

Coagulation and diffusion: a probabilistic perspective on the Smoluchowski PDE

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Abstract

The Smoluchowski coagulation-diffusion PDE is a system of partial differential equations modelling the evolution in time of mass-bearing Brownian particles which are subject to short-range pairwise coagulation. This survey presents a fairly detailed exposition of the kinetic limit derivation of the Smoluchowski PDE from a microscopic model of many coagulating Brownian particles that was undertaken in [10]. It presents heuristic explanations of the form of the main theorem before discussing the proof, and presents key estimates in that proof using a novel probabilistic technique. The survey's principal aim is an exposition of this kinetic limit derivation, but it also contains an overview of several topics which either motivate or are motivated by this derivation.

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1 Introduction

1.1 Microscopic particles and macroscopic descriptions

An important aim in statistical mechanics is to explain how the huge amount of information available in a microscopic description of a physical object, such as the positions and momenta of all the molecules comprising the air in a room, may be accurately summarised by first specifying a small number of physical parameters which are functions of macroscopic location, such as the density, temperature and pressure of this body of air at different points in the room, and then determining how these parameters evolve in space and time.

1.1.1 The elastic billiards model and the heat equation

The microscopic system may begin out of equilibrium: for example, a still body of warm air in one room may be separated by a partition from another still body of cooler air in another, and then the partition instantaneously removed, so that air molecules from one side and the other intermingle over time, and an equilibrium is eventually approached in which the body of air in the whole room is again close to still, at a temperature which is some average of those of the two isolated systems at the original time. In such a case as this, it is a natural task to seek to summarise the evolution of a few suitable macroscopic physical quantities as the solution of partial differential equations. In the example, our object of study might be the temperature of the gas, and our aim to show that it is the heat equation, $\frac{\partial}{\partial t}T(x,t) = \Delta T(x,t)$, which models the macroscopic evolution $T(x,t)$ (with x varying over the whole room, $[-1, 1]^3$, say) of the temperature from the moment of the removal of the partition at time $t = 0$ until a late time, $t \rightarrow \infty$, at which a new equilibrium is approached. In an idealized and very classical choice of microscopic description of the gas, we might model the ensemble of air molecules as a system of tiny spheres of equal radius and mass, each moving according to some velocity, and each pair of which undergoes a perfectly elastic collision on contact, in the same manner that a pair of billiards would. On each of the walls that comprise the boundary $\partial[-1, 1]^3$ of the room, each sphere bounces elastically. The partition is modelled by the immobile sheet $\{0\} \times [-1, 1]^2$ on which spheres on either side also bounce elastically before time zero; the partition is removed instantaneously at that time. The initial instant of time may be taken to be zero, or some negative time. At that moment, we may scatter the spheres in an independent Poissonian manner throughout the room $[-1, 1]^3$ (the reader may notice that in fact some extra rule is needed here to ensure the spheres' disjointness); and on one and other side of the partition, choose their velocities independently, those to the right of partition according to a non-degenerate law of zero mean, and those on the left according to another such law of lower variance than the first; in this way, we model two bodies of still air, a warm one in the right chamber $[0, 1] \times [-1, 1]^2$, and a cooler one in the left $[-1, 0] \times [-1, 1]^2$. In the microscopic model, there are huge numbers of tiny spheres in the system. Indeed, we may seek to understand the macroscopic evolution of temperature by in fact considering a whole sequence of microscopic models indexed by total particle number N , in a limit of high N . In the N -th model, spheres are initially scattered as we described, with a Poissonian intensity N throughout $[-1, 1]^3$. To carry out this task of understanding the large-scale evolution, we would wish to specify a microscopic definition of the notion of temperature, and then

explain how it is in the high N limit that the microscopic temperature data may be meaningfully reduced to a macroscopic description, and that this latter description indeed evolves according to the heat equation. Microscopically, temperature is interpreted [28, Section 1.1] as the average kinetic energy of particles, where here the velocity of particles is measured relative to the average velocity of nearby particles. Since our particle systems are large microscopically, considering as we do a high N limit, we may specify in our N -th microscopic model a definition of temperature at any given location $x \in [-1, 1]^3$ as follows: first we may compute the mean velocity $v_{N,\delta}(x)$ of the set of spheres whose centres lie within some small distance δ of a given location x in the room, and then we are able to define the microscopic temperature $T_{N,\delta}(x)$ to be the average of the square of the particle velocity minus $v_{N,\delta}(x)$, where the average is taken over the same set of spheres. Of course, the value $T_{N,\delta}(x)$ will change in time. As N approaches infinity with δ being fixed but small, huge numbers of particles are involved in the empirical counts used for averaging. Our aim is to consider the space-time evolution of the microscopically specified temperatures after the high N limit is taken, at which point, the weak law of large numbers might suggest that the concerned empirical counts behave non-randomly to first order, so that our description becomes deterministic: the limit $T_{N,\delta}(x)$ will be some non-random function $T_\delta(x)$. In fact, since δ is fixed, we should not yet expect our system to approximate the heat equation, since there is an effect of macroscopic smearing in our calculation of microscopic temperature. Rather, one might expect the heat equation description to emerge if we take a $\delta \searrow 0$ limit of $T_\delta(x)$, after the first high N limit has been taken. Moreover, to hope to obtain this description, we will also need to scale time appropriately in the N -th microscopic model, as we take the first, high N , limit. In the scaled time coordinates, the microscopic models should make their approach to the new thermal equilibrium at the same rate, as $N \rightarrow \infty$. What rate this is in fact depends on another important consideration concerning the microscopic models which our brief description left unspecified: the radius r_N of each sphere in the N -th model must certainly be chosen to satisfy $r_N^3 \leq cN^{-1}$ for some constant $c > 0$, if only to permit all of the spheres to inhabit the room disjointly; our choice of decay rate for r_N as a function of N , subject to this constraint, will determine the factor by which we scale time in the N -th model in order to seek a heat equation description in the large.

To implement the programme proposed in the preceding paragraph is an open problem, and in all likelihood, an extremely difficult one. There is no randomness in the model except in the initial selection of particle locations and velocities: from that time on, the deterministic laws of Newtonian mechanics govern the evolution of the microscopic models. Moreover, some choices for density admissible in the above description – such as when Nr_N^3 converges in high N , to a suitably small constant – lead to rather dense systems of particles. The derivation may be less ordinarly hard were more dilute choices of limit considered, where r_N converges to zero more, and perhaps much more, quickly than does $N^{-1/3}$.

It is important to note, however, that, if a choice of r_N as a function of N is made which is too rapidly decaying, we may leave the realm in which the heat equation is the appropriate macroscopic description. For example, if $r_N = o(N^{-1/2})$, it is a simple matter to check that a typical sphere after time zero will cross traverse the entire room on many occasions before meeting any other particle. The system will reach equilibrium after the removal of the partition simply by the free motion of the particles. The heat equation is only a suitable description when a typical particle

experiences the thermal agitation caused by its collision with many other particles in short periods of macroscopic time.

1.1.2 The elastic billiards model and Boltzmann's equation

Moreover, the elastic billiards model crosses at least one interesting regime as it is diluted from the dense $r_n = \Theta(N^{-1/3})$ phase towards the trivial free motion phase $r_N = o(N^{-1/2})$. Consider the choice $r_N = N^{-1/2}$. A moment's thought shows that, in this regime, a typical sphere will travel (at unit-order velocity) for a duration before its first collision with another particle which on average neither tends to zero nor to infinity as $N \rightarrow \infty$. This is the regime of *constant mean free path*. The heat equation will not offer a suitable description for the evolution of temperature in this regime, because the mechanism providing for thermal agitation of particles occurs on a time scale which is marginally too slow. However, the programme of deriving a macroscopic description by means of a PDE does make sense, and in this case, offers a powerful model of gas dynamics. Suppose that, instead of using the microscopic data to form a description of temperature, we use it to describe the density of particles having a given velocity $v \in \mathbb{R}^d$ nearby a given location $x \in [-1, 1]^3$. Particles may be scattered in a Poissonian fashion as before at the initial time, but with inhomogeneities in the intensity of this scattering permitted in both the space and velocity variables. With the macroscopic smearing parameter now being used to approximate velocity $v \in \mathbb{R}^d$ as well as location $x \in [-1, 1]^3$, we may record a microscopic description $f_N^\delta(x, v)$ for the δ -smeared density of spheres at space-velocity location (x, v) . Taking a high N and then low δ limit as above, our macroscopic evolution is modelled by the fundamental system of equations in gas dynamics, Boltzmann's equation, valid for $t \geq 0$, $x \in \mathbb{R}^d$ and $v \in \mathbb{R}^d$:

$$(1.1) \quad \frac{\partial}{\partial t} f(x, v) = -v \cdot \nabla_x f(x, v) + Q(f, f).$$

Here, $-v \cdot \nabla_x$ is the free motion operator associated to particles of velocity v , while $Q(f, f)$ is a binary collision operator that reflects the microscopic elastic collision and whose form we will specify when we return to Boltzmann's equation in a brief discussion in Section 2. For now, note that the time evolution of the macroscopic densities is governed both by the free motion and by the collision operator. This is what is to be expected in the regime of constant mean free path, where the typical particle experiences unit-order durations free of collision and other such periods where several collisions occur.

Boltzmann carried out a derivation (1.1) as a model of gas dynamics in 1872, based on several assumptions, including one of molecule chaos that he called the Stosszahlansatz and which we will later discuss. (See [5] for an English translation of his 1872 article.) The validity of his derivation was a matter of controversy, not least due to Loschmidt's paradox concerning precollisional particle independence (see Subsection 2.3.2), and it was a fundamental advance made in 1975 by Lanford [15] when the programme of rigorously deriving Boltzmann's equation from the elastic billiards model in the regime of mean free path was successfully implemented, for a short initial duration of time. By the latter condition, we mean that the validity of the description was established for some non-zero finite period, whose value depends on the form of the initial density profile of particles in space-velocity.

Lanford derived Boltzmann's equation by establishing that the correlation functions concerning several particles in the model satisfy a hierarchy of equations called the BBGKY hierarchy, where the index of an equation in the hierarchy is the number of particles whose correlation is being considered, and by showing that when the correlation functions adhere to the BBGKY hierarchy, the density profile follows Boltzmann's equation. Illner and Pulvirenti implemented this approach in [12] in order to derive Boltzmann's equation in a similarly sense, now globally in time, but with a similar smallness condition, now on sparseness of the initial particle distribution.

1.1.3 Our main goal: coagulating Brownian particles and the Smoluchowski PDE

This survey is intended to offer a detailed overview of a programme for deriving the macroscopic description of a gas of particles in the same vein as the descriptions above propose. However, our microscopic particles will diffuse, each following a Brownian trajectory, and as such their evolution is random, not deterministic; the mechanism of interaction will be pairwise as above, but a coagulation in which only one particle survives rather than a collision in which both do. On the other hand, in an effort to provide some generality in the microscopic description and richness in the macroscopic one, each of the particles will bear a mass, which the pairwise coagulation will conserve; and, moreover, we will permit the diffusivity of the Brownian trajectory of each particle to depend on the particle's mass.

The partial differential equation which the programme seeks to obtain in this case – the analogue of the heat equation or Boltzmann's equation in our opening examples – is, like Boltzmann's equation, in fact a system of PDE, in our case coupled in the mass parameter, known as the Smoluchowski coagulation-diffusion PDE. The choice made for diluteness in the high particle number limit will be that of the regime of constant mean free path. The programme of deriving the PDE in the case of constant mean free path is sometimes called a *kinetic* limit derivation.

In the special case of mass-independent diffusion rates, the kinetic limit derivation was carried out in 1980 by Lang and Nguyen [16], who followed the method of showing that the correlation functions between several particles are described by the BBGKY hierarchy which Lanford had introduced.

Introduced to the problem of generalizing Lang and Nguyen's derivation of the Smoluchowski PDE by James Norris, the author collaborated on it with Fraydoun Rezakhanlou. The principal aim of these notes is to give an informal but fairly detailed exposition of the kinetic limit derivation of the Smoluchowski PDE that was undertaken for dimension $d \geq 3$ in [10]. The treatment also first presents heuristic arguments with the aim that the reader may understand why the main theorem should be true before beginning a presentation of the proof of the theorem, and it also uses some novel probabilistic techniques to obtain key estimates used in the proof. The survey also touches on some related topics.

1.2 The Smoluchowski coagulation-diffusion PDE

We begin by recording the form of these equations and offering a brief explanation of the phenomenon that they may be expected to describe.

Let the dimension $d \geq 2$ be given. A collection of functions $f_n : \mathbb{R}^d \times [0, \infty) \rightarrow [0, \infty)$, $n \in \mathbb{N}$, is a strong solution of the discrete Smoluchowski coagulation-diffusion PDE with initial data $h_n : \mathbb{R}^d \rightarrow [0, \infty)$, $n \in \mathbb{N}$, if, for each $n \in \mathbb{N}$ and $x \in \mathbb{R}^d$, $f_n(x, 0) = h_n(x)$; and, for each $n \in \mathbb{N}$ and $(x, t) \in \mathbb{R}^d \times [0, \infty)$,

$$(1.2) \quad \frac{\partial f_n}{\partial t}(x, t) = d(n)\Delta f_n(x, t) + Q_1^n(f)(x, t) - Q_2^n(f)(x, t),$$

where the Laplacian acts on the spatial variable $x \in \mathbb{R}^d$. The final two terms are interaction terms, a gain term given by

$$(1.3) \quad Q_1^n(f)(x, t) = \frac{1}{2} \sum_{m=1}^{n-1} \beta(m, n-m) f_m(x, t) f_{n-m}(x, t),$$

and a loss term by

$$(1.4) \quad Q_2^n(f) = f_n(x, t) \sum_{m=1}^{\infty} \beta(m, n) f_m(x, t).$$

(When $t = 0$, the partial time derivative on the left-hand side in (1.2) is interpreted as a right derivative.)

Note that the equations have two sets of parameters: the diffusion rates $d : \mathbb{N} \rightarrow (0, \infty)$ and the coagulation propensities $\beta : \mathbb{N}^2 \rightarrow [0, \infty)$. The equations have a continuous counterpart, where the mass variable is now a positive real, and the above sums are replaced in an evident way by integrals, which we will not consider in this survey except in passing.

