

Dislocation dynamics: from microscopic models to macroscopic crystal plasticity

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Abstract. In this paper we study the connection between four models describing dislocation dynamics: a generalized 2D Frenkel-Kontorova model at the atomic level, the Peierls-Nabarro model, the discrete dislocation dynamics and a macroscopic model with dislocation densities. We show how each model can be deduced from the previous one at a smaller scale.

Key words: discrete to continuum approach, Frenkel-Kontorova model, phase field model, particle systems, periodic homogenization, Hamilton-Jacobi equations, non local equations

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1 Introduction

In this paper, we focus on the modelling of dislocation dynamics. We refer the reader to the book of Hirth and Lothe [20] for a detailed introduction to dislocations. Our study ranges from atomic models to macroscopic crystal plasticity. At each scale, dislocations can be described by a suitable model. Our goal is to explain how we can deduce a model at a larger scale, from the model at a smaller scale.

Even if our derivation will be done on some simplified models (essentially 2D and 1D models), we hope that our contribution will shed light, even on some well-known models. More precisely, we will consider the following four models, from the smaller to the larger scale:

1. Generalized Frenkel-Kontorova model (FK)
2. Peierls-Nabarro model (PN)
3. Dynamics of discrete dislocations (DDD)
4. Dislocation density model (DD)

Schematically, the four models are related as shown below (see also Figure 6.1 for a more detailed diagram):

$$\boxed{(\text{FK})_{\varepsilon_1, \varepsilon_2, \varepsilon_3 > 0}} \xrightarrow{\varepsilon_1 \rightarrow 0} \boxed{(\text{PN})_{\varepsilon_2, \varepsilon_3 > 0}} \xrightarrow{\varepsilon_2 \rightarrow 0} \boxed{(\text{DDD})_{\varepsilon_3 > 0}} \xrightarrow{\varepsilon_3 \rightarrow 0} \boxed{(\text{DD})} \quad (1.1)$$

The rest of the paper is composed of four sections. Each section presents one model, and explains how this model can be deduced from the previous model at a smaller scale.

2 Generalised Frenkel-Kontorova model

2.1 Geometrical description

We call (e_1, e_2, e_3) a direct orthonormal basis of the threedimensional space. We consider a perfect crystal \mathbb{Z}^3 where each position with integer coordinates is occupied by one atom. We want to describe dislocations, which are certain “line defects” in the crystal. Even if it possible to write down a full 3D model at this level of description, the whole analysis (including the change of scales) in this paper will be carried out only for the case of straight dislocation lines. For this reason, and in order to simplify the presentation, we will assume that the material is invariant by integer translations in the direction e_3 . Because of this assumption, we can simply consider the cross section of the crystal in the plane (e_1, e_2) where each atom is now assumed to have a position $I \in \mathbb{Z}^2$ in the perfect crystal. We also assume that each atom I can have a displacement $U_I \in \mathbb{R}$ in the direction e_1 , such that the effective position of the atom I is $I + U_I e_1$.

On Figure 2.1 below is represented a view of the perfect crystal. On Figure 2.2 we can see a schematic view of an edge dislocation in the crystal. On this picture, the upper part $\{I_2 \geq 0\}$ of the crystal has been expanded to the right of a vector $\frac{1}{2}e_1$, while the lower part $\{I_2 \leq -1\}$ of the crystal has been contracted to the left of a vector $-\frac{1}{2}e_1$. The net difference between these two vectors is e_1 and is called the Burgers vector of this dislocation.

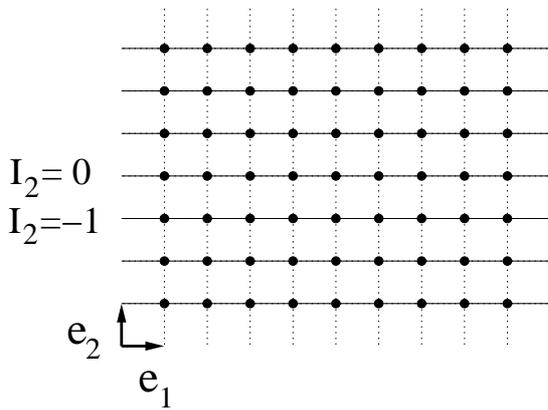


Fig. 2.1. Perfect crystal

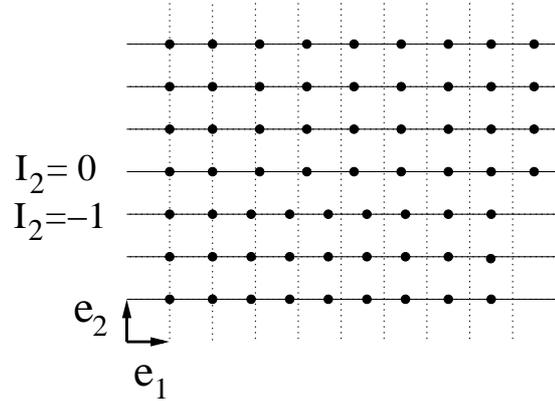


Fig. 2.2. Schematic view of an edge dislocation in the crystal

In order to describe an edge dislocation in our formalism, let us make a few assumptions. We will assume that the dislocation defects are essentially described by the mismatch between the two planes $I_2 = 0$ and $I_2 = -1$, like on Figure 2.2. For this reason, and also in order to simplify the analysis, we assume that the displacement of the crystal satisfies the following *antisymmetry property*

$$U_{(I_1, -I_2)} = -U_{(I_1, I_2-1)} \quad \text{for all } I = (I_1, I_2) \in \mathbb{Z}^2. \quad (2.1)$$

Let us also define the discrete gradient

$$(\nabla^d U)_I = \begin{pmatrix} U_{I+e_1} - U_I \\ U_{I+e_2} - U_I \end{pmatrix}.$$

Remark that defects in the crystal can be seen as regions where the discrete gradient is not small.

