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Statistical approach to dislocation dynamics: From dislocation correlations to a multiple-slip continuum theory of plasticity

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I. INTRODUCTION

Statistical mechanics provides an optimal framework and various tools for studying emergent phenomena from a complex conglomerate of bodies—may they be molecules of gases, polymer chains of rubber, or crystalline defects. The use of correlation functions in analyzing two-dimensional solids and their defects has been proven very successful in the past. For example, Mermin\(^5\) showed that the two-dimensional crystals do not have conventional long-range order, but can have “directional long-range order.” Nelson and co-worker\(^6\) applied the technique to explain dislocation-assisted melting in two dimensions. Over a decade ago, Groma\(^4\) proposed a theory to describe dislocations and their motions using distribution functions and probability arguments. Unlike the existing continuum theories at the time,\(^6\) the formalism was physically motivated and it correctly incorporated the long-range nature of dislocation interactions. Several variations of this work—all of which reduce to the same two-dimensional theory—also exist for three-dimensional dislocation systems.\(^6\)–\(^9\)

Although having laid out the foundation for possible interactions of many-dislocation configurations, Groma’s pioneering work did not investigate these correlated effects in detail. Zaiser et al.\(^10\) explicitly considered the evolution of dislocation correlations by extending Groma’s theory for systems of single-slip, parallel edge dislocations. They were able to qualitatively obtain the correct scaling behavior of the evolution equations for both single and pair correlation densities and explained some general properties of these functions. Their formulation, however, was limited to only one active slip system and the analytical forms of pair correlation functions were not derived. In a later work, Groma et al.\(^11\) included the influence of dislocation correlations in the form of a local back stress. Yefimov et al.\(^12,13\) connected this statistical description to a continuum crystal plasticity theory and applied this to various boundary value problems. While the theory successfully captured most features observed in discrete dislocation simulations, its ad hoc extension to multiple-slip systems failed to explain the size effects in single-crystal thin films.\(^14\) The main goals of this paper are (1) to correctly describe and obtain analytical expressions for dislocation pair correlations and (2) to systematically generalize the approach of Groma et al. to multiple-slip systems.

We begin in Sec. II by introducing ensembles of dislocations and deriving the partition function for multiple-slip systems. The n-th order dislocation densities and dislocation correlation functions are subsequently defined. We construct the Bogolyubov–Born–Green–Yvon–Kirkwood (BBGYK) integral equations in Sec. III. These equations link correlation functions of order n to those of order n + 1 (a technique generally used in the study of dense gases and fluids). The integral equations are expanded in powers of interaction strength (the ratio between the interaction energy and “thermal” energy). We then obtain a set of approximate integral equations for pair (n=2) correlation functions after applying a closure approximation to truncate the series. These equations are valid regardless of the form of the interaction potential, and thus are applicable to other systems, provided that this pair interaction vanishes at a large distance.

By appealing to the Peach–Koehler interaction, analytical expressions for pair dislocation densities for single- and multiple-slip systems are derived in Secs. IV and V, respectively. Our single-slip solution agrees with the result from the study of induced geometrically necessary dislocations (GNDs) in terms of a single pinned dislocation by means of a variational approach.\(^15,16\) The dislocation spacing \(1/\sqrt{\rho}\) emerges as a natural length scale in this formulation in accordance with the scaling study by Zaiser et al.\(^10\) Our analysis further shows long-range attractive correlations when more than one slip system are present, confirming the absence of dislocation patterning in single-glide systems as ob-
serve in many discrete dislocation simulations\textsuperscript{17–24} and explained in a recent three-dimensional continuum plasticity theory.\textsuperscript{7,8}

In Sec. VI, we write down the transport equations for both total dislocation densities and GND densities on each slip system under the influence of the Peach–Koehler forces from both single and pair dislocation correlations. While the former gives a self-consistent, long-range internal stress contribution, the latter exerts an additional short-range, entropic force due to a deviation away from a preferred dislocation arrangement in the form of a back stress. The formulation is a direct extension of the work by Groma co-workers\textsuperscript{10,11} for crystals with one active slip system. Using the knowledge of the pair correlation functions, we obtain a complete description of the back stress as a function of slip orientations—something that has not been achieved in the Groma–Zaiser approach but which had been previously incorporated in the multiple-slip theory using the \textit{ad hoc} phenomenological considerations.\textsuperscript{14,25}

Finally in Sec. VII, we contrast our theory with the multiple-slip theory of Yefimov and Van der Giessen.\textsuperscript{14,25} While both theories propose that interactions among slip systems solely depend on relative angles of slip orientations, the functional forms are different. We attribute the failure of the earlier theory in explaining size effects in single-crystal thin films partly to this difference and partly to the treatment of dislocation nucleation in the theory.

\section{II. DEFINITIONS OF THE BASIC QUANTITIES}

Let us first consider the equation of motion of a dislocation, which drifts along its glide plane (taken for simplicity to be along the \( \hat{x} \) direction) subject to the Peach–Koehler force \( f(x,y) = -\partial_x u(x,y) \) generated by another dislocation at the origin and to random fluctuations in the form of the Gaussian white noise\textsuperscript{26} \( \xi(t) \) (also along \( \hat{x} \)) where \( \langle \xi(t_1) \xi(t_2) \rangle = \delta(t_1 - t_2) \).

\begin{equation}
\vec{\dot{r}}(x,y,t) = \xi(f(x,y) + \sqrt{2T} \vec{\xi}(t)).
\end{equation}

The form of the fluctuation amplitude \( \sqrt{2T} \) is given afterward by the fluctuation-dissipation theorem. Thus, \( T \) in this theory is not physical temperature but a measure of the amplitude of random fluctuations. The stationary state solution of the corresponding one-dimensional Smoluchowski equation,

\begin{equation}
\frac{\partial P_1}{\partial t} = \frac{\partial}{\partial x} \left[ f(x) P_1 - \frac{\partial}{\partial x} P_1 \right],
\end{equation}

which describes that the probability \( P_1(\vec{r}) \) of finding a dislocation at \( \vec{r} \) can be described by the Boltzmann distribution \( P_1 \sim e^{-(u/2T)} \).

To generalize the above concept to a many-dislocation system, consider an arrangement of \( r \) species of dislocations and denote the coordinate of the \( i \)th dislocation of species \( s \) by \( \vec{r}_i \). The dislocation configuration \( \{N\} \) is the set of the coordinates of all dislocations, where \( \mathbf{N} = (N_1, N_2, N_3, N_4; \ldots; N_{r-1}, N_r) \) denotes the “collection” of dislocations of type \( s \). In this convention, odd and even slots, respectively, contain plus and minus dislocations on distinct slip systems.\textsuperscript{27} We introduce the notation \( \{N + 1, r \} \) to denote the addition of an extra dislocation of species \( s \) to \( \{N\} \), while similarly, a configuration \( \{N\} \) with coordinates of \( n \) removed is indicated by \( \{N - n\} \).