To interpret the solution, consider a large number of minute particles in space \mathbb{R}^d , each carrying an integer mass. In a similar manner to our opening discussion, the quantity $f_n(x, t)$ is interpreted as the density of particles of mass $n \in \mathbb{N}$ in the immediate vicinity of location $x \in \mathbb{R}^d$ at time $t \geq 0$. The form of the right-hand side (1.2) reflects the two dynamics for the particles: diffusive transport and binary coagulation. Particles of mass m diffuse at rate $2d(m)$, so that such a particle's displacement is given by $B(2d(m)t)$, $t \geq 0$, where $B : [0, \infty) \rightarrow \mathbb{R}^d$ is a standard Brownian motion. (The factor of two appears because the infinitesimal generator of standard Brownian motion is a *one-half multiple* of the Laplacian; when we call $d(n)$ the diffusion rate, this is thus strictly speaking a misnomer.) When a pair of particles are *microscopically* close, they may collide, disappearing from the model, to be replaced by a newcomer, whose mass is the sum of the two exiting particles'. The coagulation gain term (1.3) expresses the possible means by which a new particle of mass n may appear in the immediate vicinity of location x at time t : by the coagulation of some pair of particles of masses $(1, n-1)$, or $(2, n-2)$... or $(n-1, 1)$. The product form $f_m(x, t) f_{n-m}(x, t)$ in the interaction term reflects an assumption that the particles in the immediate vicinity of x are well mixed, and the coefficient $\beta(m, n-m)$ models the tendency of particles at close range of pair-type $(m, n-m)$ to coagulate in the immediate future. In the loss term (1.4), we see the means by which the density $f_n(x, t)$ may fall due to coagulation: a particle of mass n may drop out of the count for this density due to coagulation with another particle, and that other particle may have any mass $m \in \mathbb{N}$.

Our aim in this survey to explain how the system (1.2) may be derived in a kinetic limit from a collection of microscopic random models of diffusing mass-bearing particles that are liable to coagulate in pairs at close range. We now describe in precise terms the elements for this programme; for the case at hand, we are thus presenting an instance of the type of programme which we hazily sketched in our two opening examples. First, we specify the sequence of microscopic models: their initial particle distributions; and their dynamics: the free motion of individual particles, and the mechanism of pairwise coagulation at close range. In the main body of the article, we discuss only the derivation made in dimension $d \geq 3$, which was undertaken in [10]. Thus $d \geq 3$ may be assumed, except on one occasion when we make a short comment about the case when $d = 2$.

1.3 The microscopic models

The sequence of microscopic random models will be indexed by the total number N of particles initially present, at time zero. The N -indexed model will be specified by a probability measure \mathbb{P}_N . It is a measure not only on initial particle locations and masses but also on particle dynamics throughout $[0, \infty)$.

Initial particle distribution under \mathbb{P}_N . The quantity $f_n(x, 0) = h_n(x)$ may be interpreted as the density of particles of mass n in a tiny neighbourhood of $x \in \mathbb{R}^d$ at time zero. Thus, $\int_{\mathbb{R}^d} h_n(x) dx$ is interpreted as being proportional to the total number of particles of mass n and the constant $Z \in (0, \infty)$, which we define by $Z = \sum_{n \in \mathbb{N}} \int_{\mathbb{R}^d} h_n(x) dx$, as being proportional to the total number of initial particles.

We will index the time-zero particle set under \mathbb{P}_N by $[N] := \{1, \dots, N\}$; the initial mass and location of particle i will be denoted by $(x_i(0), m_i(0))$. Reflecting the above density interpretation, we choose $(x_i(0), m_i(0))$ independently, so that $(x_i(0), m_i(0))$ has density $Z^{-1}h_n(x)$ at $(x, n) \in \mathbb{R}^d \times \mathbb{N}$.

Notation for particle trajectories under \mathbb{P}_N . We wish to describe the subsequent evolution of each of the initial particles under \mathbb{P}_N . The trajectory of the i -th particle will be described by $(x_i, m_i) : [0, \infty) \rightarrow (\mathbb{R}^d \times \mathbb{N}) \cup \{c\}$, where here c is an element called a *cemetery* state whose role, which we will shortly describe in precise terms, is to house particles that have disappeared from the model due to being on the wrong side of a pairwise collision.

As such, at any given time $t \geq 0$, the particle configuration under \mathbb{P}_N is described by a map $[N] \rightarrow (\mathbb{R}^d \times \mathbb{N}) \cup \{c\}$, where $i \in [N]$ maps to $(x_i(t), m_i(t))$ (or to c).

To define the Markov process \mathbb{P}_N precisely, we will specify its Markov generator, which acts on test functions $F : \left((\mathbb{R}^d \times \mathbb{N}) \cup \{c\} \right)^{[N]} \rightarrow \mathbb{R}$. The action will be comprised of two parts: free motion of individual particles, and pairwise collision. We discuss our choice of each of these in words before providing the definition of the Markov generator.

Free motion. A particle of mass $n \in \mathbb{N}$ follows, independently of other particles, the trajectory $t \rightarrow B(2d(n)t)$ relative to its starting point, where B is a standard Brownian motion on \mathbb{R}^d .

Pairwise collision. Any two particles will be liable to collide when their locations differ by order ϵ . Here, $\epsilon = \epsilon_N$, the interaction range, is determined by N in a manner that we explain shortly.

We introduce a compactly supported smooth interaction kernel $V : \mathbb{R}^d \rightarrow [0, \infty)$ and a collection of microscopic interaction strengths $\alpha : \mathbb{N}^2 \rightarrow (0, \infty)$, and declare that, at time t , particles i and j collide at infinitesimal rate $\alpha(m_i, m_j)V_\epsilon(\cdot)$, where we adopt the convention that $V_\epsilon(\cdot) = \epsilon^{-2}V(\cdot/\epsilon)$. The argument \cdot/ϵ for V indeed entails that collision may occur only between particles whose locations differ by order ϵ ; the prefactor of ϵ^{-2} is introduced because, in dimension $d \geq 3$, once a pair of particles have approached to distance of order ϵ , they are liable to remain at such a displacement for order ϵ^2 of time, since their relative displacement evolves as a Brownian motion of rate $d(n) + d(m)$; thus, the role of this prefactor is to ensure that the proportion of instances of particle pairs approaching into the interaction range that result in collision is of unit order, uniformly in N . The role of the factor $\alpha(m_i, m_j)$ is to control whether this proportion is close to one for a given particle mass pair (which would be ensured by choosing the value of α in question to be high) or closer to zero.

The precise mechanism of collision. On collision of (x_i, m_i) and (x_j, m_j) at time t , each of the pair of particles disappears, to be replaced by a new particle of mass $m_i + m_j$ in the vicinity. As a matter of convenience for the ensuing proofs, the precise rule we pick for the appearance of the new particle is to choose its location to be x_i or x_j , with probabilities $\frac{m_i}{m_i+m_j}$ and $\frac{m_j}{m_i+m_j}$. This rule permits the interpretation that, when two particles collide, one survives the collision and the other perishes; the probability of survival is proportional to incoming particle mass; the particle surviving collects the mass of the perishing particle, and the perishing particle vanishes from space.

In a formal device, the perishing particle's location and mass are each sent to the cemetery state c , where they remain forever. As such, for each $i \in [N]$, the i -th particle's trajectory is described by setting the vanishing time $v_i \in [0, \infty]$ equal to the first time at which particle i experiences a collision in which it perishes. The trajectory is then given by $(x_i, m_i) \rightarrow \mathbb{R}^d \times \mathbb{N}$ on $[0, v_i)$ and $(x_i, m_i) = c$ on $[v_i, \infty)$.

The Markov generator of the dynamics. For any configuration $q \in \left((\mathbb{R}^d \times \mathbb{N}) \cup \{c\} \right)^{[N]}$, write I_q , the surviving particle set, for those $i \in [N]$ such that (x_i, m_i) lies in $\mathbb{R}^d \times \mathbb{N}$ (rather than equalling c). Let $F : \left((\mathbb{R}^d \times \mathbb{N}) \cup \{c\} \right)^{[N]} \rightarrow \mathbb{R}$ be smooth (in each hyperplane given by specifying the c -valued coordinates of the argument of F). Then the Markov generator \mathfrak{M} for \mathbb{P}_N is given as follows. For each $q \in \left((\mathbb{R}^d \times \mathbb{N}) \cup \{c\} \right)^{[N]}$, $\mathfrak{M}F(q) = \mathfrak{A}_F F(q) + \mathfrak{A}_C F(q)$, with the free-motion operator being given by

$$\mathfrak{A}_F F(q) = \sum_{i \in I_q} d(m_i) \Delta_{x_i} F(q),$$

where Δ_{x_i} is the d -dimensional Laplacian acting on F viewed as a function of $x_i \in \mathbb{R}^d$; and, recalling that $V_\epsilon(\cdot) = \epsilon^{-2}V(\cdot/\epsilon)$, with the collision operator being given by

$$\mathfrak{A}_C F(q) = \frac{1}{2} \sum_{i,j \in I_q} \alpha(m_i, m_j) V_\epsilon(x_i - x_j) \left[\frac{m_i}{m_i+m_j} F(S_{i,j}^1 q) + \frac{m_j}{m_i+m_j} F(S_{i,j}^2 q) - F(q) \right].$$

Here, $S_{i,j}^1(q)$, the configuration adopted in the event that particle i survives collision with particle j ,

is given by

$$(1.5) \quad S_{i,j}^1(q)(k) = \begin{cases} q(k) & \text{for } k \in [N] \setminus \{i, j\}, \\ (x_i, m_i + m_j) & \text{for } k = i, \\ c & \text{for } k = j, \end{cases}$$

while $S_{i,j}^2(q)$ is given by the same formula with the roles of i and j being reversed.

1.4 The regime of bounded mean free path and the choice of interaction range

It remains to specify how the interaction range ϵ is determined by total initial particle number N . This choice is made to be in the regime of constant mean free path: for dimension $d \geq 3$, $\epsilon = \epsilon_N$ will be chosen to satisfy

$$(1.6) \quad N = Z\epsilon^{2-d}.$$

To explain why this regime for the length of the free path is suitable, note that, since diffusion and coagulation terms are each present in the Smoluchowski PDE (1.2), we expect that the evolution of a typical particle will be determined both by its free motion and its collision with other particles. It will neither diffuse without collision nor collide repeatedly before diffusing a macroscopic distance.

The consideration that this regime be adopted forces the choice of scaling of ϵ as a function of N : picking a uniformly random particle index $i \in [N]$ at the outset, ϵ should be chosen so that the mean time to first collision of particle i converges as $N \rightarrow \infty$ to some strictly positive and finite constant.

A heuristic argument explains why (1.6) produces this outcome. We anticipate that, at any given time $t \geq 0$, a positive (although t -dependent) proportion of particles are surviving (rather than in the cemetery state). Assume that the surviving particles at time t are distributed so that the location and mass of each is chosen independently; the law of the location-mass statistic (x, n) of any given particle is equal to $f_n(x, t)$ (normalized to make the integral of this density equal to one). In other words, we are assuming a very strong sense that the density profile of particles under \mathbb{P}_N mimics the solution of (1.2).

Pick a particle uniformly at random at the initial time and call the selected particle the *tracer* particle. We would like to estimate the mean number of collisions suffered by the tracer particle during $[0, 1]$ in terms of N and ϵ . As we briefly discussed in the paragraph under the heading “pairwise collision” in the preceding section, this quantity is expected to have the same order as the number of other particles which enter the ϵ -neighbourhood of the given particle during $[0, 1]$. At any given time, our assumption on the distribution of other particles means that the probability that there is some other particle at distance less than ϵ from the tracer particle is of order $N\epsilon^d$. Thus, the mean total amount of time during $[0, 1]$ that some other particle is at distance less than ϵ from the tracer particle is also of order $N\epsilon^d$. Whenever another particle approaches the tracer particle to distance ϵ , it remains at the order of that distance for time of order ϵ^2 (since $d \geq 3$). Thus, the mean number of different particles which during $[0, 1]$ approach to within distance ϵ the tracer particle is of order $N\epsilon^d \cdot \epsilon^{-2} = N\epsilon^{d-2}$.

We thus see that imposing the relation (1.6) may indeed be expected to ensure that the mean number of collisions suffered by the tracer particle in unit time is bounded away from zero and infinity uniformly in N .

1.5 The recipe for the macroscopic coagulation propensities

The macroscopic coagulation propensities $\beta : \mathbb{N}^2 \rightarrow (0, \infty)$ that appear in the limiting system (1.2) depend in a non-trivial fashion on the microscopic parameters $V : \mathbb{R}^d \rightarrow [0, \infty)$, $\alpha : \mathbb{N}^2 \rightarrow (0, \infty)$ and $d : \mathbb{N} \rightarrow (0, \infty)$. Here is the recipe for obtaining $\beta(n, m)$ from these ingredients. As we will later explain, there exists a unique solution $u = u_{n,m} : \mathbb{R}^d \rightarrow (0, \infty)$ of the equation

$$(1.7) \quad -\Delta u_{n,m}(x) = \frac{\alpha(n, m)}{d(n) + d(m)} V(x) [1 - u_{n,m}(x)]$$

that satisfies $u_{n,m}(x) \rightarrow 0$ as $x \rightarrow \infty$. In fact, $0 \leq u_{n,m}(x) \leq 1$ for all $x \in \mathbb{R}^d$, and $u_{n,m}(x) = O(\|x\|^{2-d})$ as $x \rightarrow \infty$. (Here, as we will later, we write $\|\cdot\|$ for the Euclidean norm on \mathbb{R}^d .)

The quantities $\beta : \mathbb{N}^2 \rightarrow (0, \infty)$ in (1.2) are then specified by the formula

$$(1.8) \quad \beta(n, m) = \alpha(n, m) \int_{\mathbb{R}^d} V(x) (1 - u_{n,m}(x)) dx.$$

We mention that the minus sign appearing on the left-hand side of (1.7) was not used in the original treatment in [10]. A positive choice for $u_{n,m}$ permits an attractive probabilistic interpretation as we shall shortly discuss.

1.6 The weak formulation of the Smoluchowski PDE

We now recast the Smoluchowski PDE (1.2) in weak form, since it is to this form of the equations that we will prove convergence. To do so, let \mathfrak{J} be the space of sequences $J = \{J_n : n \in \mathbb{N}\}$ of smooth compactly supported functions $J_n : \mathbb{R}^d \times [0, \infty) \rightarrow [0, \infty)$. Then we say that $f = \{f_n : n \in \mathbb{N}\}$, with $f_n : \mathbb{R}^d \times [0, \infty) \rightarrow [0, \infty)$ measurable for each $n \in \mathbb{N}$, is a weak solution of (1.2) if, for each $J \in \mathfrak{J}$, it satisfies the formula obtained from (1.2) by multiplication by J_n , integration in space-time, and integration by parts. Namely, such an f solves (1.2) weakly if, for each $J \in \mathfrak{J}$ and $T \in (0, \infty)$,

$$(1.9) \quad \begin{aligned} & \int_{\mathbb{R}^d} \left(J_n(x, T) f_n(x, T) - J_n(x, 0) f_n(x, 0) \right) dx \\ &= \int_{\mathbb{R}^d \times [0, T)} \frac{\partial J_n(x, t)}{\partial t} f_n(x, t) dx dt \\ & \quad + \int_{\mathbb{R}^d \times [0, T)} \left(d(n) f_n(x, t) \Delta J_n(x, t) + (Q_1^n(f)(x, t) - Q_2^n(f)(x, t)) J_n(x, t) \right) dx dt. \end{aligned}$$

1.7 Empirical densities

In our opening discussion of the programme for deriving macroscopic limiting PDE, we suggested the use of δ -macroscopically smeared particle counts as candidates to approximate the limiting

evolution. Such counts play an important role in our derivation, and we will introduce them under the name *microscopic candidate densities* when we give an overview of the derivation of our main theorem, Theorem 1.1, in Section 4.