Formalism for an edge dislocation with Burgers vector e_1

In our formalism, an edge dislocation like the one of Figure 2.2, can be represented by a displacement U_I satisfying

$$\begin{cases} U_{(I_1, 0)} = -U_{(I_1, -1)} \rightarrow 0 & \text{as } I_1 \rightarrow -\infty \\ U_{(I_1, 0)} = -U_{(I_1, -1)} \rightarrow \frac{1}{2} & \text{as } I_1 \rightarrow +\infty. \end{cases}$$

Because we assume that the dislocation core lies in the two planes $I_2 = 0$ and $I_2 = -1$, it is reasonable to assume that all the components of the discrete gradient are small, except components $U_{I+e_2} - U_I$ for $I = (I_1, I_2)$ with $I_2 = -1$. More precisely, we assume that there exists a small $\delta > 0$ such that

$$\begin{cases} |U_{I+e_1} - U_I| \leq \delta & \text{for all } I = (I_1, I_2) \in \mathbb{Z}^2 \\ |U_{I+e_2} - U_I| \leq \delta & \text{for all } I = (I_1, I_2) \in \mathbb{Z}^2 \text{ with } I_2 \neq -1. \end{cases} \quad (2.2)$$

Moreover, if there is no applied stress on the crystal, then it is reasonable to assume that

$$\text{dist}((\nabla^d U)_I, \mathbb{Z}^2) \rightarrow 0 \quad \text{as } |I| \rightarrow +\infty.$$

2.2 The energy and the dynamics

We assume that the energy of a configuration $U = (U_I)_{I \in \mathbb{Z}^2}$ of the crystal can be formally written as

$$E(U) = \frac{1}{2} \sum_{I \neq J} \widetilde{W}(U_I - U_J)$$

where $\widetilde{W} : \mathbb{R} \rightarrow \mathbb{R}$ is a potential describing nearest neighbors interactions satisfying

Assumption (A1)

$$\begin{cases} \text{i) (Regularity)} & \widetilde{W} \in C^3(\mathbb{R}) \\ \text{ii) (Periodicity)} & \widetilde{W}(a+1) = \widetilde{W}(a) \quad \text{for all } a \in \mathbb{R} \\ \text{iii) (Minimum on } \mathbb{Z}) & \widetilde{W}(\mathbb{Z}) = 0 < \widetilde{W}(a) \quad \text{for all } a \in \mathbb{R} \setminus \mathbb{Z} \\ \text{iv) (Local harmonicity of } \widetilde{W}) & \widetilde{W}(a) = \frac{1}{2}a^2 \quad \text{for all } |a| < \delta \end{cases}$$

where $\delta > 0$ is introduced in (2.2). Remark that the periodicity of the potential \widetilde{W} reflects the periodicity of the crystal, while the minimum property of \widetilde{W} is consistent with the fact that the perfect crystal \mathbb{Z}^2 is assumed to minimize its energy. Assumption iv) will be used for later simplification. Then we assume that we are in a regime where the crystal reaches very quickly the equilibrium in the regions where there is no defects, i.e. satisfies

$$0 = -\nabla_{U_I} E(U) \quad \text{for all } I = (I_1, I_2) \in \mathbb{Z}^2 \text{ with } I_2 \neq 0, -1. \quad (2.3)$$

Remark that the equilibrium condition (2.3) may look artificial. Indeed this condition is also assumed in order to carry out the analysis, and can be seen as a drastic simplification of reality.

In the two planes where the dislocation lives, we have the following fully overdamped dynamics (describing the average friction of the lattice on the effective dissipative motion of the dislocations):

$$\frac{d}{dt} U_I = -\nabla_{U_I} E(U) \quad \text{for all } I = (I_1, I_2) \in \mathbb{Z}^2 \text{ with } I_2 = 0, -1. \quad (2.4)$$

Let us mention that we do not have a fundamental justification of this dynamics, but we think that one of the main justification of this model is that other known models at larger scales can be deduced from this particular model. For physical justifications of the dissipative effects in the motion of dislocations, see [2, 20]. See also [22–24] for a fundamental justification of the overdamped dynamics based on explicit computations in a 1D Hamiltonian model.

Taking into account the local harmonic assumption ($\widetilde{A1}$) iv), applied where the components of the discrete gradient are small (see (2.2)), joint to the antisymmetry property defined in (2.1), we can rewrite system (2.3)-(2.4) as follows for all $t > 0$:

$$\begin{cases} 0 = \sum_{J \in \mathbb{Z}^2, |J-I|=1} (U_J - U_I) & \text{for all } I = (I_1, I_2) \in \mathbb{Z}^2 \text{ with } I_2 \geq 1 \\ \frac{d}{dt} U_I = -\widetilde{W}'(2U_I) + \sum_{J \in \mathbb{Z}^2, |J-I|=1, J_2 \geq 0} (U_J - U_I) & \text{for all } I = (I_1, I_2) \in \mathbb{Z}^2 \text{ with } I_2 = 0. \end{cases} \quad (2.5)$$

We call this model a generalised Frenkel-Kontorova model. Even if this system of equations is not standard, it is nevertheless possible to define a unique solution under suitable assumptions, in the

framework of viscosity solutions (see [13]). We refer the reader to the book of Braun, Kivshar, [6] for a detailed presentation of the classical FK model. For homogenization results of FK models, we refer the reader to [15]. For the description of vacancy defects at equilibrium, see [19]. See also [18], where the authors study the problem involving a dislocation inside the interphase between two identical lattices. Their model corresponds to our model (2.5) at the equilibrium with the potential \widetilde{W} is a cosine function. For other 2D FK models, see [8,9].

Remark 2.1 *It is important to remark that we used condition (2.2) only to derive the model. We do not know and we do not claim that there exists solutions of system (2.5) satisfying condition (2.2). From now on, we only consider solutions of system (2.5) without requiring further assumptions on the solutions.*

Remark 2.2 *When we freeze the components $U_I = 0$ for $I_2 \geq 1$, and change the evolution equation forgetting the index J with $J_2 = 1$, this leads to the following classical fully overdamped Frenkel-Kontorova model satisfied by $V_i := U_{(i,0)}$*

$$\frac{d}{dt}V_i = V_{i+1} + V_{i-1} - 2V_i - \widetilde{W}'(2V_i).$$

2.3 The asymptotic stress created by a single dislocation

In this subsection, we will compute the asymptotic stress created by a single dislocation. To this end, we first compute the effective Hook's law of the lattice.

