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## Variational Problems with Percolation

Gradient Random Fields

May 31, 2011 BIRS, Banff

**Discrete system:** with discrete variables  $u = \{u_i\}$  indexed on a lattice (e.g.,  $\Omega \cap \mathbf{Z}^d$ )

**Discrete energy:** (e.g., pair interactions)

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**B.  $\Gamma$ -convergence for Beginners, OUP 2002**

**B. Handbook of  $\Gamma$ -convergence (Handbook of Diff. Eqns, Elsevier, 2006)**

**Cubic lattice:** variables parameterized on  $\Omega \cap \mathbf{Z}^d$

**Binary systems:** variable taking only **two values**; wlog  $u_i \in \{-1, 1\}$  (**spins**).

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**Only two possible energies** (up to affine change of variables):

$$E(u) = E_{\text{ferr}}(u) = - \sum_{\text{NN}} u_i u_j \quad (\text{ferromagnetic energy})$$

(with two trivial minimizers  $u_i \equiv 1$  and  $u_i \equiv -1$ )

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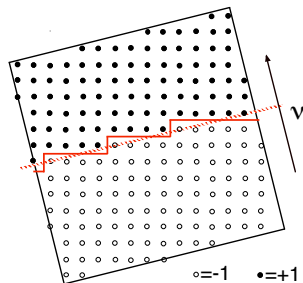
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**Note:** the change of variables  $v_i = (-1)^i u_i$  is such that  $E_{\text{anti}}(v) = E_{\text{ferro}}(u)$ , so actually we have only one energy

**Choice of the parameter:** (magnetization)  $u \in BV(\Omega; \{\pm 1\})$  continuous limit of piecewise-constant interpolations of  $\{u_i\}$

**Surface scaling:** (crystalline perimeter)

$$E_\varepsilon(u) = \sum \varepsilon^{d-1} (1 - u_i u_j) \longrightarrow 2 \int_{\Omega \cap \partial\{u=1\}} \|\nu\| d\mathcal{H}^{d-1}, \quad \text{with} \quad \|\nu\| = \sum_k |\nu_k|$$



$\nu$  = normal to the interface

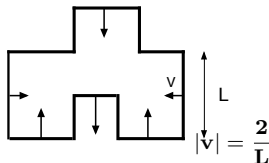
## Continuous “flows” of the perimeter

Crystalline perimeter-driven motion of sets



motion by crystalline mean curvature

(Almgren-Taylor J. Diff. Geom. 1995 in 2D)



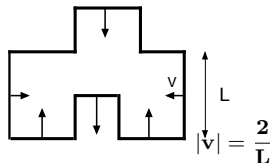
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Motion is obtained by introducing a discrete time-step  $\tau$  and initial set  $A_0$ , define a time-discrete motion by successive minimizations for fixed  $\tau$ :  $A_{k+1}$  minimizes

$$\min \left\{ P(A) + \frac{1}{2\tau} \text{“dist}(A, A_k)\text{”} \right\}$$

Define  $A^\tau(t) = A_{[t/\tau]}$  (piecewise-constant interpolation of  $\{A_k\}$ ) and pass to the limit as  $\tau \rightarrow 0$  to get a continuous  $A(t)$  (scheme by **Almgren-Taylor-Wang**, **SIAM J. Control Opt.** 1983)

## Motion of discrete interfaces

Fix  $\varepsilon$ ,  $\tau$  and  $A_0$ . Then  $A_{k+1}$  minimizes (here,  $A = \{u = 1\}$ ,  $P_\varepsilon(A) = E_\varepsilon(u)$ )

$$\min \left\{ P_\varepsilon(A) + \frac{1}{2\tau} \text{“dist}_\varepsilon(A, A_k)\text{”} \right\}$$

Define  $A^{\varepsilon, \tau}(t) = A_{[t/\tau]}$  and pass to the limit as  $\tau \rightarrow 0$  to get a continuous  $A(t)$ .

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**Pinning/depinning transition:** (B-Gelli-Novaga ARMA 2009)

- For  $\tau \ll \varepsilon$  the motion  $A(t)$  is trivial (**pinning**):

$$A(t) = A_0$$

for all (sufficiently regular) bounded initial sets  $A_0$ ;

- For  $\varepsilon \ll \tau$  the sets  $A(t)$  follow **motion by crystalline mean curvature**.