However, to state this theorem, we will not use them. Rather, we will use a close cousin, empirical density measures defined under the microscopic models \mathbb{P}_N . We now define these.

Under the law \mathbb{P}_N , let μ_N^f denote the \mathbb{P}_N -random variable, valued in measures on space-mass-time $\mathbb{R}^d \times \mathbb{N} \times [0, \infty)$ such that, for each $t \geq 0$, its time- t marginal $\mu_N^f(\cdot, t)$ is given by

$$\mu_N^f(\cdot, t) = \varepsilon^{d-2} \sum_{i \in I_q(t)} \delta_{(x_i(t), m_i(t))}.$$

For given $n \in \mathbb{N}$, let $\mu_{N,n}^f$ denote the \mathbb{P}_N -random variable, valued in measures on space-time $\mathbb{R}^d \times [0, \infty)$ such that, for each $t \geq 0$, its time- t marginal $\mu_{N,n}^f(\cdot, t)$ is given by

$$\mu_{N,n}^f(\cdot, t) = \varepsilon^{d-2} \sum_{i \in I_q(t)} \delta_{x_i(t)} \mathbf{1}_{x_i(t)=n}.$$

Let \mathcal{M} denote the space of measures μ on $\mathbb{R}^d \times \mathbb{N} \times [0, \infty)$ such that $0 \leq \mu(\mathbb{R}^d \times \mathbb{N} \times [0, T]) \leq TZ$, and note that μ_N^f is \mathbb{P}_N -a.s. valued in \mathcal{M} . Equip \mathcal{M} with the topology of vague convergence, under which this space is compact and metrizable.

1.8 Hypotheses on microscopic parameters

Our microscopic parameters are $d : \mathbb{N} \rightarrow (0, \infty)$, $\alpha : \mathbb{N}^2 \rightarrow [0, \infty)$ and $V : \mathbb{R}^d \rightarrow [0, \infty)$. In the original paper [10] and in the detailed overview of proof that we give in this survey, some hypotheses on these parameters must be imposed to enable the derivation to be made. We make some comments about the hypotheses made in [10] and then specify and discuss those we make here. The two sets of assumptions will be called the *original* and the *survey* assumptions throughout.

1.8.1 Original assumptions

On the diffusion rate and the microscopic interaction strengths. Suppose that there exists a function $\gamma : \mathbb{N}^2 \rightarrow (0, \infty)$ such that $\alpha(n, m) \leq \gamma(n, m)$ for all $(n, m) \in \mathbb{N}^2$, with γ satisfying

$$(1.10) \quad n_2 \cdot \gamma(n_1, n_2 + n_3) \cdot \max \left\{ 1, \left[\frac{d(n_2 + n_3)}{d(n_2)} \right]^{\frac{3d-2}{2}}, \left[\frac{d(n_2 + n_3)}{d(n_2)} \right]^{2d-1} \right\} \leq (n_2 + n_3) \cdot \gamma(n_1, n_2).$$

On the initial condition. A slightly technical assumption is needed, of the membership in local L^∞ space of some sums over n of certain averages of h_n ; the sums are weighted by certain powers of d and the above function γ . The reader may consult [10, Section 1] for the precise form. Although the assumption is a little technical to state, it is fairly weak: it is certainly satisfied if h_n is non-zero for only finitely many n , and each h_n is compactly supported with bounded supremum; and in fact numerous blowups in h_n are also permitted.

It is physically reasonable to think that the Brownian motion that is the free trajectory of the constituent particles in the models \mathbb{P}_N arises due to thermal agitation caused by many collisions with the constituents of an ambient environment of much smaller air molecules. Viewed in these terms, it is very natural to suppose that the diffusion rate $d(n)$ will decrease as a function of the mass. Accepting this, the assumption (1.10) is rather weak. If the diffusion rate is indeed decreasing, then (1.10) is satisfied provided that there exists a function $C : \mathbb{N} \rightarrow \infty$ for which $\alpha(n, m) \leq C(n)m$ for all $(n, m) \in \mathbb{N}$. Also, if the microscopic interaction strength α is identically constant, then the condition (1.10) is equivalent to the function $d(n)n^{1/(2-3d)}$ being non-increasing; when $d = 3$, then, we are permitted choices of $d(n)$ that grow as quickly as $n^{1/7}$.

1.8.2 Survey assumptions

These are:

On V . We will assume that $V : \mathbb{R}^d \rightarrow [0, \infty)$ is a compactly supported smooth function.

On $h_n : \mathbb{R}^d \rightarrow [0, \infty)$. There is a uniform bound on the L^∞ -norm of each of these functions, and each is supported in a given compact region of \mathbb{R}^d .

On $d : \mathbb{N} \rightarrow (0, \infty)$. The function $d : \mathbb{N} \rightarrow (0, \infty)$ is non-increasing and $\sup_{m \in \mathbb{N}} m^{-1}d(m)^{-d/2} < \infty$.

On $\alpha : \mathbb{N}^2 \rightarrow (0, \infty)$. We must suppose that $\sup_{(n,m) \in \mathbb{N}^2} \alpha(n, m) < \infty$.

Among these, the assumptions on the diffusion rates are genuinely restrictive: we must suppose that $d(m)$ grows less slowly than $m^{-2/d}$, which is not a particularly fast decay in any dimension $d \geq 3$. No such imposition was made in the original assumptions. The uniform bound demanded on $\alpha(\cdot, \cdot)$, it must also be admitted, is also a significant restriction. The survey assumptions will permit us however to offer a method of proof of key estimates needed for the main result which is largely self-contained, as well as being novel and very probabilistic in nature; since it serves our expository purpose, we have decided to accept the more limited domain of validity demanded by these assumptions.

1.9 Statement of main theorem

Here is our main result.

Theorem 1.1 *Let $d \geq 3$ and suppose that either of the above set of assumptions are in force. Let \mathcal{P}_N denote the law on \mathcal{M} given by the law of μ_N^f under \mathbb{P}_N ; recall that ϵ is related to N by means of the formula $N\epsilon^{d-2} = Z$, with the constant $Z \in (0, \infty)$ being given by the expression $Z = \sum_{n \in \mathbb{N}} \int_{\mathbb{R}^d} h_n$.*

The sequence $\{\mathcal{P}_N\}$ is tight in \mathcal{M} . Moreover, any limit point \mathcal{P} of the sequence $\{\mathcal{P}_N\}$ is concentrated on the space of measures taking the form $\sum_{n=0}^{\infty} f_n(x, t) dx \times \delta_n \times dt$ where $\{f_n : n \in \mathbb{N}\}$ ranges over weak solutions of (1.2) that satisfy the initial condition $f_n(\cdot, 0) = h_n(\cdot)$; recall that the collection of constants $\beta : \mathbb{N}^2 \rightarrow [0, \infty)$ is given by (1.8).

The reader may wonder what the meaning of this theorem is if it is not known that (1.2) has a weak (global in time) solution for the relevant parameter choices of $d(\cdot)$ and $\beta(\cdot, \cdot)$. In fact, the

method of proof furnishes the existence of at least one weak solution. In any case, Laurençot and Mischler [17] have established the existence of a global in time weak solution of (1.2) whenever $\lim_m m^{-1}\beta(n, m) = 0$, and $d(n) > 0$, for each $n \in \mathbb{N}$, conditions which are significantly weaker than those demanded by the theorem.

Theorem 1.1 describes the evolution of the density profiles of particles of various masses in the limit of large particle number by means of the Smoluchowski PDE, and in this way it realizes the derivation programme that we began this article by outlining, for the diffusive coagulating system in question. The derivation has the merit of being global in time. However, note that, in general, there are limitations in the description offered of the large scale behaviour of the system. If the weak solution of this system of PDE is not known to be unique, we merely demonstrate convergence in a subsequential sense to the space of solutions. For example, admitting the possibility that the system (1.2) has two distinct weak solutions $\{f'_n : n \in \mathbb{N}\}$ and $\{\hat{f}_n : n \in \mathbb{N}\}$ with initial condition $f_n(\cdot, 0) = h_n(\cdot)$, each of the following behaviours is consistent with Theorem 1.1:

- the empirical densities under the microscopic models \mathbb{P}_N may converge weakly to the solution $\{f'_n : n \in \mathbb{N}\}$ as $N \rightarrow \infty$ along the subsequence of even integers, and to $\{\hat{f}_n : n \in \mathbb{N}\}$ as $N \rightarrow \infty$ along the subsequence of odd integers;
- it may be that evolution of these densities is accurately approximated by flipping a fair coin, with the densities converging weakly to $\{f'_n : n \in \mathbb{N}\}$ as $N \rightarrow \infty$ should the outcome be heads, and to $\{\hat{f}_n : n \in \mathbb{N}\}$ as $N \rightarrow \infty$ should the outcome be tails.

These peculiar scenarios are excluded if uniqueness of solutions to (1.2) is known. Some conditions for uniqueness are furnished by [30, Proposition 2.6]; after deriving the kinetic limit of the PDE in [10], Fraydoun Rezakhanlou and the author in [11] provided uniqueness under rather weaker hypotheses. Indeed, as [11, Remark 1.2] discusses, the next proposition is a consequence of Theorems 1.1, 1.2, 1.3 and 1.4 of [11].

Proposition 1.2 *Let the dimension satisfy $d \geq 1$. For $a, b > 0$ such that $a + b < 1$, and for positive constants c_1 and c_2 , assume that $\beta(n, m) \leq c_1(n^a + m^a)$ and $d(n) \geq c_2 n^{-b}$ for all $n, m \in \mathbb{N}$. Also assume that $d : \mathbb{N} \rightarrow (0, \infty)$ is non-increasing. There exists $e > 0$ such that $\sum_n n^e \|h_n\|_{L^\infty(\mathbb{R}^d)} < \infty$ and $\|\sum_n n^e h_n\|_{L^1(\mathbb{R}^d)} < \infty$ imply that (1.2) has a unique weak solution.*

Note that the survey assumptions in fact imply the hypotheses of Proposition 1.2. This means that, in working with these assumptions, we automatically obtain the simpler statement of convergence available when uniqueness of the PDE system is known (and which we are about to state).

It is a simple corollary of Theorem 1.1 and Proposition 1.2 that convergence to (1.2) in fact holds in the following stronger sense.

Corollary 1.3 *Let $d \geq 3$ and suppose that the original assumptions, and the assumptions of Proposition 1.2, are in force. Let $J : \mathbb{R}^d \times [0, \infty) \rightarrow \mathbb{R}$ be a bounded and continuous test function. Then, for each $n \in \mathbb{N}$ and $T \in (0, \infty)$,*

$$(1.11) \quad \limsup_{N \rightarrow \infty} \mathbb{E}_N \left| \int_{[0, T)} \int_{\mathbb{R}^d} J(x, t) (\mu_{N, n}^f(dx, t) - f_n(x, t) dx) dt \right| = 0,$$

where again $N\epsilon^{d-2} = Z$, with $Z = \sum_{n \in \mathbb{N}} \int_{\mathbb{R}^d} h_n$. In (1.11), $\{f_n : \mathbb{R}^d \times [0, \infty) \rightarrow [0, \infty), n \in \mathbb{N}\}$ denotes the unique weak solution to the system of partial differential equations (1.2), with $\beta : \mathbb{N}^2 \rightarrow [0, \infty)$ again given by (1.8).

1.10 A simple computation about the collision of two particles

The basic mechanism of interaction in our model concerns a pair of particles. Here, we explain a brief computation concerning such a pair, which offers a probabilistic interpretation of the function $u_{n,m} : \mathbb{R}^d \rightarrow [0, 1]$ used in the recipe (1.8) for the macroscopic coagulation propensity β .

Suppose at a certain time, a particle of mass n is located at 0 and another, of mass m , is located at $x\epsilon$, where $x \in \mathbb{R}^d$. The pair are thus prone to interact shortly, in the next order ϵ^2 of time. Note also that, assuming uniform and independent placement of other particles in a compact region (in order to make an inference which we may find plausible for the actual model \mathbb{P}_N at any given time), the typical distance from a particle to the set of other particles is of order $N^{-1/d} = \epsilon^{1-2/d}$, which is far greater than the ϵ distance between the two particles in question. This means that in discussing the possible upcoming collision of this particle pair, we may harmlessly remove all other particles from the model.

Left with a two particle problem, we set $u_{n,m}^\epsilon(x)$ equal to the probability of subsequent collision of the pair. We may now use Brownian scaling, zooming in by a factor of ϵ^{-1} and slowing down time by a factor of ϵ^{-2} , to obtain a particle of mass n at the origin, one of mass m at x , with the trajectories $X_1, X_2 : [0, \infty) \rightarrow \mathbb{R}^d$ being Brownian motions of speeds $d(n)$ and $d(m)$, and collision occurring at rate $\alpha(n, m)V(X_1 - X_2)$. That is, $u_{n,m}^\epsilon(x)$ is independent of $\epsilon > 0$, and we may take $\epsilon = 1$.

As our notation suggests, $u_{n,m}^1$ is nothing other than $u_{n,m}$ from (1.7):

Lemma 1.4 *If $d \geq 3$, then $u_{n,m}^1$ is the unique solution $u_{n,m} : \mathbb{R}^d \rightarrow [0, 1]$ of (1.7).*

For occasional later use, we further define $u_{n,m}^{[t]} : \mathbb{R}^d \rightarrow [0, 1]$ for each $t > 0$ to be the probability that the two particles specified in the definition of $u_{n,m}^1(x)$ collide during $[0, t)$. Thus, $u_{n,m}^{[\infty]} = u_{n,m}^1$. In the expository discussion in Section 3 (though not for the proof of Theorem 1.1), we will need the next result.

Lemma 1.5 *Suppose that $d \geq 3$. Then $\|u_{n,m}^{[t]} - u_{n,m}^1\|_\infty \rightarrow 0$ as $t \rightarrow \infty$.*

We present the proofs of these two lemmas by using a more general notation which we now present.

1.11 Killed Brownian motion and the Feynman-Kac formula

In our two particle problem after scaling, the displacement of the particles performs a Brownian motion at rate $d(n) + d(m)$ until a random collision time. Slowing time by a factor of $d(n) + d(m)$, this process is Brownian motion *killed at rate* $\frac{\alpha(n,m)}{d(n)+d(m)}V$ in a sense we now explain.