The interacting Hamiltonian \( U_N \) of the system can be written as the sum of potentials \( u(t_1, t_2) \) of all pairs of dislocations,

\begin{equation}
U_N(\{N\}) = \sum_{t_1 \neq t_2} \sum_{i \not\in j} u(\vec{r}_i, \vec{r}_j).
\end{equation}

We can define a canonical partition function of configuration \( \mathbf{N} \) by

\begin{equation}
Z_N = \int e^{-U_N(\{N\})} dN,
\end{equation}

where the integrations are taken over the “volume” measure \( dN = \prod_{i=1}^{r} d^2 \vec{r}_i d^2 \vec{R}_i \) of the dislocation configuration at \( \{N\} \). Consider the coordinates of a particular set \( \{n\} \), the probability of observing the configuration \( \mathbf{n} \) in \( d(\mathbf{n}) \) about the points in \( \{\mathbf{n}\} \) irrespective of the remaining collection \( \mathbf{N} - \mathbf{n} \) is

\begin{equation}
P_N(\mathbf{n}) d(\mathbf{n}) = \frac{d(\mathbf{n})}{2^N} \int e^{-U_N(\{N\})} d\{N - \mathbf{n}\},
\end{equation}

where \( \int P_N(\mathbf{n}) d(\mathbf{n}) = 1 \). The probability density of observing any statistically equivalent possible collection \( \mathbf{n} \) within the volumes \( d(\mathbf{n}) \) about the points \( \{\mathbf{n}\} \) is therefore

\begin{equation}
\rho(\mathbf{n}) = \prod_{i=1}^{r} \frac{N_i!}{(N_i - n_i)!} P_N(\mathbf{n}).
\end{equation}

Consider now an \textit{open} system (which could be realized, say, by allowing for nucleation and annihilation of dislocations as the system relaxes); a grand canonical partition function is given by

\begin{equation}
\Xi = \sum_{N=0}^{r} \prod_{i=1}^{r} \frac{N_i!}{N_i^{N_i}} Z_N.
\end{equation}

where \( z_s \) is the activity of species \( s \). The prefactor arises from integrating away the momentum degrees of freedom in the Hamiltonian, which are irrelevant to this problem. The probability \( \mathcal{P} \) of the occurrence of configuration \( \mathbf{N} \) in the open system is therefore

\begin{equation}
\mathcal{P}(\mathbf{n}) = \prod_{i=1}^{r} \frac{N_i!}{N_i^{N_i}} \frac{Z_N}{\Xi}.
\end{equation}

Finally, the probability density of observing any \( n_i \) dislocations of species 1, \( n_2 \) dislocations of species 2, etc., (any collection \( \mathbf{n} \) in \( d(\mathbf{n}) \) at \( \{\mathbf{n}\} \) is

\begin{equation}
\rho(\mathbf{n}) = \sum_{\mathbf{N}=\mathbf{n}} \mathcal{P}(\mathbf{n}) \Xi(\mathbf{n}).
\end{equation}

The summation is taken over all collections \( \mathbf{N} \) greater than or
equal to \( n \), i.e., for all \( N_i \geq n_i \), \( N_2 \geq n_2 \), etc. We take Eq. (9) as the definition of dislocation density of order \((n)\). Explicitly, we have

\[
\rho^{(n)}(\{n\}) = \frac{1}{\Xi} \sum_{N \geq n} \left[ \prod_{i=1}^{r} \frac{z_{s_i}^{N_i}}{(N_i - n_i)!} \right] e^{-\bar{U}_N(\{N\})/T} d\{N - n\}.
\]

(10)

This definition of an \((n)\)th-order dislocation density is equivalent to the ones used by Groma and Zaiser in the realization of an open system. Finally, we define the \((n)\)th-order correlation function \(g^{(n)}(\{n\})\) through

\[
\rho^{(n)}(\{n\}) = \prod_{i=1}^{r} \rho^{(1)}(\vec{1}_i) \rho^{(1)}(\vec{2}_i) \cdots \rho^{(1)}(\vec{n}_i) \] \[g^{(n)}(\{n\}).
\]

(11)


The Bogolyubov–Born–Green–Yvon–Kirkwood integral equations first appeared in the study of classical fluids with a total potential energy given by the sum of pair interactions. They provided a set of relations between distribution functions of fluid density at different orders. Here, we extend the BBGYK formalism to include the non-central interactions of dislocations in a multicomponent system. We proceed in three steps: (1) take a derivative of the \((n)\)th-order dislocation density with respect to the glide direction of one particle of the interested species; (2) express the result in terms of the next higher order densities; and (3) convert the integral equations of densities into those of correlation functions.

Differentiating \(\rho^{(n)}(\{n\})\) as expressed in Eq. (10) with respect to the glide direction of dislocation \(1 \) of species \(1 \) located, we find

\[
\nabla_i \rho^{(n)}(\{n\}) = -\frac{1}{\Xi} \sum_{N \geq n} \left[ \prod_{i=1}^{r} \frac{z_{s_i}^{N_i}}{(N_i - n_i)!} \right] \int e^{-\bar{U}_N(\{N\})/T} \nabla_i \bar{U}_N(\{N\}) d\{N - n\},
\]

(12)

where we absorb \(1/T\) into the definition \(\bar{U}_N = U_N/T\). The derivative of the potential can be separated into two parts:

\[
\nabla_i \bar{U}_N = \sum_{i=1}^{r} \sum_{(i) \neq (1)} \nabla_i \bar{u}(1, \vec{i}_1) + \sum_{i=n+1}^{N} \nabla_i \bar{u}(1, \vec{i}_s).
\]

(13)

Direct substitution of Eq. (13) into the integrand of Eq. (12) splits the expression into two integrals \(I_1\) and \(I_2\). Notice in the first integral that the derivative of the potential does not depend on coordinates \(\{N - n\}\) and thus can be taken out of the integral, which yields

\[
I_1 = -\frac{1}{\Xi} \sum_{i=1}^{r} \sum_{(i) \neq (1)} \nabla_i \bar{u}(1, \vec{i}_1) \int \sum_{i=n+1}^{N} \nabla_i \bar{u}(1, \vec{i}_s) e^{-\bar{U}_N(\{N\})} d\{N - n\}.
\]

(14)

with the aid of Eq. (10). The second integral \(I_2\) requires a little more work,

\[
I_2 = -\frac{1}{\Xi} \sum_{i=1}^{r} \sum_{(i) \neq (1)} \int e^{-\bar{U}_N(\{N\})} d\{N - n\}.
\]

(15)

The expression involves integrating \(i_s\) over the sample size. Since each integral over \(i_s\) between \(n_s+1 \leq i \leq N_s\) is equivalent in infinite space, the summation therefore gives a factor of \((N_s - n_s)\). The remaining integrals over all other dislocation coordinates are unaffected.