Computation of the Hook's law

Let us consider an affine displacement

$$U_I = a \cdot I + C \quad \text{with} \quad a = (a_1, a_2) \in \mathbb{R}^2$$

where $C \in \mathbb{R}$ is a constant. Then the energy by unit cell is

$$\mathcal{E} = \widetilde{W}(U_{I+e_1} - U_I) + \widetilde{W}(U_{I+e_2} - U_I) = \frac{1}{2}(a_1^2 + a_2^2)$$

for $|a| < \delta$. Reminding the fact that U is the displacement in the e_1 direction, we get that the strain e (i.e. the symmetric part of the gradient of the displacement) is given by

$$e = \begin{pmatrix} e_{11} & e_{12} \\ e_{21} & e_{22} \end{pmatrix} = \frac{1}{2}(\nabla U \otimes e_1 + e_1 \otimes \nabla U) = \begin{pmatrix} a_1 & a_2/2 \\ a_2/2 & 0 \end{pmatrix}.$$

Therefore

$$\mathcal{E}(e) = \frac{1}{2}e_{11}^2 + 2e_{12}^2.$$

Recalling that the stress is given by $\sigma^0 = \frac{\partial \mathcal{E}}{\partial e}$, we get the Hook's law:

$$\sigma^0 = \begin{pmatrix} e_{11} & 2e_{12} \\ 2e_{21} & 0 \end{pmatrix}.$$

Computation of the stress created by a single dislocation

Remark that when there is no dislocations, the energy associated to a continuous displacement $U(X)$ for $X = (X_1, X_2)$ is formally

$$E = \int_{\mathbb{R}^2} \frac{1}{2} |\nabla U|^2.$$

Therefore the Euler-Lagrange equation (which is the corresponding equation of elasticity for this model) is

$$\Delta U = 0. \tag{2.6}$$

Let us now consider the following function

$$U_0(X) = \frac{1}{2\pi} \arctan\left(\frac{X_1}{X_2}\right) + \frac{1}{4} \text{sgn}(X_2)$$

where sgn is the sign function. This function satisfies

$$\begin{cases} U_0(X_1, X_2) = -U_0(X_1, -X_2) \\ U_0(X_1, 0^+) = -U_0(X_1, 0^-) \rightarrow 0 \quad \text{as } X_1 \rightarrow -\infty \\ U_0(X_1, 0^+) = -U_0(X_1, 0^-) \rightarrow \frac{1}{2} \quad \text{as } X_1 \rightarrow +\infty. \end{cases}$$

Moreover we can easily check that

$$\text{div} (\nabla U_0 - H(X_1)\delta_0(X_2)e_2) = 0 \quad \text{in } \mathcal{D}'(\mathbb{R}^2)$$

where H is the Heavyside function and δ_0 is the Dirac mass. This equation is the analogue of equation (2.6) when there is a dislocation. This shows that in a continuum mechanics framework associated to the particular lattice that we consider, the function U_0 is the displacement corresponding to a dislocation with Burgers vector e_1 . In particular, the stress created by this dislocation is then given by

$$\sigma^0 = \frac{1}{2\pi} \begin{pmatrix} \frac{X_2}{X_1^2 + X_2^2} & -\frac{X_1}{X_1^2 + X_2^2} \\ -\frac{X_1}{X_1^2 + X_2^2} & 0 \end{pmatrix}$$

and then

$$\sigma_{12}^0(X_1, 0) = -\frac{1}{2\pi X_1} \quad (2.7)$$

which is the asymptotic shear stress at the point $(X_1, 0) \in \mathbb{R}^2$ created by a single dislocation positioned at the origin, and with Burgers vector e_1 .

2.4 Rescaling of the generalised FK model

In this subsection, in order to simplify the notation we denote by $\varepsilon := \varepsilon_1 > 0$ the small parameter in the first passage of the scheme (1.1). We are interested in the case of asymptotically small potential \widetilde{W} for which we expect an asymptotically large dislocation core. This means that in this limit, we expect to be able to describe the discrete displacement U_I by a continuous function.

More precisely, we first define the rescaled integer coordinates:

$$\Omega^\varepsilon = (\varepsilon\mathbb{Z}) \times \varepsilon(\mathbb{N} \setminus \{0\}), \quad \partial\Omega^\varepsilon = (\varepsilon\mathbb{Z}) \times \{0\}.$$

Then we write the potential as

$$\widetilde{W} = \frac{\varepsilon}{2} W^\varepsilon$$

and define the rescaled function

$$u^\varepsilon(X, t) = 2U_{\frac{X}{\varepsilon}} \left(\frac{t}{\varepsilon} \right) \quad \text{for } X = (X_1, X_2) \in \overline{\Omega}^\varepsilon, \quad t \in [0, +\infty).$$

Remark that the factor 2 in the definition of u^ε permits to interpret u^ε as the jump of the displacement in the direction e_1 , when we pass from hyperplane $X_2 = -\varepsilon$ to the hyperplane $X_2 = 0$.

We can easily check that u^ε solves the following system of equations (with the particular value $\sigma = 0$)

$$\begin{cases} 0 = \frac{1}{\varepsilon^2} \sum_{J \in \mathbb{Z}^2, |J|=1} (u^\varepsilon(X + \varepsilon J, t) - u^\varepsilon(X, t)) & \text{for all } (X, t) \in \Omega^\varepsilon \times (0, +\infty) \\ u_t^\varepsilon(X, t) = 2\varepsilon_2 \sigma(\varepsilon_2 X_1) - (W^\varepsilon)'(u^\varepsilon(X, t)) + I^\varepsilon[u^\varepsilon](X, t) & \text{for all } (X, t) \in \partial\Omega^\varepsilon \times (0, +\infty) \\ \text{with } I^\varepsilon[u^\varepsilon](X, t) = \frac{1}{\varepsilon} \sum_{J \in \mathbb{Z}^2, |J|=1, J_2 \geq 0} (u^\varepsilon(X + \varepsilon J, t) - u^\varepsilon(X, t)). \end{cases} \quad (2.8)$$

The scalar function σ has been introduced to take into account the possible external applied shear stress on the material. Here the small parameter $\varepsilon_2 > 0$ has been introduced to take into account the slow variation of the stress field which will allow a second change of scale as ε_2 will go to zero later. We will also assume that the initial data satisfies

$$u^\varepsilon(X, 0) = u_0(X) \quad \text{for all } X \in \partial\Omega^\varepsilon \quad (2.9)$$

where u_0 is a given function independent on ε and smooth enough.

In order to identify a limit model as ε goes to zero, we also make the following assumption

$$\|W^\varepsilon - W\|_{C^1(\mathbb{R})} \rightarrow 0 \quad \text{as } \varepsilon \rightarrow 0 \quad (2.10)$$

for some new potential W satisfying the following assumption:

Assumption (A1)

$$\left\{ \begin{array}{l} \text{The potential } W \text{ satisfies } (\widetilde{\text{A1}}) \text{ i), ii), iii)} \\ \text{iv) (Non degenerate minima): } \alpha := W''(0) > 0. \end{array} \right.$$

In (2.10), we use the C^1 norm, because this is the first derivative of the potential that appears in the equations. Remark that condition (2.10) can be fulfilled, if we assume for instance that \widetilde{W} satisfies assumption $(\widetilde{\text{A1}})$ with $\delta = \delta_\varepsilon \ll \varepsilon$. We also make the following assumption on the stress:

Assumption (A2)

There exists a constant $C > 0$ such that

$$|\sigma| + |\sigma_x| + |\sigma_{xx}| \leq C \quad \text{for all } x \in \mathbb{R}.$$