Let $W : \mathbb{R}^d \rightarrow [0, \infty)$ denote a smooth and compactly supported function. Let $x \in \mathbb{R}^d$. Brownian motion in \mathbb{R}^d begun at x and killed at rate W is the stochastic process X that we now specify. The

process X maps $[0, \infty)$ into $\mathbb{R}^d \cup \{c\}$ where, as before, c is a formal cemetery state. To define X , let $B : [0, \infty) \rightarrow \mathbb{R}^d$ denote rate one Brownian motion with $B(0) = x$. Define its interaction until time t , I_t , to be equal to $\int_0^t W(B(s)) ds$. Let E denote an independent exponential random variable of rate one, and set the killing time $K_x \in [0, \infty]$ equal to $\inf \{t \geq 0 : I_t \geq E\}$, with the convention that $\inf \emptyset = \infty$. Then

$$X(s) = \begin{cases} B(s) & \text{for } s < K_x, \\ c & \text{for } s \geq K_x. \end{cases}$$

We say that killing occurs if $K_x < \infty$ and let $u_W : \mathbb{R}^d \rightarrow [0, \infty)$ be such that $u_W(x)$ is the probability that killing occurs.

Lemma 1.6 *For $d \geq 3$, $u = u_W$ is a solution of the modified Poisson equation*

$$(1.12) \quad -\Delta u(x) = W(x)(1 - u(x)).$$

satisfying $u \rightarrow 0$ as $x \rightarrow \infty$.

Remark. The solution is unique subject to $u \rightarrow 0$ as $x \rightarrow \infty$. In a formal sense, this is verified by observing that the difference v of two solutions solves $\Delta v = Wv$ on \mathbb{R}^d and then noting that

$$(1.13) \quad \int_{\mathbb{R}^d} \|\nabla v\|^2 dx = - \int_{\mathbb{R}^d} v \Delta v dx,$$

whose right-hand side is $-\int_{\mathbb{R}^d} Wv^2 dx$ and is thus at most zero. Hence, $\int_{\mathbb{R}^d} \|\nabla v\|^2 dx = 0$ and so ∇v is identically zero on \mathbb{R}^d . We thus see that v is a constant function, and, since $v \rightarrow 0$ as $\|x\| \rightarrow \infty$, v is identically equal to 0. This would prove uniqueness, except that (1.13) is a formal identity; if we integrate instead over the Euclidean ball B_R and take $R \rightarrow \infty$, then the boundary term in Green's theorem vanishes in the limit provided a decay condition on the solution u of (1.12) such as $\|x\|^{d-1} u(x) \nabla u(x) \rightarrow 0$ uniformly as $x \rightarrow \infty$ is made; thus, the solution is unique among those satisfying this decay condition. See [10, Section 6] for a proof of existence and uniqueness of the solution of (1.12) (subject only to $u \rightarrow 0$ as $x \rightarrow \infty$) that uses Fredholm theory and compactness arguments.

Proof. Let $v : \mathbb{R}^d \times [0, \infty) \rightarrow [0, 1]$ given by $v(x, t) = \mathbb{E} e^{-\int_0^t W(x+B(s)) ds}$, where the mean is taken over trajectories of standard Brownian motion $B : [0, \infty) \rightarrow \mathbb{R}^d$. The Feynman-Kac formula [24, Section III.19] shows that v satisfies the partial differential equation

$$(1.14) \quad \frac{\partial}{\partial t} v(x, t) = \Delta v(x, t) - W(x)v(x, t)$$

for $x \in \mathbb{R}^d$ and $t > 0$. Note that for any $s > 0$,

$$|v(t+s, x) - v(x, t)| \leq s \|W\|_\infty \sup_{r \in [t, t+s]} \mathbb{P}(x + B(r) \in \text{supp}(W));$$

as $t \rightarrow \infty$, this probability tends to zero uniformly in x , so that we find that $\frac{\partial}{\partial t} v(x, t) \rightarrow 0$ as $t \rightarrow \infty$, uniformly in $x \in \mathbb{R}^d$.

Note then that $1 - u_W(x)$, which is the probability that Brownian motion X begun at x and killed at rate W is never killed, is equal to $v(x, \infty)$. That u_W solves $-\Delta u(x) = W(x)(1 - u(x))$ in a distributional sense follows by taking a high t limit of (1.14), since v_t converges to 0 locally in L^1 . Since W is smooth, u being in local L^2 implies that Δu is also in this space; thus, u is locally in H^2 . Iterating, we find that in fact $u \in C^\infty$, and so u solves (1.12) in strong sense. \square

Proof of Lemma 1.4. Note that, by the spatial-temporal scaling satisfied by Brownian motion, $u_{n,m}^1$ equals u_W where $W = \frac{\alpha(n,m)V}{d(n)+d(m)}$. Hence Lemma 1.6 and the remark that follows it yield the result. \square

Proof of Lemma 1.5. Moreover, $u_{n,m}^{[t]}(x)$ equals $1 - v(x, t)$, and $u_{n,m}^1(x)$ equals $1 - v(x, \infty)$. Note that $|v(\infty, x) - v(t, x)|$ is at most the probability that Brownian motion begun at x visits the support of W after time t . With μ denoting d -dimensional Lebesgue measure, this probability is at most a constant multiple of $\mu(\text{supp}W) \cdot t^{-d/2}$, independently of $x \in \mathbb{R}^d$. \square

1.12 A guide to the rest of the survey

We have now set up the microscopic models \mathbb{P}_N and laid out the programme for deriving their macroscopic evolution, including our main result, Theorem 1.1. Our principal goal is to explain at a rather high, though not complete, level of detail, the proof of this theorem. We pause from pursuing this goal to explore two other directions first, however. First, in Section 2, we offer a glimpse of several topics which are tangentially related to this principal goal; this discussion is intended only to whet the reader's appetite for perhaps some of these topics and problems, and, for this reason as well as owing to limitations in the author's knowledge, it is brief and very inexhaustive. Second, in Section 3, we offer a leisurely heuristic overview of our kinetic limit derivation in a very simplified special case, of annihilating constant diffusivity Brownian particles on a torus with a translation invariant initial condition. The argument is not rigorous at each step here, and in its method it does not provide a template for the derivation of Theorem 1.1; rather, its main goal is to provide an intuitive explanation for the form (1.8) of the recipe for the macroscopic coagulation rates $\beta : \mathbb{N}^2 \rightarrow (0, \infty)$. The form of β is related to a repulsion effect present in the models \mathbb{P}_N at positive times, whereby the microscopic neighbourhood of a given particle is less likely to contain other particles than typical points that lie beyond this region but in a macroscopic vicinity. We thus hope that, at the end of Section 3, the reader will have a fuller understanding of why the statement of Theorem 1.1 is true, if not yet of how it was proved in [10].

We then return to the survey's principal goal. Section 4 explains how Theorem 1.1 will be proved, and the reader whose main interest is to see this proof explained may turn directly to this section. Therein, we introduce δ -smeared approximations to the particle densities defined in the microscopic models \mathbb{P}_N , called microscopic candidate densities. We state a fundamental estimate, the Stosszahlansatz, which expresses total coagulation propensity in \mathbb{P}_N approximately in terms of integrated products of microscopic candidate densities.

The next Section 5 describes the method of proof of the Stosszahlansatz. While so doing, it gives an alternative explanation to that of Section 3 for the form (1.8) of $\beta : \mathbb{N}^2 \rightarrow (0, \infty)$.

The actual proof of the Stosszahlansatz is given in Section 6. The proof relies on several estimates concerning various integrated sums of test functions over pairs and triples of particle indices. These bounds in turn are reduced to two key estimates. The first are particle concentration bounds, which state that if we have L^∞ -control at the initial time for the joint behaviour of k -tuples of particles in the models \mathbb{P}_N (with $k \in \mathbb{N}$ fixed, such as k equal to two or three), then this control propagates to all later times; it is here that the more restrictive aspect of the survey assumptions, on the decay rate of the diffusion rates $d : \mathbb{N} \rightarrow (0, \infty)$, is invoked. The second key estimate are bounds on killing probabilities u_W (that we introduced in Section 1.11), which are uniform over $W : \mathbb{R}^d \rightarrow [0, \infty)$ with given compact support. The proofs for this second key estimate also appear in this section.

Section 7 provides a proof of the first key estimate, the particle concentration bounds. The proof occupies several pages, but we hope that it is probabilistically interesting and intuitive.

Finally, in Section 8, we provide a summary of those points in our derivation where some steps were skipped, and mention where these omissions are treated in the original derivation in [10]. We also take this opportunity to explain how the proof in [10] is obscure at a certain moment, and highlight how in the present paper we have endeavoured to structure the arguments to shed light on this obscurity.

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2 A short foray into some neighbouring topics

2.1 The Smoluchowski coagulation ordinary differential equation

The spatially homogeneous analogue of the PDE we study, the Smoluchowski coagulation equation, may have been originally formulated in Smoluchowski's seminal work [25, (67)], and has been an object of attention for theoretical probabilists for a long time. The equation is now an ODE; it has a discrete and continuous mass version, as does the PDE. The solution $f : \mathbb{N} \times [0, \infty) \rightarrow [0, \infty)$ of the discrete form of this ODE may be written

$$(2.1) \quad \frac{d}{dt} f(n, t) = \frac{1}{2} \sum_{m=1}^{n-1} K(m, n-m) f(m, t) f(n-m, t) - f(n, t) \sum_{m=1}^{\infty} K(m, n) f(m, t).$$

In the continuous counterpart, the space of masses is now $[0, \infty)$ rather than \mathbb{N} , and the two sums are replaced by integrals with the natural ranges of integration. Existence and uniqueness results were obtained by McLeod [18], White [29] and Ball and Carr [3] in the discrete case, the latter also addressing the mass conservation of solutions. For such results in the continuous case, see McLeod [19], and Norris [20] who also shows an example of non-uniqueness.

Aldous' 1999 survey [1] discusses how special choices of the interaction kernel K have interesting probabilistic interpretations, in terms of point processes and related constructions on the complete

graph (with $K(x, y) = xy$, discrete), the uniform measure on large trees ($K(x, y) = x + y$, discrete), and on the continuum random tree ($K(x, y) = x + y$, continuous).

It is natural to pose the question of whether the ODE may be derived from a sequence of random models with diverging initial particle number. In such a model, the analogue of the microscopic law \mathbb{P}_N in the programme we discuss would consist of a probability measure under which N particles carry integer or non-negative real valued masses, and any pair coalesces at infinitesimal rate $K(\cdot, \cdot)$, where the arguments are the masses of the concerned pair; at the moment of their coalescence, the two particles leave the system, to be replaced by a newcomer whose mass is the sum of the departing pair's. In [20, 21], Norris derives the Smoluchowski ODE from this model, considers more general mechanisms for coalescence and carries out corresponding derivations for them.

In the case where K is identically constant, this random system of coalescing particles is called Kingman's coalescent [13, 14]. Although we have defined it only with an initial condition with a finite number of particles, the system *comes down from infinity*, in the sense that there is a well defined stochastic process S that maps $(0, \infty)$ to \mathbb{N} with the property that, for any $t > 0$, the process $[0, \infty) \rightarrow \mathbb{N} : s \rightarrow S(t + s)$ has the distribution of the process of the number of surviving particles in Kingman's coalescent given that the initial number is $S(t)$.

See [4] for a recent survey of coalescence theory, including a treatment of Kingman's coalescents, a more general class of coalescents, Λ -coalescents, in which several particles may combine simultaneously, spatial models and their applications to population genetics.

2.2 Annihilating random walkers on \mathbb{Z}^d

Suppose that at time zero, a finite collection of walkers are located, each at some site in \mathbb{Z}^d ; there may be several walkers at any given site. Each walker performs a continuous time simple random walk, staying at her present site for a duration which is exponentially distributed, of mean one, independently of other decisions, and then jumping with equal probability to one of her $2d$ neighbours. For any instance of a pair of walkers occupying the same site at any given moment of time, both walkers in the pair are annihilated at exponential rate one. The description is informal and we do not provide a precise formulation here.

2.2.1 Many walkers initially: coming down from infinity

These models have some interesting parallels with coagulating diffusive systems. Defining this model on a singleton set, where all walkers must occupy the same site, note that the model reduces to Kingman's coalescent, which, begun with infinitely many particles, has at any positive time only finitely many. Does this phenomenon also take place in \mathbb{Z}^d ? A natural starting condition for the model on \mathbb{Z}^d is to begin with N walkers, all located at the origin, and consider high N behaviour. In [2], it is shown that, in contrast to the non-spatial case, infinitely many walkers survive asymptotically: if there are N initially, there are of the order of $\log^* N$ at any given positive time, where the function $\log^* N$ in essence denotes the number of iterations of the logarithm which reduces the value of N to below zero. Since the model asymptotically manufactures an arbitrarily large number of surviving particles, however slowly it does so in the high N limit, it is a natural extension of the derivation of Theorem 1.1 to enquire as to whether a counterpart result holds for

this model. We must expect the recipe (1.8) for the macroscopic coagulation rates to be altered so that a discrete Laplacian appears instead.

2.2.2 The model on a discrete torus: Smoluchowski and Kingman

Similarly to the model on \mathbb{Z}^d , we may consider the model of annihilating random walkers on the discrete torus \mathbb{T}_N^d , formed by quotienting \mathbb{Z}^d by $N\mathbb{Z}^d$. Imagine for the sake of simplicity that, at time zero, the walkers sit one apiece at each of the N^d sites of \mathbb{T}_N^d . In the first few units of time, a positive fraction of walkers are annihilated, and the system becomes sparser. Writing $s_N : [0, \infty) \rightarrow [0, 1]$ for the proportion of surviving particles in \mathbb{T}_N^d , we may define $s : [0, \infty) \rightarrow [0, 1]$, $s = \lim_N s_N$, (an almost sure limit which clearly exists and is non-random), and ask questions about the rate at which $s(t) \searrow 0$ as unscaled time t tends to infinity.

On the other hand, by scaling space and time, we may find a regime analogous to that identified in the principal aim of this survey. Scaling \mathbb{T}_N^d to the unit torus by squeezing space by a factor of N , a respect for Brownian scaling leads us to speed time up by a factor of N^2 . In the new coordinates, there is an instantaneous scourge of walkers that reduces their number from N^d to an order of N^{d-2} (at any given positive time). Note that the relation (1.6), that $n\epsilon^{d-2}$ is of unit order, with n the surviving particle number, is satisfied in the sense that n is order N^{d-2} and ϵ , the interaction range, is N^{-1} . In scaled coordinates then, the system crashes down from infinity, and naturally slows down into the regime of constant mean free path. Our translation invariant choice of initial condition should manifest itself in this regime by convergence of the surviving particle number to a solution of the Smoluchowski coagulation ODE (2.1); since particles are indistinguishable, the kernel K is identically equal to one.

At time scales beyond the N^2 rescaling, surviving particles cross the torus many times between collisions. If we choose a speeding up of time by a factor of N^d , then we enter a regime where only finitely many particles survive. Indeed, Cox [6] proved that, when dimension $d \geq 3$, the rescaled process of surviving particle number, $[0, \infty) \rightarrow \mathbb{N}$, $t \rightarrow s_N(2N^d t/G)$, converges in law to the process of surviving particle number in Kingman's coalescent. The constant factor G is the mean total amount of time that a continuous-time simple random walk in \mathbb{Z}^d begun at the origin spends at the origin in all positive time. This regime is one where a finite population of walkers each mixes spatially at an infinite rate, thus becoming indistinguishable; the presence of the factor G is explained by noting that the time of any pair of such particles to meet should be gauged in a clock which advances at a speed which is double (since a displacement between two walks is considered) that at which the walker on \mathbb{Z}^d at late time encounters previously unvisited vertices at unit rate. Cox also studied the problem in dimension two, and noted the implications of the solution for voter model consensus times.