Equation (15) thus becomes

\[
I_2 = -\sum_{i=1}^{r} \int e^{-\bar{U}_N(\{N\})} d\{N - n\}.
\]

(16)

The symbol \(d\{N - n\}/(n+1)\) represents the volume measure of \(\{N - n\}\) without \(d^2(n+1)\). Collecting both \(I_1\) and \(I_2\) from Eq. (14) and Eq. (16), we arrive at the BBGYK equations for the \((n)\)th-order dislocation density.
\[ \nabla_1 \rho^{(n)}(\mathbf{n}) = - \rho^{(n)}(\mathbf{n}) \sum_{s=1}^{r(n)} \nabla_1 \tilde{u}(\mathbf{i}_s, \mathbf{i}_s) \times ((\mathbf{n} + 1)_s) d^2(\mathbf{n} + 1)_s. \]  

One can obtain a series of integrodifferential equations for the correlation functions \( g^{(n)} \) from Eq. (17) by expanding out \( \rho^{(n)}(\mathbf{n}) \) using Eq. (11). All but two of the single-dislocation densities on the left-hand side (LHS) and right-hand side (RHS) of the equality cancel, which results in

\[ \nabla_1 [\rho(\tilde{1})] g^{(n)}(\mathbf{n}) = - \rho(\tilde{1}) g^{(n)}(\mathbf{n}) \sum_{s=1}^{r(n)} \nabla_1 \tilde{u}(\mathbf{i}_s, \mathbf{i}_s) \times ((\mathbf{n} + 1)_s) d^2(\mathbf{n} + 1)_s. \]  

The first-order densities \( \rho(\tilde{1}) \) that plague the expression can be removed by first using the product rule to the LHS, then dividing both sides by \( \rho(\tilde{1}) \). The LHS becomes

\[ \text{LHS} = \nabla_1 g^{(n)}(\mathbf{n}) + g^{(n)}(\mathbf{n}) \nabla_1 \rho(\tilde{1}) \rho(\tilde{1}). \]

The ratio of the derivative of the first-order density with itself can be rewritten using Eq. (17) specialized to first order, which gives

\[ \frac{\nabla_1 \rho(\tilde{1})}{\rho(\tilde{1})} = - \sum_{s=1}^{r(n)} \nabla_1 \tilde{u}(\mathbf{i}_s, \mathbf{i}_s) \rho(\tilde{\zeta}_s) g^{(2)}(\mathbf{i}_s, \mathbf{i}_s) d^2 \tilde{\zeta}_s, \]

where \( \tilde{\zeta}_s = (\mathbf{n} + 1)_s \) is the position of the \((\mathbf{n} + 1)_s\)th dislocation of species \( s \) and \( g^{(2)}(\mathbf{i}_s, \mathbf{i}_s) \) represents the pair correlation function between the first dislocation of species \( 1 \) at \( \tilde{\mathbf{1}}_1 \) and the \((\mathbf{n} + 1)_s\)th dislocation of species \( s \) at \( \tilde{\mathbf{1}}_s \). This expression could be seamlessly incorporated into the right-hand side of Eq. (18). The final result is

\[ \nabla_1 g^{(n)}(\mathbf{n}) = - g^{(n)}(\mathbf{n}) \sum_{s=1}^{r(n)} \nabla_1 \tilde{u}(\mathbf{i}_s, \mathbf{i}_s) \times ((\mathbf{n} + 1)_s) d^2(\mathbf{n} + 1)_s \]

\[ \times g^{(+1)}(\mathbf{n} + 1)_s. \]  

For the remainder of this paper, we shall restrict our attention to the Peach–Koehler interaction. Recall that the interaction energy between two parallel edge dislocations of length \( L \) (over energy \( T \) due to random forces) in an infinite medium is

\[ \mathcal{U}(\tilde{\mathbf{i}}, \tilde{\mathbf{i}}') = - \Gamma \psi(\tilde{\mathbf{i}}, \tilde{\mathbf{i}}'), \]

where \( \Gamma = \frac{\mu b^2 L}{2\pi(1-\nu)} \), and

\[ \psi(\tilde{\mathbf{i}}, \tilde{\mathbf{i}}') = \left\{ \begin{array}{ll} (\mathbf{m}_i^\ast \cdot \mathbf{m}_{i'})(\ln(|\tilde{\mathbf{i}}_i - \tilde{\mathbf{i}}_i'|) \right. \\ \left. + [\mathbf{m}_i^\ast \cdot (\tilde{\mathbf{i}}_i - \tilde{\mathbf{i}}_{i'})][\mathbf{m}_{i'}^\ast \cdot (\tilde{\mathbf{i}}_i - \tilde{\mathbf{i}}_{i'})] \right\}. \]  

Here, \( \mathbf{m}_i^\ast \) denotes the slip-plane normal of species \( s \). The relative strength \( \Gamma \) represents the ratio between dislocation interaction energy versus energy from random fluctuation due to noise. Note that the latter originates from the use of the Boltzmann distribution in Eq. (4) to describe the equilibrium configuration of systems with noise. As the dislocation configuration becomes more and more correlated, \( \Gamma \) becomes smaller.

For an explicit dependence on \( \Gamma \) to use as an expansion coefficient, we rescale the distance by the square root of the relative strength, \( \sqrt{\Gamma} \tilde{r} \rightarrow \tilde{r} \). Equation (19) specialized to second order gives

\[ \nabla_1 g^{(2)}(\tilde{1}, \tilde{2}) = \Gamma g^{(2)}(\tilde{1}, \tilde{2}) \nabla_1 \psi(\tilde{1}, \tilde{2}) + \sum_{s=1}^{r} \nabla_1 \psi(\tilde{1}, \tilde{3}_s) \rho(\tilde{3}_s) \times [g^{(3)}(\tilde{1}, \tilde{2}, \tilde{3}_s) - g^{(2)}(\tilde{1}, \tilde{2}) g^{(2)}(\tilde{1}, \tilde{3}_s)] d^2 \tilde{3}_s. \]

Here, we have simplified the notation even further by suppressing all irrelevant subscripts: vectors \( \tilde{1}, \tilde{2} \) and \( \tilde{3} \) simply denote the positions of dislocations 1 and 2 with their corresponding species. The summation \( \Sigma_r \) is taken over all \( s \) species present in the system.