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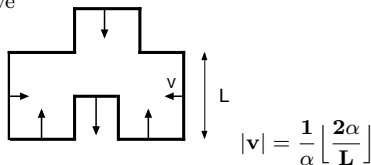
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- At the **critical scale**  $\tau = \alpha\varepsilon$  we have ‘**quantized**’ crystalline motion

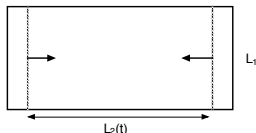


# Discreteness effects at the critical scale

(i) (**critical pinning side-length**) If all  $L > 2\alpha$  then the motion is trivial:

$$A(t) = A_0;$$

(ii) (**partial pinning and non strict inclusion principle**; e.g for rectangles) If  $L_1 < 2\alpha$  and  $L_2 > 2\alpha$  only one side is (initially) pinned



(iii) (**quantized velocity**)

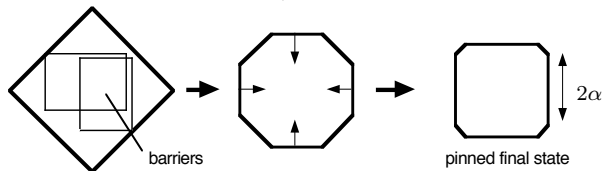
$2\alpha/L(t) \notin \mathbb{N} \Rightarrow$  velocity integer multiple of  $1/\alpha$ ;

(iv) (**non-uniqueness**)

$2\alpha/L(t) \in \mathbb{N} \Rightarrow$  velocity not uniquely determined  $\Rightarrow$  non-uniqueness

(v) (**non-convex pinned sets**)

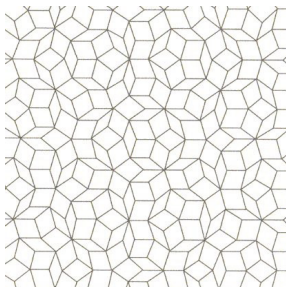
(vi) (**pinning after initial motion**)





With the due changes the process can be repeated on more general periodic lattices (e.g. triangular, hexagonal, FCC, BCC, etc.); even though we do not have in general a duality between ferro- and anti-ferromagnetic energies (**frustration**). For ferromagnetic energies we still have the **same continuous parameter**  $u \in BV(\Omega; \{\pm 1\})$ . The form of the *surface tension* changes accordingly.

Techniques must be refined to take care of **a-periodic lattices** (e.g. Penrose tilings or quasicrystals)



(B-Solci M<sup>3</sup>AS 2011)

We may have more complex interactions:

$$- \sum_{i,j} \sigma_{ij} u_i u_j$$

Conditions of the type

- **(uniform minimal states)**  $\sigma_{ij} \geq 0$
- **(coerciveness conditions)**  $\sigma_{ij} \geq c > 0$  for  $|i - j| = 1$
- **(decay conditions)**  $\sum_j \sigma_{ij} \leq C < +\infty$  for all  $i$

guarantee that (up to subsequences) the **continuous parameter** is still  $u \in BV(\Omega; \{\pm 1\})$  and

$$\sum_{ij} \varepsilon^{d-1} \sigma_{ij} (1 - u_i u_j) \longrightarrow \int_{\Omega \cap \partial\{u=1\}} \varphi(x, \nu) d\mathcal{H}^{d-1}$$

i.e., the limit is still a (possibly inhomogeneous) interfacial energy.

The integrand  $\varphi$  is determined by a family of discrete (non-local) **minimal-surface problems**. In the **2D case** and if only **nearest-neighbours** are considered ( $\sigma_{ij} = 0$  if  $|i - j| > 1$ ) equivalently it is given by an **asymptotic distance** on the lattice  $\mathbb{Z}^2$  (where the distance between the nodes  $i$  and  $j$  is  $\sigma_{ij}$ ) (**B-Piatnitsky 2010**)

When not only nearest neighbours are taken into account we do not have a correspondence between ferromagnetic and anti-ferromagnetic energies.

1) **Anti-ferromagnetic spin systems in 2D** (B-Alicandro-Cicalese NHM 2006)

$$E(u) = c_1 \sum_{\text{NN}} u_i u_j + c_2 \sum_{\text{NNN}} u_k u_l \quad u_i \in \{\pm 1\}$$

(NNN = next-to-nearest neighbours)

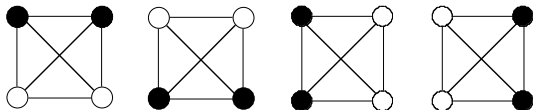
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For suitable positive  $c_1$  and  $c_2$  the ground states are 2-periodic



(representation in the unit cell)

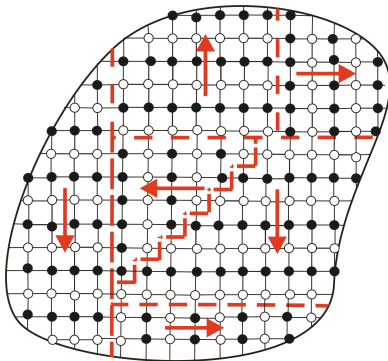
The correct order parameter is the **orientation**  $v \in \{\pm e_1, \pm e_2\}$  of the ground state.