2.3 The elastic billiards model and Boltzmann's equation

Our discussion draws heavily on Chapter 1 of Villani's review [28] of collisional kinetic theory.

2.3.1 The form of the equations

In Subsection 1.1.2, we mentioned Lanford's 1975 derivation for short times of Boltzmann's equation (1.1) in a kinetic limit from a system of elastically colliding billiards. In the form of the equation suitable for such a billiard model in dimension $d = 3$, the collision operator in (1.1) is given by

$$(2.2) \quad Q(f, f)(t, x, v) = C \int_{\mathbb{R}^d} dv_* \int_{S^{d-1}} d\sigma |v - v_*| \left(f'(t, x, v') f'_*(t, x, v'_*) - f(t, x, v) f_*(t, x, v_*) \right),$$

where $C \in (0, \infty)$ is a constant. To explain the notation, consider a collision that a sphere of velocity $v \in \mathbb{R}^d$ may undergo. The particle with which it collides has some velocity $v_* \in \mathbb{R}^d$. Impact may occur over a hemisphere in a surface of the velocity v sphere. Denoting the outgoing velocities of the two spheres by v' and v'_* , conservation of momentum and kinetic energy imply that

$$\begin{cases} v' + v'_* = v + v_* \\ |v'|^2 + |v'_*|^2 = |v|^2 + |v_*|^2. \end{cases}$$

The form of the outgoing velocities is determined by the angle of impact. The possibilities may be parameterized as follows:

$$\begin{cases} v' = \frac{v+v_*}{2} + \frac{|v+v_*|}{2} \sigma \\ v'_* = \frac{v+v_*}{2} - \frac{|v+v_*|}{2} \sigma, \end{cases}$$

as σ varies over S^{d-1} .

We mention in passing one notable feature of (1.1): the term $|v - v_*|$, which is the *Boltzmann collision kernel*, and which in a more general setting may depend non-trivially on σ , has no such dependence in the present case of elastic collisions and dimension $d = 3$. The reason for this is as follows. Imagine a torch shone from far away on a billiard ball floating in space. The torch's rays fall on one hemisphere of the billiard ball and are then reflected (elastically, as if the boundary of the ball were perfectly smooth and reflective). Consider the illumination provided by the reflected rays alone. These rays travel in every direction in S^2 , and, as a simple exercise verifies, their intensity is independent among all these directions.

The collision operator $Q(f, f)$ may be written as a difference of non-negative gain and loss terms, $Q^+(f, f) - Q^-(f, f)$, by splitting (2.2) across the minus sign in the big bracket. We obtain a phenomenological description of (1.1) akin to that offered for the Smoluchowski PDE appearing after the equations in Section 1.2. Particles of velocity v near location x are subject to collision, and contribute to the loss term $Q^-(f, f)$; collisions of particles with other velocity pairs may occur which produce new velocity v particles near x , as the gain term $Q^+(f, f)$ records.

2.3.2 Loschmidt's paradox and the Stosszahlansatz

A system of elastically colliding billiards in a box is at equilibrium a reversible continuous-time Markov chain. However, Boltzmann's equation begun from generic initial data does not share this reversibility. Indeed, Boltzmann's H theorem shows that the Boltzmann H functional,

$$H(f) = \int_{(x,v) \in \mathbb{R}^d \times \mathbb{R}^d} f \log f,$$

satisfies $\frac{d}{dt}H(f(t, \cdot, \cdot)) \leq 0$. The quantity H may be viewed as a measure of information; information dissipates monotonically as time evolves, in accordance with the second law of thermodynamics. At some late time, this rate of dissipation may slow as the system approaches equilibrium. However, for generic initial data for Boltzmann's equation, H may decrease in a strictly monotonic fashion as time advances. This irreversible property of the macroscopic evolution seems to be in tension with the reversible nature of the basic collision event that two billiards may undergo revealed by reversing in time a viewing of the collision. Concerns such as these caused Boltzmann's claim that the equation offered an accurate macroscopic description of classical many body systems such as elastic billiards to be treated with much scepticism. Loschmidt found a paradox which brought these concerns into a sharper relief. Accepting that a large system of elastic billiards is accurately modelled by Boltzmann's equation for all typical choices of initial data, begin with some such data and run the deterministic dynamical rules for the billiard system for some fixed time t . The density profiles will approximately follow the solution of Boltzmann's equation, and the H functional will drop from its initial value. Stop the evolution at time t and then reverse the velocity of each particle, leaving each particle's location unchanged. Then run the system for a further t units of time. Clearly the resulting evolution will be a rerun of the dynamics we just witnessed in the sense of reversed time. At the end of this second dynamics, the collection of billiards has its original set of locations, with reversed velocities. Note however that during this second dynamics, the H functional was rising, not falling. However, this is impossible for a system which is accurately approximating a solution of Boltzmann's equation.

Loschmidt's paradox indicates that not all microscopic data consistent with a given macroscopic density profile may result in an evolution for which Boltzmann's equation is an accurate model. The velocity-reversed time- t particle data is a counterexample to the hypothesis that Boltzmann's equation may be so derived from all such microscopic data. However, there is no contradiction to the hypothesis that all but a tiny fraction of particle configurations approximating a given density profile begin a dynamics whose evolution is accurately described by Boltzmann's equation.

The paradox also has implications for methods of proof that may be proposed for deriving Boltzmann's equation from microscopic models. In Boltzmann's own derivation, he invoked an assumption of independence on the part of colliding particles, which he called the *Stosszahlansatz*, or the *collision number hypothesis*. This asserts roughly that, in the neighbourhood of a location x at any time t , the distribution of the numbers of collisions of particles of two given velocities, v and v_* , in a many body system of elastic billiards, is accurately specified by knowing the densities (macroscopically denoted by $f(x, v, t)$ and $f(x, v_*, t)$) of velocity v and v' particles near x at time t ; the particles' histories until this time does not significantly bias the local structure of these families of particles away from that of a Poissonian system of such particles at these two densities; so that, for example, the rate of collision of a randomly picked particle of velocity v near x at time t with a particle of velocity v' , and the distribution of the impact parameter on collision, are accurately modelled by the Poisson systems at these densities. Loschmidt's paradox indicates a subtlety about the *Stosszahlansatz*. It may be valid for precollisional particle velocities, but in cannot be for postcollisional ones; for the latter, the history of the concerned particles has a great deal that biases their distribution from a Poissonian cloud model for the two velocity types. In other words, the mechanism of elastic collision may propagate chaos, taking independent randomness present at

an initial time and preserving it at given later times, but the chaos propagated is one-sided, not double-sided, referring to statistical inferences about the particles' future, and not their past.

In our more humble setting of coagulating Brownian particles, a key role is played by a result, Proposition 4.1, concerning collision propensity for the microscopic models, which we interpret as the Stosszahlansatz, as we will see in Section 4. However, the random and reversible nature of the free motion of the individual particles means there is no analogue of Loschmidt's paradox and no need to formulate the Stosszahlansatz as a statement concerning merely one-sided, rather than double-sided, chaos.

2.4 Gelation and mass conservation

2.4.1 Mass conservation for the Smoluchowski PDE

The collision event in the microscopic models \mathbb{P}_N conserves mass. How does mass conservation manifest itself macroscopically, in a solution of the Smoluchowski PDE? For a solution $\{f_n : n \in \mathbb{R}^d\}$ of (1.2), we may interpret $M_n(t) := n \int_{\mathbb{R}^d} f_n(x, t) dx$ as the total mass among particles of mass $n \in \mathbb{N}$ at time $t \in [0, \infty)$, and thus $M(t) := \sum_{n \in \mathbb{N}} M_n(t)$ to be the cumulative mass of particles at this time. For any $T > 0$, a solution of (1.2) is said to conserve mass during $[0, T]$ if $M(s) = M(0)$ for all $s \in [0, T]$. The passage from the microscopic to the macroscopic might lead one to expect solutions to be mass conserving on all such intervals. In fact, only the inference that $M : [0, \infty) \rightarrow [0, \infty)$ is non-increasing is readily available. We may define then the gelation time $t_{\text{gel}} \in [0, \infty]$, $t_{\text{gel}} = \inf \{t \geq 0 : M(t) < M(0)\}$, with $\inf \emptyset = \infty$. It is shown in [11] that the unique weak solution of the PDE which Proposition 1.2 provides is mass conserving in the sense that $t_{\text{gel}} = \infty$. Certainly under the survey assumptions, the resulting solution of the PDE satisfies the hypotheses of Proposition 1.2, and so is mass conserving. Indeed, this is true in every circumstance under which a kinetic limit derivation of the Smoluchowski PDE has been carried out.

2.4.2 The meaning of gelation for the microscopic models

Nonetheless, it is natural to ask what behaviour we would expect to see under the laws \mathbb{P}_N in a system which converges to a solution of the Smoluchowski PDE with gelation. After the gelation time t_{gel} , a positive fraction of the initial particle mass in a high- N indexed model \mathbb{P}_N will be present in particles above any given $K \in \mathbb{N}$; this fraction is independent of the value of K , although, for high values of K , we will have to increase N in order to witness this effect. A gel, composed of super-heavy particles, is forming microscopically beyond the gelation time.

Does this phenomenon actually take place in a model \mathbb{P}_N for some choice of parameters V , $d(\cdot)$ and $\alpha(\cdot, \cdot)$, or in some variant of this model?

To prepare to answer this, we first discuss a natural extension to our definition of microscopic model. Under \mathbb{P}_N , all particles have an equal interaction range $\epsilon = \epsilon_N$, irrespective of their mass. It is natural to introduce a mass-dependent interaction range, of the form $r_n \epsilon$ for particles of mass n ; presumably r_n would be increasing, and the choice $r_n = n^{1/d}$ would correspond to solid ball particles composed of a common material which instantaneously merge to form a larger such ball on collision. Other choices $r_n = n^\chi$, for $\chi \in [1/d, 1]$, may be possible, corresponding to fractal

geometries for the internal particle structure (as we will discuss in Section 2.6). Without the assumption of additional and non-local attractive inter-particle forces, it is hard to see, however, how a choice of the form $\chi > 1$ would be physically meaningful. The choice $\chi = 1$ is already a little beyond the border of the plausible realm: in a farfetched effort to justify this choice, we may model each particle as a long and very thin bar, and imagine that each bar rotates rapidly and chaotically about its centre of mass, while diffusing on a slower time-scale; when two bars touch, they instantaneously and rigidly join. Because of their rapid rotation, this will tend to happen when they are closely aligned, so that the new particle also resembles a long thin bar.

Whatever the physically reasonable range of choices for radial parameters $\{r_n : n \in \mathbb{N}\}$ may be, the natural adaptation of particle dynamics when they are introduced is a change in the pairwise collision rule discussed in Section 1.3. Where before particles x_i and x_j of mass n and m coagulate at rate $\alpha(n, m)\epsilon^{-2}V((x_i - x_j)/\epsilon)$, we now stipulate that this rate is $\alpha(n, m)\epsilon^{-2}(r_n + r_m)^{-2}V(\frac{x_i - x_j}{(r_n + r_m)\epsilon})$; the presence of the term $(r_n + r_m)^{-2}$ allows the microscopic coagulation propensity $\alpha(n, m)$ to retain its interpretation of determining the proportion of particle pair overlaps that lead to coagulation (uniformly as the masses of the pair vary).

The kinetic limit derivation of Theorem 1.1 is undertaken after these changes are made by Rezakhanlou [23]. When $d \geq 3$ and the relation $r_n = n^\chi$ is imposed, the derivation is made when $\chi \in [0, 1/(d - 2))$. The macroscopic coagulation rates $\beta : \mathbb{N}^2 \rightarrow (0, \infty)$ are then found to satisfy

$$\beta(n, m) \leq C(d(n) + d(m))(r_n + r_m)^{d-2} \cdot \text{Cap}(\text{supp}V),$$

where the latter term denotes the Newtonian capacity of the support of V .

When dimension d equals three, and we suppose, very reasonably, that $\sup_{n \in \mathbb{N}} d(n) < \infty$, we see that $\beta(n, m) = O(n + m)$ whenever $r_n = O(n)$. In such a regime for β , it is reasonable to believe that the Smoluchowski PDE is mass conserving for all time; indeed, Proposition 1.2 comes close to showing this if $d(\cdot)$ decreases gradually enough.

We may tentatively conclude then that the perturbation of our model which includes radial dependence of particles without making more profound changes to inter-particle interaction is not a suitable physical context to study the phenomenon of gelation.

2.4.3 A weaker notion of gelation: an analogue of weak turbulence

A weaker notion of solution blowup than finite gelation time is the condition that

$$(2.3) \quad \int_0^\infty \int_{\mathbb{R}^d} m^r f_m(x, t) dx dm \rightarrow \infty \text{ as } t \rightarrow \infty,$$

for some $r > 1$. In [31, Appendix], an analogy is drawn between the non-linear Schrödinger equation and the Smoluchowski PDE under which, in the case of the cubic defocussing NLS, the criterion above corresponds to weak turbulence. The condition (2.3) corresponds to ongoing coagulation under which a positive fraction of mass reaches arbitrarily high mass nodes at sufficiently late time. It is argued non-rigorously in [31] on the basis of scaling considerations for the PDE that, modelling $\beta(n, m) = n^\eta + m^\eta$ and $d(n) = n^{-\phi}$, the behaviour (2.3) is not expected to occur provided that $\eta + \phi < 1$.

2.5 The kinetic limit derivation when $d = 2$ and with other variants

In [9], the kinetic limit derivation of the PDE from the models \mathbb{P}_N was undertaken in dimension $d = 2$. We mention here the key changes needed in the models \mathbb{P}_N , and the changes in the recipe for determining $\beta : \mathbb{N}^2 \rightarrow (0, \infty)$ from the microscopic parameters. We will also return to the discussion of case $d = 2$ in Section 8, in order to discuss how the proofs in this case differ from when $d \geq 3$.

Interaction range. The relation (1.6) becomes $N|\log \epsilon|^{-1} = Z$, for a given constant $Z \in (0, \infty)$. The interaction range is now exponentially small in N , far smaller than it was in the case $d \geq 3$. It is the same regime of constant mean free path that dictates the scale, but now particles are readily available to each other due to Brownian recurrence; small interaction range acts as a countervailing effect.