We proceed by assuming that the correlation functions have the following forms:

\[ g^{(2)}(\tilde{1}, \tilde{2}) = 1 + \Gamma f^{(2)}(\tilde{1}, \tilde{2}), \]

\[ g^{(3)}(\tilde{1}, \tilde{2}, \tilde{3}) = 1 + \Gamma [f^{(2)}(\tilde{1}, \tilde{2}) + f^{(2)}(\tilde{1}, \tilde{3}) + f^{(2)}(\tilde{2}, \tilde{3})] \]

\[ + \Gamma^2 f^{(3)}(\tilde{1}, \tilde{2}, \tilde{3}) \]

for any vectors \( \tilde{1}, \tilde{2} \), and \( \tilde{3} \). The functions \( f^{(2)}(\tilde{1}, \tilde{2}) \) and \( f^{(3)}(\tilde{1}, \tilde{2}, \tilde{3}) \) should asymptotically vanish along the boundaries of the sample or as \( |\tilde{1} - \tilde{2}|, |\tilde{1} - \tilde{3}|, |\tilde{2} - \tilde{3}| \rightarrow \infty \) for an infinite system. Note, in particular, that

\[ g^{(3)}(\tilde{1}, \tilde{2}, \tilde{3}) - g^{(2)}(\tilde{1}, \tilde{2}) g^{(2)}(\tilde{1}, \tilde{3}) = \Gamma f^{(2)}(\tilde{2}, \tilde{3}) \]

\[ + \Gamma^2 [f^{(3)}(\tilde{1}, \tilde{2}, \tilde{3}) - f^{(2)}(\tilde{1}, \tilde{2}) f^{(2)}(\tilde{1}, \tilde{3})]. \]

So far, no approximation has been made. Equation (22) governing the second-order correlations naturally involves the third-order correlations. To systematically close the chain at the second order, we substitute Eqs. (23) and (24) into Eq. (22) to produce a set of integrodifferential equations of \( f^{(2)} \) and \( f^{(3)} \) for each power of \( \Gamma \). This technique was introduced.
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by Bogolyubov\textsuperscript{37} in the study of correlations in the Coulomb interactions and has been widely used in both high-energy and condensed matter communities in renormalization group theory. It should be mentioned that, in principle, $\Gamma$ needs to be small for the expansion to be valid,\textsuperscript{38} but this is not true here. The values of $\Gamma$ in Secs. IV and V turn out to be greater than 1—as often happens in diagrammatic expansion calculations in high-energy physics. The validity of the approach can be judged by comparisons between the theory and numerical simulation results.

The equation of power $\Gamma^0$ gives an identity. After integrating away $\nabla_i$ because $f^{(2)}$ and $\psi$ vanish at infinity, the equation of power $\Gamma$ becomes

$$f^{(2)}(\vec{1}, \vec{2}) = \psi(\vec{1}, \vec{2}) + \sum_{s=1}^{r} \phi(\vec{1}, \vec{3}_s) \rho(\vec{3}_s) f^{(2)}(\vec{2}, \vec{3}_s) d^2 \vec{3}_s.$$  

(25)

This equation is the key result of the analysis. We shall use it to obtain dislocation pair correlation functions for systems with one in Sec. IV and many in Sec. V active slip systems.

IV. PAIR CORRELATION FUNCTIONS FOR SINGLE SLIP

To illustrate the use of Eq. (25), we first apply it to the case of one slip system containing two species of dislocations (denoted as + and −). According to Eq. (21) valid for an infinite sample, $\psi(1, \vec{2}) = \psi(-1, \vec{2}) = \psi(\vec{2} - 1)$ which implies that $f^{(2h)}(\vec{1}, \vec{2}) = f^{(2h)}(\vec{1}, \vec{-2}) = f^{(2h)}(2 \vec{-1})$. Without loss of generality, we can take the origin to be at 2 and thus, from Eq. (25), we obtain the following set of integral equations:

$$f^{++}(\vec{r}) = \psi_1(\vec{r}) + \int d^2 \vec{r}' \psi_1(\vec{r} - \vec{r}') [\rho^+ (\vec{r}') f^{++}(\vec{r}') - \rho^- (\vec{r}') f^{++}(\vec{r}')],$$  

(26a)

$$f^{+-}(\vec{r}) = - \psi_1(\vec{r}) - \int d^2 \vec{r}' \psi_1(\vec{r} - \vec{r}') [\rho^- (\vec{r}') f^{+-}(\vec{r}') - \rho^+ (\vec{r}') f^{+-}(\vec{r}')],$$  

(26b)

$$f^{-+}(\vec{r}) = \psi_1(\vec{r}) + \int d^2 \vec{r}' \psi_1(\vec{r} - \vec{r}') [\rho^- (\vec{r}') f^{-+}(\vec{r}') - \rho^+ (\vec{r}') f^{-+}(\vec{r}')],$$  

(26c)

$$f^{--}(\vec{r}) = - \psi_1(\vec{r}) - \int d^2 \vec{r}' \psi_1(\vec{r} - \vec{r}') [\rho^+ (\vec{r}') f^{--}(\vec{r}') - \rho^- (\vec{r}') f^{--}(\vec{r}')].$$  

(26d)

In the current context, Eq. (21) reduces to

$$\psi_1(\vec{r}) = \psi^+ (\vec{r}) = - \psi^- (\vec{r}) = \ln(|\vec{r}|) + \frac{\sqrt{\pi} \Gamma}{|\vec{r}|},$$  

(27)

where we orient our $(x, y)$ coordinate system in such a way that the slip direction points along the $x$ direction. The minus signs in Eqs. (26a)–(26d) arise from a sign difference in the interactions between plus-plus dislocations versus plus-minus dislocations, as shown in Eq. (27). By comparing Eq. (26a) against Eq. (26d) and Eq. (26b) against Eq. (26c), we find that $f^{++}(\vec{r}) = - f^{--}(\vec{r})$ and $f^{+-}(\vec{r}) = - f^{--}(\vec{r})$. These symmetries further imply that $f^{++}(\vec{r}) = f^{--}(\vec{r})$. Finally, we obtain

$$f^{++}(\vec{r}) = \psi_1(\vec{r}) + \int \psi_1(\vec{r} - \vec{r}') [\rho^+ (\vec{r}') + \rho^- (\vec{r}')] d^2 \vec{r}'.$$  

(28)