3 The Peierls-Nabarro model

3.1 Description of the PN model

In this section we introduce the Peierls-Nabarro model, which is a phase field model (see [20] for a presentation of this model). In this model, phase transitions describe the dislocation cores. We set

$$\Omega = \{X = (X_1, X_2) \in \mathbb{R}^2, \quad X_2 > 0\}.$$

A function $u^0(X, t)$ is said to be a solution of the PN model, if it satisfies the following system

$$\left\{ \begin{array}{ll} 0 = \Delta u^0 & \text{on } \Omega \times (0, +\infty) \\ u_t^0 = 2\varepsilon_2 \sigma(\varepsilon_2 X_1) - W'(u^0) + \frac{\partial u^0}{\partial X_2} & \text{on } \partial\Omega \times (0, +\infty) \end{array} \right. \quad (3.1)$$

with initial data

$$u^0(X, 0) = u_0(X) \quad \text{for all } X \in \partial\Omega. \quad (3.2)$$

The stationary version of this model has been originally introduced in order to propose a method to compute at the equilibrium a finite stress created by a dislocation. In this model, u^0 is the phase transition. For instance, for an edge dislocation with Burgers vector e_1 as presented in Section 2, u^0 is a transition between the value 0 on the left to the value 1 on the right (see Figure 3.1). In the special case $u_t^0 = 0 = \sigma$ and for sinusoidal potentials W , the stationary solution u^0 is known explicitly (see for instance [7]), which makes the PN model very attractive. Let us mention that a physical and numerical study of the evolution problem (3.1) has been treated in [25].

Remark 3.1 *Remark that system (3.1)-(3.2) and also system (2.8) are equations which have the nice property to enjoy a comparison principle (see for instance related works [15, 7]), which can be made rigorous in the framework of viscosity solutions. Remark also that when we consider system (3.1)-(3.2) in this framework, the evolution equation on the boundary $\partial\Omega$ appears to be a boundary condition of the system. For this reason, as it is usual for viscosity solutions (see for instance [3, 4]), this boundary condition has to be understood technically in the sense that on $\partial\Omega$ the function u^0 solves pointwisely either $0 = \Delta u^0$ or the evolution equation.*

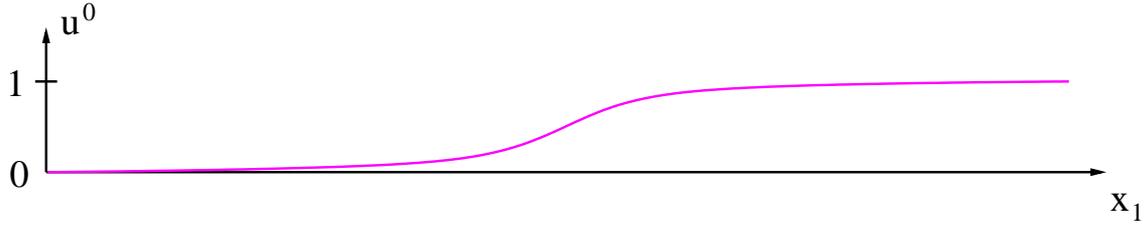


Fig. 3.1. Phase transition for an edge dislocation with Burgers vector e_1 for $X_2 = 0$

3.2 Convergence of the generalised FK model to the PN model

We have the following result

Theorem 3.2 (Formal convergence of FK to PN, $\varepsilon = \varepsilon_1 \rightarrow 0$)

Let $\varepsilon > 0$. For the initial data $u_0 \in W^{2,\infty}(\overline{\Omega})$ which is assumed harmonic on Ω , and under assumption (A1) on $\varepsilon W^\varepsilon$, and (A2) on σ , there exists a unique viscosity solution u^ε of system (2.8)-(2.9). Moreover assuming (2.10) with the potential W satisfying assumption (A1), then, as ε goes to zero, the solution u^ε formally converges to a viscosity solution of system (3.1)-(3.2).

The proof of Theorem 3.2 is done in full details in [13].

Sketch of the proof of Theorem 3.2

One way to guess the limit model (3.1)-(3.2) is to pass to the limit formally in system (2.8)-(2.9) assuming that the solution u^ε (and its derivatives) converges to a limit u^0 . The convergence in the system is then obtained using a simple Taylor expansion. The existence of a solution u^ε to system (2.8)-(2.9) is technically delicate and is based on the proof of a suitable comparison principle for this system.

3.3 Reformulation of the PN model

In this subsection and in Subsection 3.4, in order to simplify the notation we denote by $\varepsilon := \varepsilon_2 > 0$ the small parameter in the second passage of the scheme (1.1). We recall that it is well known that for bounded smooth functions u^0 defined on $\overline{\Omega}$ which are harmonic on Ω , we can write

$$\frac{\partial u^0}{\partial X_2}(X_1, 0) = L(u^0(\cdot, 0))(X_1) \quad \text{for all } (X_1, 0) \in \partial\Omega$$

where for a general bounded smooth function w , the linear operator L is given by the Levy-Khintchine formula (see Theorem 1 in [11]):

$$(Lw)(x) = \frac{1}{\pi} \int_{\mathbb{R}} \frac{dz}{z^2} (w(x+z) - w(x) - zw'(x)1_{\{|z| \leq 1\}}). \quad (3.3)$$

Then for smooth solutions u^0 , system (3.1) can be rewritten for $V(X_1, t) = u^0(X_1, X_2, t)|_{X_2=0}$ with $x = X_1 \in \mathbb{R}$ as

$$V_t = 2\varepsilon\sigma(\varepsilon x) - W'(V) + LV \quad \text{on } \mathbb{R}. \quad (3.4)$$

We also recall (see [7]) that there exists a unique function ϕ solution of

$$\begin{cases} 0 = L\phi - W'(\phi) & \text{on } \mathbb{R} \\ \phi' > 0 & \text{and } \phi(-\infty) = 0, \quad \phi(0) = \frac{1}{2}, \quad \phi(+\infty) = 1. \end{cases} \quad (3.5)$$

The function ϕ is called the layer solution and a translation of ϕ is pictured on Figure 3.1.

3.4 Rescaling of the PN model

We now consider the following rescaling

$$v^\varepsilon(x, t) = V\left(\frac{x}{\varepsilon}, \frac{t}{\varepsilon^2}\right).$$

Then system (3.4) can be rewritten as

$$v_t^\varepsilon = \frac{1}{\varepsilon} \left\{ Lv^\varepsilon - \frac{1}{\varepsilon} W'(v^\varepsilon) + 2\sigma(x) \right\} \quad \text{on } \mathbb{R} \quad (3.6)$$

with initial condition

$$v^\varepsilon(x, 0) = v_0^\varepsilon(x) \quad \text{for } x \in \mathbb{R}. \quad (3.7)$$

Again, a good notion of solution for system (3.6)-(3.7) is the notion of viscosity solution for non local equations (see for instance [5]).