Pairwise collision rule. The infinitesimal rate of coagulation between two particles of mass n and m located at x_i and x_j is now taken to be $\alpha(n, m)\epsilon^{-2}|\log \epsilon|^{-1}V\left(\frac{x_i - x_j}{\epsilon}\right)$, for a collection of microscopic interaction strengths $\alpha : \mathbb{N}^2 \rightarrow (0, \infty)$. The change from the case $d \geq 3$ is the appearance of the factor $|\log \epsilon|^{-1}$. Its role to preserve the interpretation of $\alpha(\cdot, \cdot)$ as specifying the proportion of particle overlaps leading to coagulation: were it absent, Brownian recurrence would offer overlapping particles endless opportunities to coagulate, and the proportion of coagulation would be one, for any positive value of α .

The recipe for the macroscopic coagulation rates. With a choice of compactly supported interaction kernel $V : \mathbb{R}^2 \rightarrow [0, \infty)$ for which $\int_{\mathbb{R}^2} V = 1$, the formula (1.8) becomes

$$(2.4) \quad \beta(n, m) = \frac{2\pi \cdot (d(n) + d(m)) \cdot \alpha(n, m)}{2\pi \cdot (d(n) + d(m)) + \alpha(n, m)}.$$

Thus, the nature of V is manifest in the macroscopic evolution only through the value of its L^1 norm. The reason for this is that, having accepted the presence of a new factor of $|\log \epsilon|^{-1}$ into the formula for $\alpha(\cdot, \cdot)$, any overlapping pair of particles are now not likely to coagulate in any particular excursion into each other's interaction range of duration of order ϵ^2 . Rather, many such opportunities to visit occur for the particles before they move to a large distance from one another, and one among these many visits may cause collision. During all the visits, the details of the form of V no longer really matter, except in a weak large of large numbers' sense, where the average rate of interaction is determined by the L^1 norm of V .

2.6 Diffusion and coagulation: two effects from one random dynamics

In our microscopic models, the Brownian motion that is the free motion of individual particles is simply a definition, as is the binary coalescence mechanism. Might it be possible to find a microscopic model in which these two phenomena emerge from one microscopic description? Here are two possible answers.

2.6.1 Physical Brownian motion

Physically, Brownian motion arises by the thermal agitation of a particle caused by many random collisions with its neighbours, in a similar vein to the way that the heat equation is expected to arise as a macroscopic evolution equation which we discussed in this survey's opening paragraphs. A physically natural but mathematically presumably intractable microscopic model might suspend comparatively large spheres in an ambient environment of much smaller particles, with a dynamic of elastic collision, and a mechanism of sticking of the large particles on mutual contact. In this sense, the Smoluchowski PDE is sometimes called a model of a colloid. Regarding the important question of the physical derivation of Brownian motion, we mention the recent advance [8], concerning the long time behaviour of a tracer billiard in a system of elastic billiards, in a dilute, constant mean free path limit.

2.6.2 Random walker clusters

A less classical but probabilistically interesting model is the following. Consider a Markov chain whose state space consists of a finite collection of *occupied* sites in \mathbb{Z}^d . Think that each site is occupied by one walker. Each walker decides to make a transition at the ring times of independent Poisson rate one clocks. Any given walker's transition takes place instantaneously. If a walker is selected to make a transition, he may not move – his transition is in place – if his removal from the lattice disconnects a connected component of occupied sites in the nearest neighbour structure. Otherwise, the walker considers making a uniformly random move to one of the nearest-neighbour or diagonally adjacent sites. He does so if the move is to an unoccupied site and the move does not disconnect any two occupied sites that were connected before the move. Otherwise, he stays in place.

Suppose for a moment that initially the occupied sites are nearest-neighbour connected. The rules are rigged so that this remains so at later times. The ergodicity of the system indicates that the centre of mass of the connected component diffuses in the long term. It is natural to pose the question as to how the diffusivity depends on the mass. Anyway, we obtain a collection of mass-dependent diffusion rates $d : \mathbb{N} \rightarrow (0, \infty)$, where now mass means the number of occupied sites in the cluster.

Suppose instead that we begin with a collection of comparatively well separated pairs of nearest neighbours. Each pair begins a random journey which in the large is Brownian. When two clusters meet, they combine, and never break apart subsequently.

All in all, then, it would seem that with a suitable initial condition and a parabolic scaling of space-time, the model may converge to a solution of the Smoluchowski PDE for some choice of its parameters. Note that the appropriate form of the equation may include the mass-dependent interacting range which we discussed in Subsection 2.4.2. One may speculate that r_n should be chosen to scale as $n \rightarrow \infty$ according to the scaling satisfied by the typical diameter of an isolated cluster with n occupied sites at equilibrium. Presumably the cluster has a fractal structure that contributes an exponent of the form $r_n = n^{\chi(1+o(1))}$.

3 Homogeneously distributed particles in the torus

We now study the convergence of the microscopic models \mathbb{P}_N to the limiting system (1.2) in a very particular special case. The choice is made so that, while most of the technical subtleties of definition and proof in the convergence are eliminated, the recipe (1.8) for the macroscopic coagulation propensities will be maintained. The main aim of our study of the special case is to explain why the relation (1.8) holds; intimately tied to this is a certain microscopic repulsion experienced by the particles at positive macroscopic times, which we also take the opportunity to discuss. Despite these various simplifications, our discussion here is heuristic, with the derivation of several intuitively plausible steps only sketched or omitted entirely; the model in this section is a special case of the annihilating system studied by Sznitman [26], and the reader may wish to refer there for a rigorous treatment.

In the special case, under the microscopic models \mathbb{P}_N , there will initially be particles of only of unit mass, and each will diffuse at unit rate. Moreover, as time evolves and particles collide in pairs, the concerned particles will disappear, without the appearance of any new particle. Thus we discuss an interaction of annihilation rather than coagulation. The particles will initially be placed not in \mathbb{R}^d but rather in the unit d -dimensional torus, each placed independently and uniformly with respect to Lebesgue measure. This choice forces the whole dynamics of \mathbb{P}_N to be invariant under any given spatial translation.

Let \mathbb{T}^d denote the d -dimensional unit torus, namely the quotient of \mathbb{R}^d by \mathbb{Z}^d , or the unit cube in \mathbb{R}^d with periodic boundary conditions. In this section, \mathbb{P}_N refers to the annihilating particle dynamics described in the preceding paragraph; we write $\alpha = \alpha(1, 1)$ for the microscopic interaction strength of the single particle mass pair in question. In the formal language specifying the Markov generator that we saw for our main object of study in Section 1.3, we are instead setting the free-motion operator equal to $\mathfrak{A}_F F(q) = \sum_{i \in I_q} \Delta_{x_i} F(q)$, and the collision operator is equal to $\mathfrak{A}_C F(q) = -\frac{\alpha}{2} \sum_{i, j \in I_q} \epsilon^{-2} V\left(\frac{x_i - x_j}{\epsilon}\right) F(q)$.

We also make a further minor simplification, which is to choose the constant Z in (1.6) equal to one.

The task of making the kinetic limit derivation in this case is to explain how it is that statistics summarising the densities of particles in the microscopic model converge to the appropriate macroscopic evolution, which in this case is given by a function $f : \mathbb{T}^d \times [0, \infty) \rightarrow [0, \infty)$ satisfying the PDE

$$(3.1) \quad \frac{\partial}{\partial t} f(x, t) = \Delta f(x, t) - \beta f(x, t)^2$$

with initial condition $f(x, 0) = 1$ for all $x \in \mathbb{T}^d$. One simplification in our analysis is readily apparent: the initial condition has no dependence on the spatial parameter, and this property will be maintained in time. So we may define $h : [0, \infty) \rightarrow [0, \infty)$ by setting $h(t) = f(x, t)$ for any choice of $x \in \mathbb{T}^d$ and thereby recast (3.1) as an ordinary differential equation

$$(3.2) \quad \frac{d}{dt} h(t) = -\beta h(t)^2,$$

with $h(0) = 1$.

In what sense is the evolution of the microscopic models \mathbb{P}_N approximately but accurately summarised by the ODE (3.2)?

The spatial homogeneity present in the special case in question offers a simple form for the answer to this question. We introduce a *microscopic candidate density* $h_N : [0, \infty) \rightarrow [0, \infty)$, a quantity which summarises the density of particles in the microscopic model \mathbb{P}_N and which we hope to show approximates $h : [0, \infty) \rightarrow [0, \infty)$ when N is high. We set $h_N(t) = N^{-1}\mathbb{E}s_N(t)$, where $s_N(t)$ is the number of *surviving* particles at time t , namely, the mean number of particles which have not been annihilated before time t . (In the general, spatially inhomogeneous, setting, we will also define a microscopic candidate density, but its definition will be a little more involved, it will be random rather than deterministic, and it will depend not only on the time variable but also on the macroscopic location.)

In the special case, our kinetic limit derivation amounts to explaining how it is that $h_N : [0, \infty) \rightarrow [0, \infty)$ converges as $N \rightarrow \infty$ to the unique solution $h : [0, \infty) \rightarrow [0, \infty)$ of (3.2). The principal aim of this section is to justify heuristically the relation (1.8) between the quantity β appearing in (3.2) and the microscopic parameters. In the present case, this relation takes the following form.

Proposition 3.1 *Assume that $V : \mathbb{R}^d \rightarrow [0, \infty)$ is continuous and compactly supported. Then the functions $h_N : [0, \infty) \rightarrow [0, \infty)$ converge pointwise as $N \rightarrow \infty$ to the unique solution of (3.2), where β is specified by*

$$(3.3) \quad \beta = \alpha \int_{\mathbb{R}^d} (1 - u(x))V(x)dx,$$

with $u : \mathbb{R}^d \rightarrow [0, 1]$ being the unique solution (provided by Lemma 1.6) subject to $u(x) \rightarrow 0$ as $x \rightarrow \infty$ of the modified Poisson equation

$$-2\Delta u(x) = \alpha V(x)(1 - u(x)).$$

As we prepare to justify the proposition, it is convenient to recast the definition of h_N in terms of the evolution of a particle, which we will call the tracer particle, picked uniformly at random at time zero. Since the distribution of particles at the initial time is invariant under particle reindexing, the next definition is suitable.

Definition 3.2 *The tracer particle is the particle whose time zero index is 1.*

The relationship between h_N and tracer particle survival probability is straightforward.

Lemma 3.3 *Let $N \in \mathbb{N}$ and $t \geq 0$. Then the microscopic candidate density $h_N(t)$ is equal to the \mathbb{P}_N -probability that the tracer particle has survived until time t .*

Proof. Note that

$$h_N(t) = N^{-1} \sum_{i=1}^N \mathbb{P}_N \left(\text{the } i\text{-th indexed particle survives until time } t \right).$$

However, the summand is independent of $i \in [1, N]$ due to the symmetry in both the initial placement of particles and in their dynamics. \square

In seeking to argue that h_N converges to the unique solution h of (3.2) in some appropriate sense, it is natural to try to find an expression for $\frac{d}{dt}h_N(t)$. We may hope to show that in fact this derivative equals $-\beta h_N(t)^2$ up to some error term which in some way tends to zero in the high N limit. Since the initial conditions $h_N(0)$ (for $N \in \mathbb{N}$), and $h(0)$, all coincide (with 1), we might then argue that $h_N \rightarrow h$ in some sense as $N \rightarrow \infty$.

With this aim in mind, we find an expression for the derivative of h_N in terms of the behaviour of the tracer particle in the microscopic model \mathbb{P}_N :

Lemma 3.4 *Let $N \in \mathbb{N}$ and $t \geq 0$. Let \mathbf{S}_t denote the event that the tracer particle under \mathbb{P}_N survives until time t . Then*

$$(3.4) \quad \frac{d}{dt}h_N(t) = -\lim_{\delta \searrow 0} \delta^{-1} \mathbb{P}_N(\mathbf{S}_t) \mathbb{P}_N(\mathbf{S}_{t+\delta}^c | \mathbf{S}_t),$$

should the limit on the right-hand side exist.

Proof. By Lemma 3.3, $h_N(t) = \mathbb{P}_N(\mathbf{S}_t)$, and thus $h_N(t + \delta) - h_N(t) = \mathbb{P}_N(\mathbf{S}_t \cap \mathbf{S}_{t+\delta}^c)$. \square

The expression (3.4) gives us a probabilistic means of thinking about the derivative of h_N . We should consider

- the \mathbb{P}_N -probability that the tracer particle survives until time t ; and
- given that it does so, the conditional probability that it is instantaneously annihilated in a collision.

There is a particular value of t for which the probability of this event is easier to evaluate: $t = 0$. In this case, the survival probability $\mathbb{P}_N(\mathbf{S}_0)$ is trivially equal to one. What then is the \mathbb{P}_N -probability that the tracer particle is annihilated before a very short time δ has passed? As our discussion will now tend to be heuristic rather than rigorous, we write subsequent statements as claims rather than lemmas.

Claim 3.5 *Let $t = t_N$ satisfy $t = o(\epsilon^2)$ as $N \rightarrow \infty$; equivalently, by (1.6), $t = o(N^{-2(2-d)})$. Then $\mathbb{P}_N(\mathbf{S}_t^c) = t\alpha \int_{\mathbb{R}^d} V(x) dx (1 + o(1))$ as $N \rightarrow \infty$.*

Sketch of proof. We begin by estimating the probability that the tracer particle collides with a given other particle in a very short interval $[0, t]$. In the case of the particle with index two, this probability is by definition given by

$$(3.5) \quad 1 - \exp \left\{ -\alpha \int_0^t V_\epsilon(X_2(s) - X_1(s)) \mathbf{1}_{\{1,2\} \subseteq I_{q(s)}} ds \right\},$$

since the exponential term here is the probability that the Poisson process for collision of particles indexed by 1 and 2 has yet to ring by time t . Note the presence of the indicator function for the

event $\{1, 2\} \subseteq I_{q(s)}$ that two particles have yet to be annihilated by time s . In fact, this term may be dropped from the expression at the expense of a lower order term as $t \searrow 0$ because, as we explained in Section 1.4, there is asymptotically zero probability that either of the two particles are annihilated during $[0, s]$ as $s \searrow 0$. Recalling that $V_\epsilon(\cdot) = \epsilon^{-2}V(\cdot/\epsilon)$ and that $V : \mathbb{R}^d \rightarrow [0, \infty)$ is supposed to be continuous, and noting that the difference $X_2 - X_1$ is a rate two Brownian motion, we see that

$$\int_0^t V_\epsilon(X_2(s) - X_1(s)) ds = \int_0^t V_\epsilon(X_2(0) - X_1(0))(1 + o(1)) ds$$

provided that $t = t_N$ is chosen so that $t\epsilon^{-2} \searrow 0$ as $N \rightarrow \infty$. That is to say, as the total particle number N tends to infinity, the collision probability (3.5) on $[0, t]$ is accurately approximated by $1 - e^{-t\alpha V_\epsilon(X_2(0) - X_1(0))}$ provided that t tends to zero more quickly than ϵ^2 , because, in this limiting regime, the locations X_1 and X_2 are asymptotically static on scale ϵ . Since $tV_\epsilon(x)$ converges to zero uniformly in $x \in \mathbb{R}^d$ in this regime, our asymptotic expression for (3.5) is $t\alpha V_\epsilon(X_2(0) - X_1(0))$. Since $X_2(0) - X_1(0)$ is simply uniformly distributed in \mathbb{T}^d , this quantity, after averaging over $X_1(0)$ and $X_2(0)$, equals $t\alpha \int_{\mathbb{R}^d} V_\epsilon(s) ds$, which is $t\alpha\epsilon^{d-2} \int_{\mathbb{R}^d} V(x) dx$.