Our general formulation in Sec. III allows for spatial variation of an uncorrelated density $\rho(\vec{r})$. Without externally applied force, $\rho(\vec{r}) = (N / \lambda)\gamma$ is constant in space. An analytical solution to Eq. (28) can be obtained for constant $\rho^+$ and $\rho^-$. The dimensionless nature of the interaction potential $\psi_1$ suggests a change in variable $\sqrt{\rho^+ + \rho^-} \vec{r} \rightarrow \vec{r}$ (note that $\rho^+$ and $\rho^-$ are always positive). The resulting dimensionless integral equation

$$f^{++}(\vec{r}) = \psi_1(\vec{r}) + \int \psi_1(\vec{r} - \vec{r}') f^{++}(\vec{r}') d^2 \vec{r}'$$  

(29)

can be directly solved by applying $\Delta^2 = (\partial^2 + \partial^2)$ on both sides of the equation and using the identity

$$\Delta^2 \psi_1(\vec{r}) = 2 \pi \Delta \partial (\vec{r}) + 2 \pi (\partial^2 - \partial^2) \partial (\vec{r}) = 4 \pi \partial^2 \partial (\vec{r}).$$  

(30)

Equation (29) then becomes

$$\Delta^2 f^{++} = 4 \pi \partial^2 [f^{++} + \partial (\vec{r})],$$  

(31)

whose explicit solution is

$$f^{++} = \frac{\gamma}{r} \sinh (\sqrt{\pi} r) K_1(\sqrt{\pi} r) - \cosh (\sqrt{\pi} r) K_0(\sqrt{\pi} r),$$  

(32)

where $K_0(\cdot)$ and $K_1(\cdot)$ are the zeroth- and first-order modified Bessel functions of the second kind. With the aid of Eqs. (3a) and (3b), the correlation functions $g^{(++)} = g^{(--)}$ and $g^{(+-)} = g^{(-+)}$, correct to $O(\Gamma^2)$, can be expressed in the original coordinates,

$$g^{(++)}(\vec{r}) = 1 + \frac{\gamma}{r} \sinh (\sqrt{\pi} r) K_1(\sqrt{\pi} r) - \cosh (\sqrt{\pi} r) K_0(\sqrt{\pi} r),$$  

(33a)

$$g^{(+-)}(\vec{r}) = 1 - \frac{\gamma}{r} \sinh (\sqrt{\pi} r) K_1(\sqrt{\pi} r) - \cosh (\sqrt{\pi} r) K_0(\sqrt{\pi} r),$$  

(33b)

where $\sigma_0 = \sqrt{\pi} \Gamma (\rho^+ + \rho^-)$ gives an inverse “Debye radius” of the dislocation cloud. The third-order correlation functions correct up to $O(\Gamma^2)$ straightforwardly follow from Eq. (3b). The validity of Eq. (32) can be verified by comparing $g^{(++)}(\vec{r}) - g^{(--)}(\vec{r})$ to the dislocation difference, or GND, field $\kappa(\vec{r})$ in Eq. (15) of Ref. 15. In this latter work, the same expression is obtained for the induced GND due to a single pinned dislocation, which was interpreted by the authors as the pair correlation of dislocations in a relaxed system.

The values of $g^{(++)}$ and $g^{(+-)}$ in Eq. (32) diverge—and hence are unphysical—at very small radii. The reason arises
from the omission of a core energy term in Eqs. (20) and (27) which would otherwise offset the logarithmic divergence. Physically speaking, when two dislocations approach each other, if they are of the same sign, they merge into one dislocation; otherwise they cancel. In both cases, nothing is diverging. Note also that the pair correlation functions depend only on the scaled space coordinate $\sqrt{pr}$ ($p = \rho^+ + \rho^-$ being the total dislocation density) in agreement with the scaling argument given by Zaiser et al. This dependence also holds in the multiple-slip case to be discussed in Sec. V.

V. PAIR CORRELATION FUNCTIONS FOR MULTIPLE SLIP

The procedure to obtain the correlation functions for a system with multiple slips follows the same types of arguments and expansions as those for single slip. We shall further develop the integral of Eq. (25) for a system of $N$ slip systems, each with two charges, and subsequently give an explicit analytical solution for the pair correlation function in the case wherein the difference in slip-orientation angle between adjacent slip planes is constant.

For an $N$-slip system with both types of charges, we have $4N^2$ coupled integral equations for different pairs of 1 and 2 in Eq. (25). To reduce the number of equations, and essentially decouple them, some symmetry arguments can be employed. For an infinite system,

$$\tilde{\psi}_{ij}^+=\tilde{\psi}_{ij}^-=-\tilde{\psi}_{ji}^+, \quad \tilde{\psi}_{ij}^{ab} = \tilde{\psi}_{ji}^{ab},$$

where the superscripts denote the charges of the first and second dislocations, while the subscripts show the slip systems in which they live. Equation (25) can be recast by using the convolution operator $*$ and the symmetry of $\tilde{\psi}_{ij}^{ab}$ as

$$f_{ij}^{ab} = \tilde{\psi}_{ij}^{ab} + \sum_{k=1}^{N} \tilde{\psi}_{ik}^{ab} * [\rho_k f_{jk}^{ab} - \tilde{\rho}_k f_{jk}^{ab}].$$

By direct substitution of $+$ and $-$ into $a$ and $b$, it is immediate that $f_{ij}^{+-}(\hat{r}) = -f_{ij}^{-+}(\hat{r})$ and $f_{ij}^{ab}(\hat{r}) = -f_{ij}^{ba}(\hat{r})$, which further implies that

$$f_{ij}^{+-} = f_{ij}^{-+} = \sum_{k=1}^{N} \tilde{\psi}_{ik}^{ab} * [\rho_k f_{jk}^{ab} + \tilde{\rho}_k f_{jk}^{ab}].$$

With this, Eq. (34) reduces to

$$f_{ij} = \tilde{\psi}_{ij} + \sum_{k=1}^{N} \tilde{\psi}_{ik} * [\rho_k f_{jk}],$$

where the superscripts have been omitted and $\rho_k = \rho^+ + \rho^-$. The total dislocation density of both types on slip $k$ is effectively reduced the number of coupled equations to $N^2$. Also note that because of $\tilde{\psi}_{ij} = \tilde{\psi}_{ji}$, there are only $N(N+1)/2$ independent $\tilde{\psi}_{ij}$’s.