Here we will choose carefully the initial condition v_0^ε as follows

Assumption (A3)

$$\begin{cases} x_1^0 < x_2^0 < \dots < x_N^0 \\ v_0^\varepsilon(x) = \frac{\varepsilon}{\alpha} \cdot 2\sigma(x) + \sum_{i=1}^N \phi\left(\frac{x - x_i^0}{\varepsilon}\right) \end{cases}$$

where we recall that $\alpha = W''(0) > 0$ and ϕ is defined in (3.5).

4 Dynamics of discrete dislocations

4.1 Description of the DDD model

In this section we assume that the phase transition reduces to a sharp interface where the transition is localized at the position $x = x_1^0 \in \mathbb{R}$. For a dislocation associated to a Burgers vector e_1 , the sharp interface is associated to a non-decreasing step function like the one of Figure 4.1.

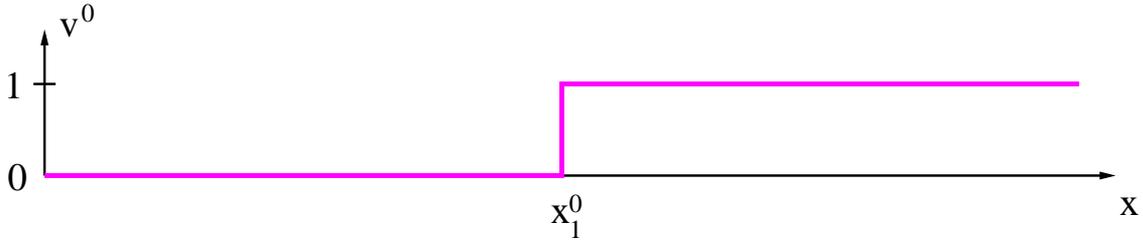


Fig. 4.1. Sharp interface describing a discrete dislocation at $x = x_1^0$

More generally, we can consider the case of N dislocations (or particules) of positions $(x_i(t))_{i=1, \dots, N}$ solving the following system

$$\frac{dx_i}{dt} = -\gamma \left(\sigma(x_i) + \sum_{j \neq i} V'(x_i - x_j) \right) \quad \text{on } (0, +\infty) \quad \text{for } i = 1, \dots, N \quad (4.1)$$

with the two-body interaction potential

$$V(x) = -\frac{1}{2\pi} \ln |x|$$

with initial data

$$x_i(0) = x_i^0 \quad \text{for } i = 1, \dots, N. \quad (4.2)$$

Here the constant $\gamma > 0$ is the inverse of the damping factor. It is related to the layer solution ϕ defined in (3.5) and is given by

$$\gamma = 2 \left(\int_{\mathbb{R}} (\phi')^2 \right)^{-1}.$$

The function σ is the applied shear stress and $V'(x - x_j)$ is the (singular) shear stress created at the point x by the dislocation x_j . This corresponds exactly to the shear stress already computed in (2.7). The total stress $\sigma(x_i) + \sum_{j \neq i} V'(x_i - x_j)$ is called the resolved Peach-Koehler force acting on the dislocation x_i .

4.2 Convergence of the PN model to the DDD model

We have

Theorem 4.1 (Convergence of PN to DDD, $\varepsilon = \varepsilon_2 \rightarrow 0$)

Let $\varepsilon > 0$. Under assumptions (A1)-(A2)-(A3), there exists a unique viscosity solution v^ε of (3.6)-(3.7). Moreover there exists a unique solution of (4.1)-(4.2), and we define

$$v^0(x, t) = \sum_{i=1, \dots, N} H(x - x_i(t))$$

where H is the Heavyside function. Then as ε goes to zero, the function v^ε converges to v^0 in the following sense

$$\limsup_{(x', t') \rightarrow (x, t), \varepsilon \rightarrow 0} v^\varepsilon(x', t') \leq (v^0)^*(x, t)$$

and

$$\liminf_{(x', t') \rightarrow (x, t), \varepsilon \rightarrow 0} v^\varepsilon(x', t') \geq (v^0)_*(x, t).$$

The proof of this result is done in full details in [17].

Remark 4.2 We recall that the semi-continuous envelopes of a function v are defined as

$$v^*(x, t) = \limsup_{(x', t') \rightarrow (x, t)} v(x', t') \quad \text{and} \quad v_*(x, t) = \liminf_{(x', t') \rightarrow (x, t)} v(x', t').$$

Sketch of the proof of convergence

The existence of a solution for all time of the ODE system (4.1)-(4.2) comes from the fact that $V(x)$ is a convex potential outside the origin. This property allows to show that the minimal distance between particles

$$d(t) = \inf_{i \neq j} |x_i(t) - x_j(t)|$$

satisfies

$$d(t) \geq d(0)e^{-C\gamma t} \tag{4.3}$$

which prevents the meeting of the particles at any finite time.

Then the main idea to prove the convergence is to approximate the solution v^ε by the following ansatz

$$\tilde{v}^\varepsilon(x, t) = \frac{\varepsilon}{\alpha} \cdot 2\sigma(x) + \sum_{i=1}^N \left\{ \phi \left(\frac{x - x_i}{\varepsilon} \right) - \varepsilon \dot{x}_i(t) \psi \left(\frac{x - x_i}{\varepsilon} \right) \right\} \quad \text{with} \quad \dot{x}_i(t) = \frac{dx_i}{dt}(t)$$

where $\alpha = W''(0)$ and the corrector ψ solves the following equation

$$L\psi - W''(\phi)\psi = \phi' + \eta(W''(\phi) - W''(0)) \quad \text{with} \quad \eta = \frac{1}{W''(0)} \int_{\mathbb{R}} (\phi')^2.$$

The stress created in x by a dislocation positioned at the origin, comes from the following property

$$\phi(x) - H(x) \sim -\frac{1}{\alpha\pi x} \quad \text{as} \quad |x| \rightarrow +\infty.$$

The rest of the proof of convergence of v^ε is done by construction of sub and super solutions based on the ansatz \tilde{v}^ε .