By symmetry of the particle indices, this estimate applies equally to the probability of collision between particles with any two given indices in $[1, N]$. Since the probability that the tracer particle experiences two collisions during $[0, t]$ (with $t = o(\epsilon^2)$) behaves as $\binom{N}{2} (t\alpha\epsilon^{d-2} \int_{\mathbb{R}^d} V(x) dx)^2 = O(\epsilon^4)$, which is much smaller than the ϵ^2 -order probability of a single such collision, the probability that the tracer particle experiences some collision during $[0, t]$ is well approximated by the mean number of collisions that it experiences, which is

$$(N-1)t\alpha\epsilon^{d-2} \int_{\mathbb{R}^d} V(x) dx.$$

Recalling that $N = \epsilon^{2-d}$, we obtain the claim. \square

We are ready to return to (3.4) and record a limiting expression in high N for the time zero derivative of the microscopic candidate density:

Claim 3.6 *Let $N \in \mathbb{N}$. Then*

$$(3.6) \quad \lim_N \frac{dh_N}{dt}(0) = -\alpha \int_{\mathbb{R}^d} V(x) dx.$$

Proof. Since the expression $\lim_N \frac{d}{dt} h_N(0)$ equals $\lim_{N \rightarrow \infty} \lim_{\delta \searrow 0} \delta^{-1} \mathbb{P}_N(\mathbf{S}_\delta^c)$, the claim follows from Claim 3.5. \square

Recall that our aim to show that, in some appropriate sense, h_N converges to the solution h of (3.2) as $N \rightarrow \infty$. Claim 3.6 gives us a guess for the derivative at time zero of h : if it is the limit of the derivatives of its anticipated approximates, then $h'(0) = -\alpha \int_{\mathbb{R}^d} V(x) dx$. However, by its definition (3.2), $h'(0)$ is also $-\beta h(0)^2$, which is simply $-\beta$. In other words, the preceding argument has given us a guess for the recipe by which the macroscopic coagulation propensity β is

to be computed from the parameters in the underlying microscopic models. Namely, the argument points to the conclusion that

$$(3.7) \quad \beta = \alpha \int_{\mathbb{R}^d} V(x) dx.$$

3.1 Surviving and ghost particles

However, this guess is wrong. The formula (3.7) is not the correct relation between the microscopic and macroscopic coagulation propensities. To see why this is so, it is useful to introduce a coupling of our annihilating Brownian dynamics \mathbb{P}_N with a system of independent non-interacting Brownian particles \mathbb{P}'_N .

Definition 3.7 *Let \mathbf{C} denote a coupling of the law \mathbb{P}_N with a further law \mathbb{P}'_N . Under \mathbb{P}'_N , N particles are scattered in \mathbb{T}^d at time zero with the same law as in \mathbb{P}_N , and under \mathbf{C} the two initial conditions are always equal. In both marginals under \mathbf{C} , each particle pursues a given rate one Brownian motion, independently of the others. In the \mathbb{P}_N marginal, particles disappear on collision according to the rule for that dynamics; in the \mathbb{P}'_N marginal, the collision has no effect on either particle, and each continues its Brownian trajectory undisturbed.*

We say that under \mathbf{C} each particle is initially *surviving*. When a collision event occurs between two surviving particles, each becomes a ghost. In this way, the collection of all particles has the law \mathbb{P}'_N while the collection of surviving particles has the law \mathbb{P}_N .

For now, we need only one consequence of the coupling, namely that, for any $t \geq 0$, the time- t marginal of \mathbb{P}_N is stochastically dominated by its time zero marginal. To see this, note that this time- t marginal is dominated by the time- t marginal of all particles which has the law of the time 0 marginal of \mathbb{P}_N .

3.2 The naive guess is wrong

We now give an intuitive explanation of why (3.7) is the wrong relation between β , α and V . Suppose for convenience that $\int_{\mathbb{R}^d} V(x) dx = 1$. Now choose α to be fixed but very high. We are left with the formula $\beta = \alpha$, so that the solution $h : [0, \infty) \rightarrow [0, \infty)$ of (3.2) with $h(0) = 1$ equals $h(t) = \frac{1}{1+\alpha t}$.

Our high choice of α means that h drops towards zero quickly after time zero: specifically, $h(\alpha^{-1/2}) \leq \alpha^{-1/2}$. Here, however, we encounter a difficulty when we think of h as a limit of its approximates h_N . Expecting that $h_N(t) \rightarrow h(t)$ pointwise, recall from Lemma 3.3 that $h_N(t) = \mathbb{P}_N(\mathbf{S}_t)$ is the survival probability of the tracer particle until time t under \mathbb{P}_N . If the annihilation event \mathbf{S}_t^c is occur, then it is necessary that at some time before t , some other particle enters the interaction range of the tracer particle. By the coupling \mathbf{C} with an independent system of Brownian particles, to find an upper bound on $\mathbb{P}_N(\mathbf{S}_t^c)$, it is enough to bound the probability that, among N uniformly scattered particles in \mathbb{T}^d each performing Brownian motion, a given particle comes at some time in $[0, t]$ to distance ϵ of some other. Since the probability of such an approach at any given time is $N\epsilon^d = \epsilon^2$, and any such approach occurs for a mean time during of order ϵ^2 , this probability is bounded above by a constant multiple of t for small t . (An explicit bound on this mean time

is that the time spent overlapping during $[0, \infty)$ by two rate-one radius- ϵ Brownian spheres in \mathbb{R}^d , with $d \geq 3$, which are tangent at time zero is in expectation at most $(1 + \frac{2^{1-d}}{(d-2)\Gamma(d/2+1)})\epsilon^2$.) Thus, $\mathbb{P}_N(\mathbf{S}_t) \geq 1 - Ct$, where $C > 0$ may be chosen uniformly in both $N \in \mathbb{N}$ and $\alpha \in (0, \infty)$. For all $\alpha > 0$, the pointwise convergence of h_N to h forces $h(t) \geq 1 - Ct$. However, this is inconsistent with $h(\alpha^{-1/2}) \leq \alpha^{-1/2}$ for $\alpha > (C + 1)^2$. We conclude then that the guessed formula (3.7) is in fact wrong.

What is wrong with the derivation of (3.7) is that, in fact, the initial Poissonian distribution of particles is in a certain sense unstable, making an inference based on an analysis at time zero misleading. Although at later times the particle distribution is Poissonian in the large, there is a microscopic repulsion effect which modifies this: in the order ϵ vicinity of the tracer particle at some positive time, there is a diminished probability for presence of another particle. Indeed, this other particle may have already collided with the tracer particle, in which case, the tracer particle and the other particle would not in fact be located close to each other because each would have vanished. We now turn to quantifying the effect of this mechanism of curtailment of interaction due to collision and so derive the correction needed to (3.7). To make sense of the notion of a particle in the vicinity of the tracer particle which may already have vanished at a certain time, we will make use of our coupling of surviving and ghost particles.

3.3 Quantifying the correction

The discussion in the preceding section reveals that our computation of $\frac{dh_N(t)}{dt}$ for $t = 0$ may have exceptional features which change as t increases. In fact, as we will see, this change will be apparent already when t reaches the order of ϵ^2 . To understand the change quantitatively, we want to return to Lemma 3.4 and use the right-hand side of (3.4) to compute $\frac{dh_N(t)}{dt}$ for $t > 0$. This involves computing the terms $\mathbb{P}_N(\mathbf{S}_t)$ and $\mathbb{P}_N(\mathbf{S}_{t+\delta}^c | \mathbf{S}_t)$ appearing on the right-hand side of (3.4), when $t > 0$ is fixed, and $\delta > 0$ is infinitesimally small. Reexpressing the events in terms of the coupling \mathbf{C} , \mathbf{S}_t is the event that the tracer particle survives until time t , and $\mathbf{S}_t \cap \mathbf{S}_{t+\delta}^c$ is the event that the tracer particle changes its status from surviving to ghost during the short time interval $[t, t + \delta]$.

Claim 3.8 *Suppose that $\delta\epsilon^{-2} \searrow 0$ as $N \rightarrow \infty$. Then*

$$\mathbb{P}_N(\mathbf{S}_{t+\delta}^c \cap \mathbf{S}_t) = \delta\alpha h(t)^2 \int_{\mathbb{R}^d} (1 - u(x))V(x)dx (1 + o(1)),$$

as $N \rightarrow \infty$, where $u : \mathbb{R}^d \rightarrow [0, 1]$ is specified in Proposition 3.1.

Seeking to justify this claim, we let $C_{i,j}(t, \delta)$ denote the event that

- $|X_i(t) - X_j(t)| \leq \epsilon$,
- and these two particles collide during $[t, t + \delta]$.

As we now record, the event $\mathbf{S}_t \cap \mathbf{S}_{t+\delta}^c$ is characterized up to a probability of smaller order by the survival of the tracer particle until time t and the presence at time t of some other surviving particle

in $X_1(t) + \mathbf{B}_\epsilon$, a set that contains the interaction range of the tracer particle, which during $[t, t + \delta]$ collides with the tracer particle. (Here, \mathbf{B}_ϵ denotes the Euclidean ball of radius ϵ about the origin, so that, since the support of V is contained in the Euclidean unit ball, the set $X_1(t) + \mathbf{B}_\epsilon$ indeed contains the interaction range of the tracer particle.)

Claim 3.9 *For each $t > 0$, there exists $C_t > 0$ such that*

$$\mathfrak{C}\left(\left(\mathbf{S}_t \cap \mathbf{S}_{t+\delta}^c\right) \Delta \left(\bigcup_{j=2}^N \{\{1, j\} \in I_{q(t)}\} \cap C_{1,j}(t, \delta)\right)\right) \leq C_t \mathfrak{C}\left(\mathbf{S}_t \cap \mathbf{S}_{t+\delta}^c\right) (\delta + \alpha \epsilon \delta^{1/2}).$$

Sketch of proof. The claim will emerge from two assertions. First,

$$(3.8) \quad \mathfrak{C}\left(\left(\mathbf{S}_t \cap \mathbf{S}_{t+\delta}^c\right) \Delta \left(\bigcup_{j=2}^N \{\{1, j\} \in I_{q(t)}\} \cap C_{1,j}(t, \delta)\right)\right) \leq C\delta^2 + C\alpha\epsilon\delta^{3/2},$$

and, second, for some t -dependent constant $c > 0$,

$$(3.9) \quad \mathfrak{C}\left(\mathbf{S}_t \cap \mathbf{S}_{t+\delta}^c\right) \geq c\delta.$$

The second bound holds because, as we described in Section 1.4, the tracer particle will survive to any time t with some positive, t -dependent probability, and it is then liable to collide with some other particle at a rate of order one.

Regarding the symmetric difference in (3.8), note that, if one event occurs without the other, the cause must be one of the following:

- although $\{1, j\} \in I_{q(t)}$ and $C_{1,j}$ occur for some $j \in [2, N]$, there is a third particle which collides with particle j after time t but before the collision of j with 1 that happens before time $t + \delta$;
- or, $\mathbf{S}_t \cap \mathbf{S}_{t+\delta}^c$ occurs due to some particle $j \in [2, N]$, which is not in the interaction range $X_1 + \mathbf{B}_\epsilon(t)$ at time t , entering this range and colliding with 1 during $[t, t + \delta]$.

Regarding the first possibility, for given $j \in [2, N]$, the probability of $C_{1,j}$ is at most $C\epsilon^d\delta\epsilon^{-2}$, while, given this event, the conditional probability of some third particle behaving as described is at most $C\delta$. Summing over j , the probability is at most $CN\delta^2\epsilon^{d-2} = C\delta^2$.

For given $j \in [2, N]$, the probability of the second occurrence is at most a constant multiple of $\epsilon\delta^{3/2}$. We only sketch how this bound is obtained. Should particle j at time t lie within a distance of order $\delta^{1/2}$ of the boundary of the interaction range $X_1 + \mathbf{B}_\epsilon$ of the tracer particle, there is positive probability that particle j enters this range during the ensuing δ units of time, and should this happen, there is conditional probability at most $1 - \exp\{-\alpha|V|_\infty\delta\} \leq \alpha|V|_\infty\delta \leq C\alpha\delta$ of collision between the two particles. The probability of this turn of events is thus $|\mathbf{B}_{\epsilon+\delta^{1/2}} \setminus \mathbf{B}_\epsilon|C\alpha\delta \leq C\epsilon^{d-1}\delta^{3/2}$. On the other hand, it is easily checked that there is negligible probability of such a collision should particle j at time t lie at much greater distance than $\delta^{1/2}$ from the boundary of

$X_1 + B_\epsilon$. Summing over the $N - 1 \sim \epsilon^{2-d}$ choices of $j \in [2, N]$, we see that the probability of the second listed event is at most $C\epsilon\delta^{3/2}$. \square

The next two claims estimate the probability of $\{\{1, j\} \in I_{q(t)}\} \cap C_{1,j}(t, \delta)$ for $j \in [2, N]$ and will lead to Claim 3.8.

Claim 3.10 *Suppose that $\delta\epsilon^{-2} \searrow 0$ as $N \rightarrow \infty$. Then, for each $x \in B_\epsilon$,*

$$\mathbb{C}\left(C_{1,2}(t, \delta) \mid X_1(t) - X_2(t) = x\right) = \epsilon^{-2}\delta\alpha V(x/\epsilon) (1 + o(1)).$$

Claim 3.11 *For each $x \in B_\epsilon$, we have that*

$$\mathbb{C}\left(\{1, 2\} \subseteq I_{q(t)} \mid X_1(t) - X_2(t) = x\right) = (1 - u(x/\epsilon))h(t)^2.$$

Proof of Claim 3.8. Note that, conditionally on $X_2 - X_1 \in B_\epsilon$, $X_2 - X_1$ is uniform on B_ϵ . In light of this, and Claims 3.10 and 3.11, we see that, in the limit in question,

$$\mathbb{C}(C_{1,2}(t, \delta), \{1, 2\} \subseteq I_{q(t)}) = \epsilon^{-2}\delta\alpha h(t)^2 \int_{B_\epsilon} V(x/\epsilon)(1 - u(x/\epsilon))dx (1 + o(1)),$$

whose integral term may be also written as $\epsilon^d \int_{\mathbb{R}^d} V(x)(1 - u(x))dx$. By particle symmetry and $N - 1 \sim \epsilon^{2-d}$, we obtain

$$\sum_{j=2}^N \mathbb{C}\left(C_{1,2}(t, \delta), \{1, 2\} \subseteq I_{q(t)}\right) = \delta\alpha h(t)^2 \int_{\mathbb{R}^d} V(x)(1 - u(x))dx (1 + o(1)).$$

It is easy to convince oneself that typically the occurrence of $\cup_{j=2}^N C_{1,2}(t, \delta) \cap \{\{1, 2\} \subseteq I_{q(t)}\}$ entails the occurrence of exactly one of the constituent events. For this reason, the preceding equality holds equally for the probability of $\cup_{j=2}^N C_{1,2}(t, \delta) \cap \{\{1, 2\} \subseteq I_{q(t)}\}$. Thus, Claim 3.8 follows from Claim 3.9. \square

Proof of Claim 3.10. Conditionally on $X_1(t) - X_2(t)$ being a given $x \in B_\epsilon$, the probability of collision between 1 and 2 during $[t, t + \delta]$ is $\delta V_\epsilon(x)(1 + o(1))$, since $V : \mathbb{R}^d \rightarrow [0, \infty)$ is assumed continuous, and $X_1 - X_2$ is asymptotically static on scale ϵ . \square

Proof of Claim 3.11. To reiterate the problem, given that at time t particles 1 and 2 have displacement $x \in B_\epsilon$, what is the probability that both survive to this time? There are two reasons why one or other may be a ghost particle at time t :

- it may be that, during $[0, t]$, at a moment when each of X_1 and X_2 are surviving, a collision between this pair occurs;
- it may be that one or other of X_1 and X_2 , at a moment when this particle is surviving, collides with some other surviving particle.