As seen from the single-slip case, Eq. (36) subjected to an arbitrary distribution of the local density $\rho_k(\hat{r})$ cannot be analytically solved. For spatially independent $\rho_k$, however, these equations can be decoupled. Let $\lambda_k$ be the relative population of density in slip system $k$ relative to the total density $\rho$, i.e., $\rho_k = \lambda_k \rho$, where $\sum_{k=1}^{N} \lambda_k = 1$. We can then perform a change in variable $\sqrt{pr} \rightarrow \tilde{r}$ to absorb the $\rho$ dependence. In addition, in the Fourier space (indicated by a superposed \textasciitilde), a convolution becomes a product. We can solve the Fourier transform of Eq. (36) for $f_{ij}$ by essentially performing a matrix inversion on

$$\tilde{\psi}_{ij} = \sum_{m,n} (\delta_{im} \delta_{jn} - \lambda_n \tilde{\psi}_{im} \delta_{jn}) \tilde{f}_{mn}.$$  

The Fourier representation of $\tilde{\psi}_{ij}$ in Eq. (21) can be very simply expressed in polar coordinates $(k, \phi_k)$,

$$\tilde{\psi}_{ij} = -\frac{4\pi}{k^2} \sin(\phi_k - \theta_i) \sin(\phi_k - \theta_j) = -\frac{4\pi}{k^2} (\hat{m}_i \cdot \hat{k})(\hat{m}_j \cdot \hat{k}),$$

where $\theta_i$ is the angle that slip plane $i$ makes with the $x$ axis (which can be arbitrarily chosen, so that $\theta_i = i\pi/N$). Owing to the simple form of Eq. (38), the solution to Eq. (37) is

$$\tilde{f}_{ij} = \frac{\tilde{\psi}_{ij} / \lambda_{ij}}{1 - \sum_n \tilde{\psi}_{in}},$$

where we have used $\sum_n \tilde{\psi}_{in} \tilde{\psi}_{ij} = \tilde{\psi}_{ij} \sum_n \tilde{\psi}_{in}$. Equation (39) shall be used in the derivation of the evolution equation for parallel edge dislocations in a multislip system in Sec. VI.

To verify that Eq. (39) is applicable in glide-controlled systems, we consider an ensemble of 1500 relaxed configurations of 64 plus and 64 minus dislocations randomly placed on a $1 \mu m^2$ square and restricted to move along their glide directions. The simulations were performed with periodic boundary conditions in the absence of thermal noise. The glide constraint helps prevent dislocation annihilation, and thus, to fix the total number of dislocations and to maintain the finite effective temperature. As an example, Fig. 1 shows the density plot of the theoretical correlation function $f_{ij}^{+-}$ between plus dislocations on $60^\circ$ and $120^\circ$ slip systems against (a) the simulation result. The erroneous oscillations in Fig. 1(b) along $0^\circ$ and $90^\circ$ lines are caused by the numerical inverse Fourier transform operation of Eq. (39). (The general closed form solution of a double-slip pair correlation function does not exist for an arbitrary pair of slip-orientation angles.) Overall, the theory gives accurate angular predictions except along the two slip directions where it underpredicts the same-sign anticorrelation due to the suppression of climb. The plot of the correlation function along the $\hat{x}$ axis is shown in Fig. 2. Very close to the origin, the function logarithmically diverges as does the unscreened potential. About one dislocation spacing from the core, the correlation function decays as $1/\hat{x}^2$.

The real-space solution to Eq. (39) is possible if we assume that the \emph{angle between each adjacent pair of slip planes is constant}. For any $N \in \mathbb{Z}^+$ and $N > 1$,
regardless of $\phi_k$. With the above identity, the denominator of \( \tilde{f}_{ij} \) becomes angular independent and can be directly integrated. The final result, with

\[ k_0 = \sqrt{2\pi N\Gamma}\rho, \]

reads

\[
\sum_{n=1}^{N} \frac{\sin^2\left(\phi_n - \frac{n\pi}{N}\right)}{N} = \frac{N}{2}
\]

FIG. 1. (Color online) (a) Discrete dislocation result and (b) theoretical prediction of the correlation function \( f_{12}^{\text{dis}} \) between plus dislocations on \( 60^\circ \) and \( 120^\circ \) slip systems. The values increase toward brighter regions. The coordinates are measured in units of \( 1/\sqrt{\rho} \). The dashed lines indicate the two slip directions where the plus-plus anticorrelation is underpredicted due to the glide constraint of the discrete dislocation simulations. The fitting parameter due to rescalings of length was found to be \( k_0 = 22\sqrt{\rho} \).
\( \partial_t \rho_i = - (\hat{b}_1 \cdot \hat{v})(- \rho_i (\hat{r}_i, t) \tau_{i}^{\text{ext}} + \sum_j d_i^2 \tau_j^\text{ind}(\hat{r}_i, \hat{r}_j, t)) \),

\( \tau_{ij}^{\text{ind}}(\hat{r}) = \hat{s}_i \cdot \sigma_{ij} \cdot \hat{m}_i = Gb(\hat{s}_i \cdot \hat{V})(\hat{m}_i \cdot \hat{V})[r^2 \ln r]. \)

Here, \( G = \mu / [2(1-\nu)] = E/[4\pi(1-\nu^2)] \), where \( E, \mu, \) and \( \nu \) are the Young modulus, the shear modulus, and the Poisson ratio, respectively.

Addition and subtraction of Eqs. (42a) and (42b) give the evolution equations for the total dislocation density \( \rho_i = \rho_i^s + \rho_i^t \) and the GND density \( \kappa_i = \kappa_i^s - \rho_i \).

\( \delta \rho_i = - (\hat{b}_1 \cdot \hat{v})(\kappa_i \tau_{i}^{\text{ext}} + \sum_j d_i^2 \tau_j^\text{ind}(\hat{r}_i, \hat{r}_j, t)) \),

\( \delta \kappa_i = - (\hat{b}_1 \cdot \hat{v})(\rho_i \tau_{i}^{\text{ext}} + \sum_j d_i^2 \tau_j^\text{ind}(\hat{r}_i, \hat{r}_j, t)) \).

In accordance with Eq. (11), the dislocation-dislocation density can be written as

\( \rho_{ij}^{s^2} = \rho_i^s(\hat{r}_i) \rho_j^s(\hat{r}_j) g_{ij}^{s^2} (\hat{r}_i - \hat{r}_j) = \rho_i^s(\hat{r}_i) \rho_j^s(\hat{r}_j)[1 + d_{ij}^{s^2}(\hat{r}_i - \hat{r}_j)], \)

where \( s, s' \in \{+,-\} \) and, according to Eq. (23a) and (23b),

\( d_{ij}^{s^2} = \Gamma_{ij}^{s^2} \). In terms of the single and pair correlation functions, the total dislocation density and GND are

\( \rho_{ij}^{(2)} = \rho_i(\hat{r}_i) \rho_j(\hat{r}_j) + \frac{1}{2} \{ \rho_i(\hat{r}_i) \rho_j(\hat{r}_j) d_{ij} + \rho_i(\hat{r}_i) \kappa_j(\hat{r}_i) [d_{ij}^{s^2} + d_{ij}^{t^2}] \}