4.3 Rescaling of the DDD model

In this subsection, in order to simplify the notation we denote by $\varepsilon := \varepsilon_3 > 0$ the small parameter in the third passage of the scheme (1.1). We consider a given initial data w_0 which satisfies

Assumption (A4)

$$\begin{cases} w_0 \in W^{2,\infty}(\mathbb{R}), \\ w_0' > 0, \quad w_0(-\infty) = 0. \end{cases}$$

We also introduce the integer N_ε and the position of the dislocations $x_1^0 < \dots < x_{N_\varepsilon}^0$ such that

$$\sum_{i=1,\dots,N_\varepsilon} H(x - x_i^0) = \lfloor \frac{w_0(\varepsilon x)}{\varepsilon} \rfloor \quad (4.4)$$

where $\lfloor \cdot \rfloor$ denotes the floor function which is the following integer part:

$$\lfloor x \rfloor = \max\{n \in \mathbb{Z}; n \leq x\}.$$

We also assume that the stress σ is periodic. Precisely, we make the following assumption:

Assumption (A2')

$$\sigma \in C^2(\mathbb{R}) \quad \text{and} \quad \sigma(x+1) = \sigma(x) \quad \text{for all } x \in \mathbb{R}.$$

This assumption allows to study the collective behaviour of dislocations in a landscape with periodic obstacles, and to get the effective macroscopic model by a periodic homogenization approach when $\varepsilon \rightarrow 0$.

Then we consider the solution $(x_i(t))_{i=1,\dots,N_\varepsilon}$ of the system (4.1)-(4.2) with $N = N_\varepsilon$ where N_ε is defined by (4.4). In particular, remark that

$$N_\varepsilon \sim \frac{\sup_{\mathbb{R}} w_0}{\varepsilon} \rightarrow +\infty \quad \text{as } \varepsilon \rightarrow 0.$$

We define the function

$$v^0(x, t) = \sum_{i=1,\dots,N_\varepsilon} H(x - x_i(t))$$

and the rescaling

$$w^\varepsilon(x, t) = \varepsilon v^0\left(\frac{x}{\varepsilon}, \frac{t}{\varepsilon}\right). \quad (4.5)$$

5 Dislocation density model

5.1 Description of the DD model

We first introduce a function $g : (0, +\infty) \times \mathbb{R} \rightarrow \mathbb{R}$ which satisfies

Assumption (A5)

$$\begin{cases} g \in C^0((0, +\infty) \times \mathbb{R}), \\ l \mapsto g(\rho, l) \quad \text{is nondecreasing.} \end{cases}$$

Then we consider a function $w^0(x, t)$ which is a solution of

$$w_t^0 = g(w_x^0, Lw^0) \quad \text{on } \mathbb{R} \times (0, +\infty) \quad (5.1)$$

where the operator L is defined in (3.3), and with initial data

$$w^0(x, 0) = w_0(x) \quad \text{for all } x \in \mathbb{R}. \quad (5.2)$$

Here the function w^0 is such that its derivative w_x^0 represents the macroscopic dislocation density. Moreover w^0 can be seen as the plastic strain localized in plane $x_2 = 0$ and $\frac{1}{2}Lw^0$ can be identified to the stress created by the dislocation density w_x^0 . Equation (5.1) can be interpreted as the plastic flow rule in a model for macroscopic crystal plasticity. Indeed, from a mechanical point of view, we have the following table (see also [21]) of equivalence between our homogenized model and a classical model in mechanics for elasto-visco-plasticity of crystals (see [16]).

	Crystal elasto-visco-plasticity	DD model
resolved plastic strain	$\gamma(x_1)\delta_0(x_2)$	$w^0(x_1)\delta_0(x_2)$
Nye tensor of dislocations densities	$\alpha = (e_1 \otimes e_2)\gamma'(x_1)\delta_0(x_2)$	$\alpha = (e_1 \otimes e_2)w_x^0(x_1)\delta_0(x_2)$
exterior applied stress	Σ^{ext}	
microscopic resolved shear stress		$\sigma - \int_{(0,1)} \sigma$
resolved exterior applied stress	$\Sigma^{ext} : e^0$	$\int_{(0,1)} \sigma$
displacement	$v = v_1 e_1$	
strain	$e := e(v) - e^0 \gamma \delta_0(x_2)$ with $e(v) := \frac{1}{2}(\nabla v + {}^t \nabla v)$ and $e^0 := \frac{1}{2}(e_2 \otimes e_1 + e_1 \otimes e_2)$	
total elastic energy	$E := \int_{\mathbb{R}^2} \frac{1}{2}(\Lambda : e) : e + \Sigma^{ext} : e$	$E := \int_{\mathbb{R}} -\frac{1}{2}w^0 Lw^0 - \left(\int_{(0,1)} \sigma \right) w^0$
macroscopic stress	$\Sigma := \Lambda : e + \Sigma^{ext}$	
resolved macroscopic shear stress	$\tau := \Sigma : e^0$	$\tau := Lw^0 + \int_{(0,1)} \sigma$
visco-plastic law	$\frac{\partial \gamma}{\partial t} = f(\tau)$	$\frac{\partial w^0}{\partial t} = g\left(w_x^0, Lw^0 + \int_{(0,1)} \sigma\right)$
energy decay	$\frac{d}{dt}E = \int_{\mathbb{R}} -\tau f(\tau) \leq 0$	$\frac{d}{dt}E = \int_{\mathbb{R}} -\tau g(w_x^0, \tau) \leq 0$

Remark that when we choose the microscopic stress σ in Assumption (A2') so that $\int_{(0,1)} \sigma = 0$, we expect that g satisfies

$$g(\rho, 0) = 0. \quad (5.3)$$

This equality reflects the pinning of dislocations (see [14, Th 2.6]). In the model presented in the previous table, the plastic strain velocity w_t^0 is prescribed by the function g (whose typical profile can be seen on Figure 5.1) which is assumed to satisfy (5.3).

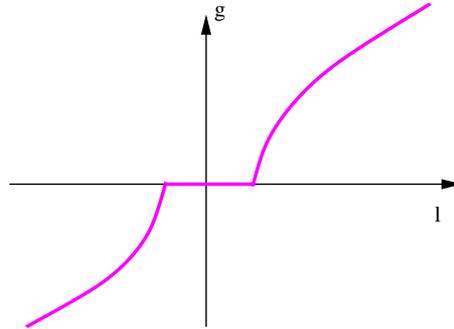


Fig. 5.1. Graph of the map $l \mapsto g(\rho, l)$ showing a threshold effect

Let us mention that a similar profile for g as a function of l only have been obtained rigorously for a different model involving the motion of a phase transition solution of Allen-Cahn type equations (see [10]).

5.2 Convergence of the DDD model to the DD model

We have the following result

Theorem 5.1 (Convergence of DDD to DD, $\varepsilon = \varepsilon_3 \rightarrow 0$)

Let us assume (A2'), (A4). Then there exists a function g satisfying assumption (A5). Moreover the function w^ε defined in (4.5) converges to the unique solution w^0 of (5.1)-(5.2), locally uniformly on $\mathbb{R} \times [0, +\infty)$.