## Surface-scaling limit

$$F(v) = \int_{S(v)} \psi(v^+ - v^-, \nu) d\mathcal{H}^1$$

$S(v)$  = discontinuity lines;  $\nu$  = normal to  $S(v)$   
 $\psi$  given by an optimal-profile problem

## Microscopic picture of a limit state with finite energy



## Ferromagnetic-anti-ferromagnetic spin systems

We can consider e.g. two-dimensional systems with NN, NNN, NNNN (next-to-next-...) interactions,  $u_i \in \{\pm 1\}$  and

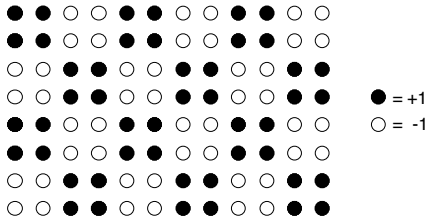
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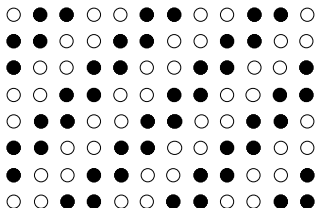
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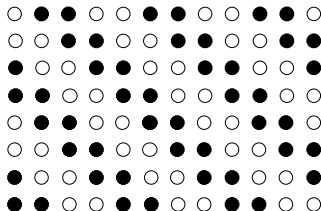
For suitable  $c_1$  and  $c_2$  again we have a non-trivial 4-periodic ground state



but also...



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(counting translations 16 different ground states)

and a description for the surface-scaling  $\Gamma$ -limit combining the two previous examples



Let  $d = 2$ . Introduce a random variable depending on an ergodic stationary discrete random process on the bonds of  $\mathbf{Z}^2$ . The simplest energy depends on its realizations  $\omega$ :

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- we can substitute 0 and  $T\nu^\perp$  with arbitrary  $x$  and  $x + T\nu^\perp$  ( $x = O(T)$ )
- oscillations of the minimal path from the segment  $[x, x + T\nu^\perp]$  are small.

**Rigid spin systems.** We may consider  $\omega$  a realization of an **i.i.d. random variable** in  $\mathbb{Z}^2$ , and the corresponding energy (surface scaling)

$$E_\varepsilon^\omega(u) = \sum_{\text{NN}} \varepsilon \sigma_{ij}^\omega (1 - u_i u_j) \quad \text{with} \quad \sigma_{ij}^\omega = \begin{cases} 1 & \text{with probability } p \\ +\infty & \text{with probability } 1 - p \end{cases}$$

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## Percolation Theorem (B-Piatnitski 2008)

In the surface scaling, the  $\Gamma$ -limit  $F_p$  of  $E_\varepsilon^\omega$  is a.s.

(1)  $F_p(u) = +\infty$  if  $u \neq 1$  or  $u \neq -1$  identically, for  $p < 1/2$

(2)  $F_p(u) = \int_{\Omega \cap \partial\{u=1\}} \varphi_p(\nu) d\mathcal{H}^1$  for  $p > 1/2$  ( $u \in BV(\Omega; \{\pm 1\})$ )

The limit is deterministic and  $\varphi_p(\nu)$  is given by an *asymptotic distance on the ‘weak cluster’* for  $p > 1/2$ .

NOTE: this is the limit case when  $\sigma_{ij}^\omega = \begin{cases} 1 & \text{with probability } p \\ T & \text{with probability } 1 - p \end{cases}$  for

$T \rightarrow +\infty$

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(1)  $F_p(u) = 0$  on all  $u \in L^1(\Omega; [-1, 1])$  for  $p \leq 1/2$

(2)  $F_p(u) = \int_{\Omega \cap \partial\{u=1\}} \varphi_p(\nu) d\mathcal{H}^1$  for  $p > 1/2$

The limit is deterministic and  $\varphi_p(\nu)$  is given by a *first-passage percolation* formula for  $p > 1/2$ .

NOTE: the parameter  $u \in BV(\Omega; \{\pm 1\})$  is the “dominant phase” (no control if  $\sigma_{ij} = 0$ )

**Ferromagnetic/antiferromagnetic interactions:** an **open problem** is when

$$E^\omega(u) = - \sum_{\text{NN}} \sigma_{ij}^\omega u_i u_j \quad \text{with} \quad \sigma_{ij}^\omega = \begin{cases} 1 & \text{with probability } p \\ -1 & \text{with probability } 1 - p \end{cases}$$

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**Deterministic ‘toy’ problem (for the case  $p \sim 0$ ):** discrete ‘perforated domain’ with well-separated ‘holes’ where  $\sigma_{ij} = -1$  (**B-Piatnitski 2010**). In this case

- need **stronger separation conditions** between the perforations
- the surface scaling is more complex and not explicit
- the  $\Gamma$ -limit may be still described by an interfacial energy  $\int_{\Omega \cap \partial\{u=1\}} \varphi(\nu) d\mathcal{H}^1$

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**Question:** How does  $p$  influence the geometry (and number) of ground states? What happens when  $p \rightarrow 1/2$ ?