+ \kappa_i(\hat{r}_i) \rho_j(\hat{r}_j) [d_{ij}^{s^2} - d_{ij}^{t^2}] \),

\( \kappa_{ij}^{(2)} = \kappa_i(\hat{r}_i) \kappa_j(\hat{r}_j) + \frac{1}{2} \{ \rho_i(\hat{r}_i) \rho_j(\hat{r}_j) [d_{ij}^{s^2} - d_{ij}^{t^2}] + \rho_i(\hat{r}_i) \kappa_j(\hat{r}_i) d_{ij}^{s^2} \}

- \kappa_i(\hat{r}_i) \rho_j(\hat{r}_j) d_{ij}^{t^2} + \kappa_i(\hat{r}_i) \kappa_j(\hat{r}_j) [d_{ij}^{s^2} + d_{ij}^{t^2}], \)

where \( d_{ij} = d_{ij}^{s^2} + d_{ij}^{t^2} \), \( d_{ij}^{s^2} = (1/2)(d_{ij}^{s^2} + d_{ij}^{t^2}) \), and \( d_{ij}^{t^2} = (1/2)(d_{ij}^{s^2} - d_{ij}^{t^2}) \). After substitution of Eq. (45) and (46), Eq. (44) becomes

\( \partial_t \rho_i = - (\hat{b}_1 \cdot \hat{v})(\kappa_i (\tau_{i}^{\text{ext}} + \tau_{i}^{s^2} - \tau_{i}^{t^2}) + \rho_i \tau_i^{s^2}), \)
The Fourier transform of $\tilde{x}_{ij}^{\text{ind}}$ can be directly computed from Eq. (43),
\[
\mathcal{F}[\tilde{x}_{ij}^{\text{ind}}](\vec{k}) = -4\pi Gb \tilde{\phi}_j \frac{1}{k^4} \left[ (\hat{s}_j \cdot \hat{k})(\hat{m}_j \cdot \hat{k}) \right] \frac{(\vec{s}_j \cdot \vec{\hat{k}})(\vec{m}_j \cdot \vec{\hat{k}})}{k^4}.
\]

Owing to the connection $d^i_j(\vec{x}) = \Gamma f_{ij}(\vec{x})$, Eq. (50) becomes, from Eqs. (39) and (51),
\[
\tilde{I}_{ij} = \frac{\Gamma^2 Gb}{\lambda_j} \int \tilde{\phi}_j N^i_j(\vec{s}_j \cdot \vec{\hat{k}}) d^2 \vec{k}.
\]

The vector $\tilde{I}_{ij}$ is most conveniently expressed in the coordinate system of slip $j$. Substitution of Eq. (38) into Eq. (52), while projecting $\hat{s}_i$ and $\hat{m}_j$ onto ($\hat{s}_j, \hat{m}_j$), gives
\[
\tilde{I}_{ij} = (4\pi)^2 \frac{\Gamma^2 Gb}{\lambda_j} \frac{\hat{s}_j}{k^2 + 4\pi \sum_n \sin^2(\phi_n - \theta_n)} \times \left\{ \int_0^{2\pi} \int_0^\infty \begin{matrix} \frac{1}{k} \sin^2(\phi_k) \sin(\phi_k + \theta_j) \sin(3\phi_k + 2\theta_j) d\phi_k \bigg/ k^2 + 4\pi \sum_n \sin^2(\phi_k - \theta_n) \end{matrix} \\
\times dk d\phi_k \right\}.
\]

where $\theta_j = (j - i) \pi / N$ is the angle between slip planes $i$ and $j$. We impose a cutoff $\epsilon$ at small $k$ to prevent the logarithmic divergence due to the long-range nature of the pair correlation functions.

Under the assumption of equal interval of successive slip orientation, as in Sec. V, we can very straightforwardly carry out the above integrals, which give
\[
\tilde{I}_{ij} = \frac{GD}{\lambda_j} \cos(\theta_j) \hat{s}_j,
\]

where $D = 2\pi^2 \Gamma^2 \ln \epsilon / N$ serves as a fitting parameter. The factor $\lambda_j$, nicely combines with $\rho$ in the denominator of Eq. (55) to make $\rho_j = \lambda_j \rho$.

Each of these functions are of an ensemble of spatially constant single-dislocation densities in thermal equilibrium. When the distributions of single-dislocation densities are nonuniform in space as is the case for systems out of equilibrium, the back stress response should depend on how much the densities locally vary.

The final result is amazingly simple,
obtain the desired fit. In their analysis, Yefimov et al.\textsuperscript{12} used the value of $D$ from their previous single-slip theory without any readjustment. There is no \textit{a priori} reason why this value should stay unaltered. The density of nucleation sources in their continuum theory was chosen to match that in the discrete dislocation simulations. The discrepancy could also arise from different ways in which the discrete dislocation theory and the continuum theory handle dislocation nucleation.

In a later publication, Yefimov and Van der Giessen\textsuperscript{14} applied their formalism to the problem of stress relaxation in single-crystal thin films on substrates subjected to thermal loading. Due to the difference in thermal expansion coefficients between film and substrate, high tensile stresses can develop in the films as the temperature decreases. Contrary to the discrete dislocation simulations by Nicola et al.,\textsuperscript{46,47} which show increasing stress built up inside a film with decreasing film thickness, the results from the continuum theory show a size-dependent hardening only during the early stage of cooling. Moreover, the theory gives identical results between some pair of slip orientations (e.g., when the angle between the two slip planes $\theta_i$ is either 60° or 120°), whereas the discrete dislocation simulations and our theory predict otherwise. Finally, in the previous continuum theory,\textsuperscript{25} dislocations nucleate when the sum of the external stress $\tau^{ext}$, the self-consistent long-range stress $\tau^c$, and back stress $\Phi$ exceed a certain value. From our analysis, we believe that, in a more correct treatment of dislocation nucleation, this back stress should be supplemented by flow stress $\Phi$ [Eqs. (48c)] which is dominant in a nucleation region where plus and minus dislocations are equally populated. Applications of the current theory to the shearing problem and the thin film problem, which shows the size-dependent hardening, will shortly appear following this publication.

**VIII. DISCUSSION AND CONCLUSIONS**

We have described $n$th-order dislocation densities and dislocation pair correlation functions in a grand canonical ensemble and obtained the relationships between different orders of the correlation functions in the form of a hierarchy of integral equations. By using the Bogolyubov ansatz instead of the more customary Kirkwood approximation, we have closed the chain of the equations at second order and solved for approximate expressions of the pair correlation functions—valid at all distances—for systems with one slip and multiple active slip systems. These solutions are invariant under simultaneous transformations $\vec{r}\rightarrow\vec{r}/\sqrt{\rho}$ and $\rho\rightarrow\rho^2$. The transformations suggest that any emergent dislocation pattern should exhibit a length scale given by $1/\sqrt{\rho}$ as pointed out by Holt\textsuperscript{48} and in agreement with the “law of similitude.”\textsuperscript{49} For a complete analysis of scaling relations, the reader is referred to Ref. 10.