The proof of this result is done in full details in [14].

Remark that Theorem 5.1 is an homogenization result in the periodic setting. In the particular case where the periodic stress σ is equal to zero, we get

$$g(\rho, l) = \frac{\gamma}{2}\rho l.$$

The presence of a non-zero 1-periodic stress with zero mean value, creates a threshold phenomenon where for a fixed dislocation density ρ , the quantity $g(\rho, l)$ can be equal to zero if $|l|$ is small enough (see for instance the numerical simulations in [12] which look like Figure 5.1).

Heuristic ideas used in the proof of Theorem 5.1

Step 1 : Splitting of the interactions

Let us first approximate the positions of the particles x_j as follows (dropping the time dependence of w^0)

$$w^0(\varepsilon x_j) = j\varepsilon. \quad (5.4)$$

For some radius R_ε satisfying:

$$\frac{1}{\varepsilon} \gg R_\varepsilon \rightarrow 0 \quad \text{as } \varepsilon \rightarrow 0,$$

we split the interactions into two contributions (short and long range interactions):

$$\sum_{j \neq i} V'(x_i - x_j) = \underbrace{\sum_{|j-i| \leq R_\varepsilon, j \neq i} V'(x_i - x_j)}_{I^\varepsilon} + \underbrace{\sum_{|j-i| > R_\varepsilon} V'(x_i - x_j)}_{II^\varepsilon}.$$

Step 2 : The long range interactions

Remark that we can rewrite (using the homogeneity of V'):

$$II^\varepsilon = \sum_{|j\varepsilon - i\varepsilon| > \varepsilon R_\varepsilon} \varepsilon V'(\varepsilon x_i - \varepsilon x_j).$$

Setting $z = z_j = \varepsilon x_j$, $dz = z_{j+1} - z_j$, $x = \varepsilon x_i$, we see for a finite difference of (5.4) that $\varepsilon \simeq (w^0)'(z)dz$. Noticing that $\varepsilon R_\varepsilon \rightarrow 0$, we see formally that we can approximate II^ε by the following integral

$$II^\varepsilon \simeq \int_{\mathbb{R}} (w^0)'(z) V'(x - z) dz = (w^0 * V'')(x) = \frac{1}{2} L w^0(x) =: \frac{l}{2}.$$

This shows in particular that, because of the slow decay of the interactions V' , the long range contribution II^ε is non zero in general.

Step 3 : The short range interactions and determination of the function g

Assume that for $|j - i| < R_\varepsilon$ and $|t - t_0| \ll \frac{1}{\varepsilon}$, the following quantity is bounded:

$$|x_j(t) - x_i(t_0) - v(t - t_0) - (j - i)/\rho| \leq C$$

where v is the mean velocity of the particles x_j , and $\rho = (w^0)'(x_i(t_0))$ is the mean density of particles. A way to guess the correct velocity v is to consider equation (4.1) with σ replaced by $\sigma + \frac{l}{2}$ (because of the long range contribution $II^\varepsilon \simeq \frac{l}{2}$), and to look for particular solutions:

$$x_j(t) = h(vt + j/\rho) \quad \text{with} \quad h(a + 1) = 1 + h(a) \quad \text{for all } a \in \mathbb{R}.$$

The function h is called the hull function. Both h and the constant v have to be determined. It can be shown that v is unique. Then we set

$$g(\rho, l) = -v\rho$$

which is known in physics as the Orowan's law.

6 Conclusion

We considered a two-dimensional Frenkel-Kontorova model in the fully overdamped regime. From this model, we derived by a scaling argument the time-dependent Peierls-Nabarro model. Looking at the sharp interface limit of the phase transitions of the Peierls-Nabarro model, we were able to identify a dynamics of particles that corresponds to the classical discrete dislocation dynamics, in the particular case of parallel straight edge dislocation lines in the same glide plane with the same Burgers vector. Considering the motion of these particles in a landscape with periodic obstacles, we were able to identify at large scale an evolution model for the dynamics of a density of dislocations. This model is a macroscopic model for crystal visco-elasto-plasticity, where we predicted a plastic flow rule. This last model shows in particular a threshold effect where dislocations can be pinned in the obstacles, if the effective stress acting on these dislocations is too small.

In order to present a summary of our approach, we give here a diagram (see Figure 6.1) that shows the links between the four models treated in this paper. Up to our knowledge, this derivation of classical models from a single microscopic model (the 2D Frenkel-Kontorova model), seems new and allows to make clear connections between different modelling of dislocation dynamics.

