtain (Eur.Phys.JE,2005)

Plots of reduced energy $E/(\gamma A_G^{1/2})$ vs $(A_L/A_G)^{1/2}$

for n-fold symmetrical clusters n=3,4,5

Analytically (curves)

Surface Evolver (symbols)



✓ Curves exhibit a minimum

✓ Surface Evolver reproduces the exact results to within 0.1%

 $\checkmark \text{Plots}$ of the excess pressure $\textbf{p}_{L}\textbf{-}\textbf{p}_{0}$ in the internal PBs exhibits a maximum

 \checkmark Energy minimum occurs when p_L-p₀ =0



Starting configuration:

-internal PBs surrounded by n quadrilateral bubbles

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-internal PBs = regular polygons with n straight sides
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-peripheral PBs = triangles
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Linear mode

Perturbations in vertices = shuffling (jiggle command) Apply T1 (neighbor switching) in short side edges (topological changes) **7**Relaxation

 $\underline{\text{MINIMISE ENERGY}} = \gamma_{L}L_{L} + \gamma_{F}L_{F}$

- shuffling/relaxation procedure = 100
- 6 levels of refinement

$$\gamma = 2 \gamma_L$$

 L_{I} and L_{F} are total lengths of the PBs surfaces and the films



SURFACE EVOLVER results:

n=3 Without and with peripheral PBs



 $E/(\gamma A_G^{1/2})$ = 8.768 =8.656

The peripheral PBs lower the cluster energy

SURFACE EVOLVER results: n=4

Surface Evolver yielded the configurations found in experiments

Adding peripheral PBs does not change the geometry/topology of the internal PBs

 $> A_L/A_G >$ that of the minimum of the energy curve \Rightarrow <u>higher</u> energy configurations (not discussed here)

 $>A_L/A_G <$ that of the minimum of the energy curve \Rightarrow <u>lower</u> energy configurations



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SURFACE EVOLVER results: n =5













3+3+3

4'+3



SURFACE EVOLVER results:



SURFACE EVOLVER results: n = 5

 ${\it \ensuremath{\mathbb R}}$ A_L/A_G < that of the minimum of the symmetric curve \Rightarrow lower energy configurations

close to the minimum configuration 5' is the one of lowest energy

ﷺ it is beaten at lower A_L/A_G first by 5[™] and then by 5[™] (consistent with experimental sequence) : 5 →5[™] →5[™] →(3+3+3)

 \Re configurations (4+3) and (4+3)' have approximately the same energy; none of these is ever that of lowest energy

(beaten by (3+3+3) or 5")

the Surface Evolver sequence on increasing A_L/A_G is the experimental sequence 2; sequence 1 was not found

Discussion

□ n=4 : the sequence of equilibrium configurations on decreasing A_L/A_G is 4 → 4' →(3+3) on experiments and with the Surface Evolver

The transitions do not occur at the A_L/A_G predicted by Surface Evolver due to deviations from 2D

□ n=5 ; various configurations of a 5-sided PB;

two experimental sequences but only <u>one</u> Surface Evolver sequence

□ 4- and 5- bubble clusters break their symmetry when the PB excess pressure vanishes. The energy is minimal at this point

For liquid fractions smaller than these minima there are alternative lower energy configurations

□ similar to petal clusters; break symmetry (Weaire, Graner , Cox) at energy minimum