Recently, Groma et al.\textsuperscript{15} have developed a mean-field variational approach to study the screening of dislocations, which is similar in spirit to the Debye–Hückel theory in the study of classical plasmas.\textsuperscript{50–52} This method is based on approximating the system’s total density matrix as a product of single-particle density matrices $\rho_i$ with the free energy given by $F=\langle H\rangle+T\Sigma_i Tr \rho_i \ln \rho_i$. Although this technique provides a complimentary approach and results in the same pair correlation expressions for a single-slip system (after some interpretation), its generalization to multiple-slip system is not obvious. In particular, one would have to supply additional cross couplings between different slips by hand. These couplings should automatically emerge from a complete theory.

In Sec. IV, we have formulated transport equations for the total dislocation and GND densities for general multiple slip. Interactions among dislocation pairs produce an additional (relatively) short-ranged “back stress” contribution to the long-range internal stress of individual dislocations. Most of the complexities of the correlation functions were integrated away, which leave only the cos($\theta_{ij}$) coupling between slip systems $i$ and $j$ [see Eq. (55)]. This dependence was also proposed by Gurtin\textsuperscript{12,43} in his strain gradient plasticity theory, but it was abandoned by Yefimov et al.\textsuperscript{14,25} We have argued in Sec. VII that this refusal was based on an unfair comparison with discrete dislocation simulations for the way in which dislocation nucleation was treated.

There is an important issue regarding the use of dislocation correlations $f_{ij}$ or $d_{ij}^*$ in Sec. V. The formalism developed in Sec. III assumes that dislocations relax along the directions dictated by the Peach–Koehler forces. This implies dislocation glide, as included in the transport equations developed in Sec. VI, but also climb which is not considered a mechanism of plastic flow here. Mathematically speaking, Eq. (17) is \textit{not} the stationary state of Eqs. (42a) and (42b). Early attempts in numerically describing dislocation correlations in glide-only, multiple-slip systems failed to produce noticeable patterns due to the need for large number of dislocations; the role of climb (or cross slip) was suggested to help overcome this difficulty.\textsuperscript{53,54} The original motivation for our approach was to find the orientation dependence of the back stress in the most straightforward way. Extracting the angular dependence from a climb-assisted relaxed state gave us a quick input to use in the glide-only, multiple-slip theory. The validity of the continuum theory will always be vindicated by comparisons against discrete dislocation results.

Finally, we believe that our multiple-slip formulation provides a framework to address a long standing challenge in explaining dislocation patterning. For single-slip systems, short-range correlations occur between two dislocations except along directions normal to their glide plane (taken to be along $\hat{y}$). It has been shown that for a small deviation away from this “dislocation wall” direction, an attractive parabolic potential produced by the correlated dislocations decays as $|y|^{-5/2}$, which is compared to $|y|^{-2}$ in the unscreened case.\textsuperscript{15} We have found in Sec. V, however, that when one or more extra slips are introduced, the effect of Debye-type screening diminishes. In this case, the attractive potential, in fact, decays like $r^{-2}$ as if it was unscreened. This could explain the necessity to introduce extra slips to see the formation of walls in discrete dislocation simulations,\textsuperscript{17–24} \textit{unless} further aided by climb motions.\textsuperscript{33,55} The latter suggests the existence of a critical exponent of the attractive potential below which structure formation cannot occur, as is the case in single-slip systems restricted to glide. A more detailed investigation of this is left for future work.
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5For a summary of various continuum theories, see, e.g., Ref. 56 and references therein.
26We assume that the noise is weak enough that dislocation climb is not possible.
27Note that this limits the analysis to two-dimensional systems of dislocations.
28One distinction due to the choice of an ensemble type can be seen from the normalization condition. In the grand canonical ensemble, according to Eqs. (8) and (9), $\int \rho^{(N)}(\mathbf{r})d\mathbf{r} = \sum_{N=1}^{\infty} N! \left[ \prod_{n=1}^{N} (N_n)! \right] \langle \mathbf{r}_{\mathbf{n}} \rangle$, where $\langle \cdot \rangle$ denotes an average over all statistically equivalent ensembles. In particular, the density of a single dislocation of species $s$ in a system with no external shear, $\rho^{(1)}(\mathbf{r})$, is independent of $\mathbf{r}$, thus, $\int \rho^{(1)}(\mathbf{r})d\mathbf{r} = \rho^{(1)}(A) = \langle N_s \rangle$. In other words, $\rho^{(N)} = \langle N_s \rangle/A$ depends on the average number of dislocation of species $s$. If one were to carry a similar analysis in a canonical ensemble where the number of dislocations of each species is fixed, $\rho^{(1)}$ would have to be replaced by $N_s/A$, where $N_s$ is fixed. For higher order density, the expression becomes quite cumbersome. For example, $\rho^{(2)}(\mathbf{r}_s, \mathbf{r}_s') = \frac{N_s N_s'}{A^2} \delta(\mathbf{r}_s - \mathbf{r}_s')$, while for $s = s'$, it is $\frac{N_s}{(N_s - 1)/A^2}$. In thermodynamic limit ($N_s \rightarrow \infty$ and $A \rightarrow \infty$ while keeping the ratio fixed), these two expressions reduce to the same thing. In this sense, it is cleaner to work in the grand canonical ensemble.
32H. S. Green, Molecular Theory of Liquids (North-Holland, Amsterdam, The Netherlands, 1952).
35In the presence of an external conservative force, it can be shown that both Eqs. (17) and (19) remain valid provided that an additional term representing the applied external force, $F(\mathbf{r}) = -(1/T)\nabla \Phi(\mathbf{r})$ generated by the external potential $\Phi(\mathbf{r})$ which acts on $\mathbf{r}$, is added to their RHS. Qualitatively speaking, the original expression is nothing but the sum of all the Peach–Koehler interactions on the dislocation at $\mathbf{r}$ due to all other dislocations in the collection $\mathbf{n}$.
38There are some systems where the critical strain is naturally small such as dislocations in vortex lattices of type II superconductor where the values of elastic moduli can be small at suitably applied magnetic field (Ref. 57). By modifying the form of the interaction potential, the present analysis can be carried over straightforwardly.
39The form of the solution is not surprising; it suggests that the solution can be written as a sum of diagrams due to the expansion $1/(1-x) = 1 + x + x^2 + \cdots$, which is often encountered in a many-body theory.