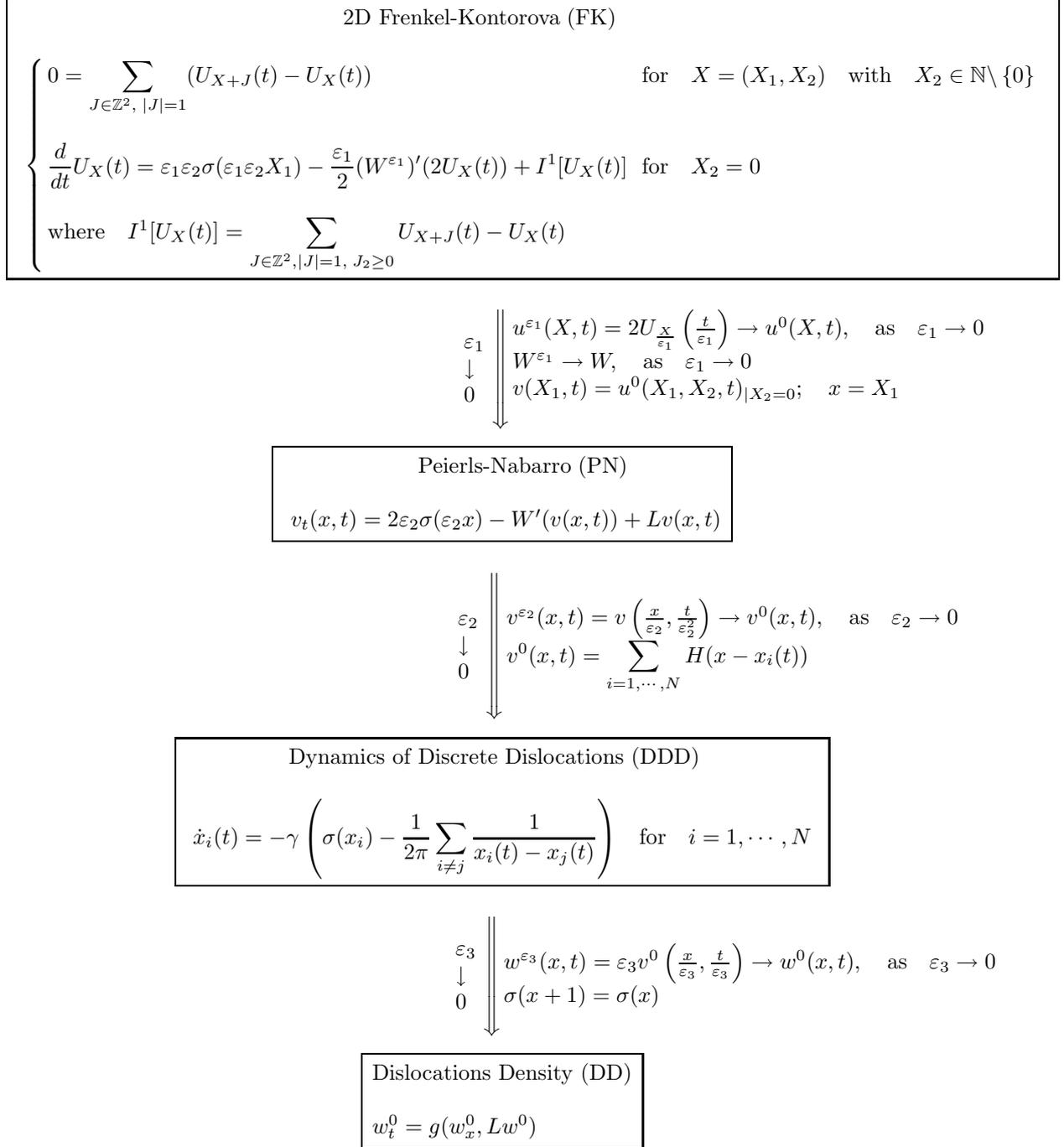


Fig. 6.1. Descriptive diagram summarizing the links between the four models

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References

1. O. Alvarez, P. Hoch, Y. Le Bouar, and R. Monneau. *Dislocation dynamics: short-time existence and uniqueness of the solution*. Arch. Ration. Mech. Anal., 181 (3), (2006), 449–504.
2. V.I. Alshits, V.L. Indenbom, *Mechanisms of dislocation drag*. In: Nabarro, F.R.N. (Ed.), Dislocations in Solids. Elsevier, Amsterdam, (1986), 43-111.
3. G. Barles, *Solutions de viscosité des équations de Hamilton-Jacobi*, vol. 17 of Mathématiques & Applications (Berlin) [Mathematics & Applications], Springer-Verlag, Paris, (1994).
4. G. Barles, F. Da Lio, *On the Boundary Ergodic Problem for Fully Nonlinear Equations in Bounded Domains with General Nonlinear Neumann Boundary Conditions*, Annales de l'Institut Henri Poincaré Analyse non linéaire 22 (5), (2005), 521-541.
5. G. Barles and C. Imbert. *Second-order elliptic integro-differential equations: viscosity solutions' theory revisited*. Annales de l'IHP 25 (3), (2008), 567-585.
6. O.M. Braun, Y.S. Kivshar, *The Frenkel-Kontorova Model, Concepts, Methods and Applications*, Springer-Verlag, (2004).
7. X. Cabré and J. Solà-Morales. *Layer solutions in a half-space for boundary reactions*. Comm. Pure Appl. Math. 58 (12), (2005), 1678-1732.
8. A. Carpio, L.L. Bonilla, *Edge dislocations in crystal structures considered as travelling waves of discrete models*, Phys. Rev. Lett. 90 (13), (2003), 135502, 1-4; 91 (2), (2003), 029901-1.
9. A. Carpio, S. J. Chapman, S. D. Howison and J. R. Ockendon, *Dynamics of line singularities*, Phil. Trans. R. Soc. Lond. A 355, (1997), p. 2013.
10. N. Dirr, N. K. Yip, *Pinning and de-pinning phenomena in front propagation in heterogeneous media*, Interfaces and Free Boundaries 8 (2006), 79-109.
11. J. Droniou, C. Imbert, *Fractal first order partial differential equations*, Arch. Rat. Mech. Anal. 182 (2006), 299-331.
12. A. El Hajj, H. Ibrahim, R. Monneau, *Homogenization of dislocation dynamics*, preprint 2008.
13. A. Fino, H. Ibrahim, R. Monneau, work in progress.
14. N. Forcadel, C. Imbert, R. Monneau, *Homogenization of the dislocation dynamics and of some particle systems with two-body interactions*, Discrete and Continuous Dynamical Systems - A, vol. 23 (3) (March 2009), 785-826.
15. N. Forcadel, C. Imbert, R. Monneau, *Homogenization of fully overdamped Frenkel-Kontorova models*, Journal of Differential Equations 246 (3) (1 February 2009), 1057-1097.
16. D. François, A. Pineau, A. Zaoui, *Comportement mécanique des matériaux;lasticité et plasticité*, Herms, Paris, (1995).
17. M. González, R. Monneau, work in progress.
18. S. Haq, A. B. Movchan, G. J. Rodin, *Analysis of lattices with non-linear interphases*, Acta Mechanica Sinica 22 (2006), 323-330.
19. S. Haq, A. B. Movchan, G. J. Rodin, *Lattice Green's functions in nonlinear analysis of defects*, Journal of applied mechanics - Transactions of the ASME 74 (4) (July 2007), 686-690.
20. J.P. Hirth, J. Lothe, *Theory of Dislocations*, McGraw-Hill, New York, (1972).
21. C. Imbert, R. Monneau, E. Rouy, *Homogenization of first order equations with (u/ϵ) -periodic Hamiltonians. II. Application to dislocations dynamics*, Comm. Partial Differential Equations 33 (2008), 479-516.
22. O. Kresse, L. Truskinovsky, *Mobility of lattice defects: discrete and continuum approaches*, Journal of the Mechanics and Physics of Solids 51 (2003), 1305-1332.
23. O. Kresse, L. Truskinovsky, *Lattice friction for crystalline defects: from dislocations to cracks*, Journal of the Mechanics and Physics of Solids 52 (2004), 2521-2543.
24. O. Kresse, L. Truskinovsky, *Prototypical lattice model of a moving defect: The role of environmental viscosity*, Izvestiya, Physics of the Solid Earth, vol. 43, no. 1 (2007) 63-66.
25. A. B. Movchan, R. Bullough, J. R. Willis, *Stability of a dislocation: Discrete model*, European Journal of Applied Mathematics, vol. 9 (1998), 373-396.