Calling these two events E_1^t and E_2^t , we want to gauge the probability of $(E_1^t)^c \cap (E_2^t)^c$.

In considering these possibilities, it is convenient to reverse time, running time backwards from t to 0. We will now use these time coordinates, where the forward time evolution from 0 to t corresponds to the usual evolution backwards from t to 0. Note that, because we do not condition on X_1 or X_2 surviving until time t , the conditional distribution of the trajectories $X_1, X_2 : [0, t] \rightarrow \mathbb{T}^d$ in the new time coordinates is a pair of independent rate one Brownian motions, where $X_1(0)$ is uniformly distributed on \mathbb{T}^d and $X_2(0) = X_1(0) + x$. Phrased in these terms, E_1^t is the event of collision between X_1 and X_2 during $[0, t]$.

In light of Lemma 1.5, we learn that

Claim 3.12 *If $t = t_N$ satisfies $t/\epsilon^2 \rightarrow \infty$ as $N \rightarrow \infty$, then*

$$\mathbb{C}\left((E_1^t)^c\right) = (1 - u(x/\epsilon))(1 + o(1)),$$

where $u : \mathbb{R}^d \rightarrow [0, 1]$ is specified in Proposition 3.1.

We also need to estimate the conditional probability that E_2^t occurs given that E_1^t does not.

Claim 3.13 *If $t = t_N$ satisfies $t/\epsilon^2 \rightarrow \infty$ as $N \rightarrow \infty$, then*

$$\mathbb{C}\left((E_2^t)^c \mid (E_1^t)^c\right) = h(t)^2(1 + o(1)).$$

Sketch of proof. Should E_1^t not occur, the two particle trajectories $X_1, X_2 : [0, t] \rightarrow \mathbb{T}^d$, begun at the points 0 and x at distance of order ϵ , will not experience collision, and will separate to a distance much greater than ϵ in a time of order large compared to ϵ^2 ; by our assumption on the time t , this separation occurs on a time scale (called s) much shorter than t , and, after it has done so, it is reasonable to think that the collision behaviour of the two trajectories will become effectively independent of one another. Any given one of these trajectories experiences collision with some other particle with probability $h(t - s)$, by the definition of h ; assuming this independence, both particles survive collision during $[0, t]$ with probability $h(t - s)^2$. However, since $s \ll t$, $h(s) = h(t)(1 + o(1))$ as $N \rightarrow \infty$. \square

Claim 3.11 follows from Claims 3.12 and 3.13. \square

We are now able to complete our sketch proof of the formula for β for the model in question.

Sketch proof of Proposition 3.1. Applying Lemma 3.4 and Claim 3.8, we find that

$$\frac{dh_N(t)}{dt} = -\alpha h(t)^2 \int_{\mathbb{R}^d} (1 - u(x))V(x)dx \left(1 + o(1)\right),$$

provided that $t/\epsilon^2 \rightarrow \infty$ as $N \rightarrow \infty$. This estimate does not control the behaviour of $\frac{dh_N(t)}{dt}$ on the small time scale where t is of order ϵ^2 . Here, however, arguments in the style of those leading to Claim 3.6 justify that this derivative is non-positive and bounded below by $-\alpha \int_{\mathbb{R}^d} V(x)dx$. Recalling that h is the unique solution of (3.2), we see that, since $h_N(0) = h(0) = 1$ for all $N \in \mathbb{N}$, h_N converges h pointwise as $N \rightarrow \infty$. \square

3.4 Bose-Einstein condensates, and a parallel in the form of macroscopic interaction

We briefly discuss the quantum mechanical problem of the dynamics of a collection of N bosons in three dimensions that interact via a short-range pair potential, based on Erdős, Schlein and Yau's derivation [7] of the macroscopic evolution of the system. The dynamics of the system is governed by the Schrödinger equation

$$i\partial_t\psi_{N,t} = H_N\psi_{N,t} ,$$

for the wave function $\psi_{N,t} \in L^2_s(\mathbb{R}^{3N})$, the subspace of $L^2(\mathbb{R}^{3N})$ consisting of all functions symmetric under permutations of the N particles. Short-range repulsive interaction is modelled by the choice of Hamiltonian $H_N = H_{\beta,N}$,

$$H_{\beta,N} = -\sum_{j=1}^N \Delta_j + \frac{1}{N} \sum_{1 \leq i < j \leq N} N^{3\beta} V(N^\beta(x_i - x_j)) ,$$

where $V : \mathbb{R}^3 \rightarrow [0, \infty)$ is a compactly supported interaction potential, and $\beta > 0$ is a parameter. (Note that this notation is in conflict with our use of $\beta(\cdot, \cdot)$.) The choice of $\beta = 1$ provides the closest parallel with the main discussion in this survey and is the principal object of study in [7]. The macroscopic evolution of the system may be summarized by a decoupling property enjoyed by the k -particle reduced density matrices; these matrices in [7] are shown by an analysis of the BBGKY hierarchy to factorize asymptotically in the high N limit, with the factor governed by the non-linear Gross-Pitaevskii equation

$$i\partial_t\psi_t = \Delta\psi_t + \sigma|\psi_t|^2\psi_t ,$$

where the coupling constant σ is given by

$$\sigma = \begin{cases} b_0 & \text{if } 0 < \beta < 1 , \\ 8\pi a_0 & \text{if } \beta = 1 . \end{cases}$$

Here, $b_0 = \int_{\mathbb{R}^3} V(x)dx$, while a_0 satisfies

$$a_0 = \frac{1}{8\pi} \int_{\mathbb{R}^3} V(x)(1 - \omega_0(x))dx ,$$

with ω_0 being the unique solution to

$$\left[-\Delta + \frac{1}{2}V(x) \right] (1 - \omega_0(x)) = 0$$

that satisfies $\lim_{x \rightarrow \infty} \omega_0(x) = 0$. (Note that ω_0 is nothing other than the solution of (1.7) if we take $\frac{\alpha(n,m)}{d(n)+d(m)}$ equal to $1/2$.)

That is, the macroscopic interaction coefficient undergoes a transition as $\beta \in (0, 1)$ changes to $\beta = 1$ in precise correspondence to the modification from naive guess $\beta = \int_{\mathbb{R}^d} V(x)dx$ to β given by (1.8) which we have devoted this section to discussing.

Indeed, we may specify a collection of random models $\mathbb{P}_{N,\beta}$, with $\beta > 0$, in such a way that our models \mathbb{P}_N coincide with $\mathbb{P}_{N,1}$, while $\mathbb{P}_{N,\beta}$, $\beta \in (0, 1)$, form counterparts to the interacting bosonic systems at such values of β . Maintaining the relation (1.6) between ϵ and N , we modify the pairwise collision rule from Section 1.3 from one under which the particles indexed by $i, j \in [1, N]$ collide at rate $\alpha(m_i, m_j)\epsilon^{-2}V\left(\frac{x_i-x_j}{\epsilon}\right)$ to one whose rate is $\alpha(m_i, m_j)\epsilon^{-2+d(1-\beta)}V\left(\frac{x_i-x_j}{\epsilon^\beta}\right)$. The rule for $\mathbb{P}_{N,\beta}$ is determined in order that a typical particle maintain a order unit interaction with all the others per unit time, so that the new models remain in the regime of constant mean free path. The derivation of Theorem 1.1 may be reprised for choices of $\beta \in (0, 1)$, with the formula for the macroscopic interaction rates $\beta(n, m)$ now given by $\beta(n, m) = \alpha(n, m) \int_{\mathbb{R}^d} V(x)dx$. The new formula holds in essence because microscopic pairwise repulsion is absent asymptotically in high N in these models.

See [27] for a blog post by Terry Tao, written after a talk by Natasa Pavlovic, which provides a more informative summary of this quantum problem, including at its end, and in the ensuing comments, mention of the dichotomy between interaction coefficient in the cases $\beta \in (0, 1)$ and $\beta = 1$.

4 The route to Theorem 1.1

In this section, we explain the overall plan for proving the main theorem, and reduce it to a fundamental proposition, the *Stosszahlansatz*, which concerns the total particle coagulation propensity in the microscopic models.

4.1 Approximating the PDE using microscopic candidate densities

4.1.1 A microscopic counterpart to the PDE in weak form

Recall the weak formulation (1.9) of the Smoluchowski PDE. Our aim is to show that the particle densities in the microscopic model \mathbb{P}_N converge in a suitable sense to this weak solution. To do so, we find a microscopic counterpart to the equation (1.9), namely an equation expressed in terms of the law \mathbb{P}_N . Note that (1.9) expresses the change in the quantity $\int_{\mathbb{R}^d} J(x, t)f_n(t)dx$ that occurs between times 0 and T as an integral over the intervening duration $[0, T]$ of the differential changes caused by variation in the test function J , and by the diffusion and coagulation of the particles being modelled. The quantity $\int_{\mathbb{R}^d} J(x, t)f_n(t)dx$ is an expression for the total number of particles of mass n at time t , where each particle is weighted by J . As such, it has a clear microscopic analogue: under the law \mathbb{P}_N , the random variable $\sum_{i \in I_q(t)} J_n(x_i, m_i)\mathbf{1}_{m_i(t)=n}$, which is the sum over mass- n particles at time t where each particle carries a weight given by J . The form of the Markov generator for the dynamics of \mathbb{P}_N now provides us with an analogue of the weak formulation (1.2)

of the Smoluchowski PDE:

$$\begin{aligned}
(4.1) \quad & \epsilon^{d-2} \sum_{i \in I_q(T)} J_n(x_i, T) \mathbf{1}_{m_i(T)=n} - \epsilon^{d-2} \sum_{i \in I_q(0)} J_n(x_i, 0) \mathbf{1}_{m_i(0)=n} \\
& = \epsilon^{d-2} \int_0^T \left(\sum_{i \in I_q(t)} \frac{\partial J_n}{\partial t}(x_i, t) \mathbf{1}_{m_i(t)=n} + \sum_{i \in I_q(t)} d(n) \Delta J_n(x_i, t) \mathbf{1}_{m_i(t)=n} \right. \\
& \quad \left. + \sum_{i, j \in I_q(t)} \alpha(m_i, m_j) V_\epsilon(x_i - x_j) \mathfrak{C}_{i, j, t, n} J_n \right) dt + M_T.
\end{aligned}$$

In the integrands on the right-hand side, we see the infinitesimal mean changes caused by the time-variation of J , by the diffusion of the individual particles, and by their collision in pairs; the final term is a martingale (which we will argue to be suitably small). In the collision term, the real-valued quantity $\mathfrak{C}_{i, j, t, n} J_n$ is the instantaneous change in the value of

$$\sum_{i \in I_q(t)} J_n(x_i, t) \mathbf{1}_{m_i(t)=n}$$

that is caused by the collision of particles (x_i, m_i) and (x_j, m_j) at time t . As such, it has the expression

$$\begin{aligned}
\mathfrak{C}_{x_i, x_j, t, n} J_n & = \frac{m_i}{m_i + m_j} J_n(x_i, t) \mathbf{1}_{m_i(t) + m_j(t) = n} + \frac{m_j}{m_i + m_j} J_n(x_j, t) \mathbf{1}_{m_i(t) + m_j(t) = n} \\
& \quad - J_n(x_i, t) \mathbf{1}_{m_i(t) = n} - J_n(x_j, t) \mathbf{1}_{m_j(t) = n} :
\end{aligned}$$

two gain terms arise from the appearance of a new particle at one or other of the locations of the disappearing pair, and two loss terms correspond to the disappearance of each element of this pair.

4.1.2 Introducing microscopic candidate densities

Our plan for the kinetic limit derivation of the Smoluchowski PDE is to argue that (1.2) emerges in a suitable sense when we take the high N limit, with the martingale term M_T vanishing in this limit. To implement this plan, we introduce *microscopic candidate densities* $f_{n, t}^{\epsilon, \delta} : \mathbb{R}^d \rightarrow [0, \infty)$ of mass n particles at time t under the law \mathbb{P}_N . Here, $\delta > 0$ is a fixed positive quantity, while the interaction radius ϵ is determined from N as usual by $N\epsilon^{d-2} = Z$. The candidate density is given by

$$(4.2) \quad f_{n, t}^{\epsilon, \delta}(u) = \epsilon^{d-2} \sum_{i \in I_q(t)} \delta^{-d} \eta\left(\frac{x_i - u}{\delta}\right), u \in \mathbb{R}^d,$$

where $\eta : \mathbb{R}^d \rightarrow [0, \infty)$ is a smooth compactly supported function for which $\int_{\mathbb{R}^d} \eta(x) dx = 1$. That is, $f_{n, t}^{\epsilon, \delta}(u)$ is a statistic reporting the number of particles in a small macroscopic region about the point $u \in \mathbb{R}^d$ at time t in the model \mathbb{P}_N ; note that the time zero microscopic candidate density $f_{n, 0}^{\epsilon, \delta}(u)$ has a high N pointwise limit which as a function of u is given by the convolution of the

initial condition h_n of (1.2) and η^δ . Taking a $\delta \searrow 0$ limit after this limit, we see that, at time zero at least, the appropriate initial condition $h_n(u)$ is obtained at all $u \in \mathbb{R}^d$.

Our aim is to argue that something similar happens at all later times $t > 0$. To do so, we will replace the various terms appearing in the expectation value of equation (4.1) with approximating terms expressed in terms of the microscopic candidate densities, and then take the high N and then the low δ limit. If we are to reach (1.9) as a result, we will need to understand that the new terms approximate the old ones appropriately.

4.1.3 Replacing old terms by new: simple cases

The first term $\sum_{i \in I_q(T)} J_n(x_i, T) \mathbf{1}_{m_i(T)=n}$ has a simple counterpart expressed in the fashion we seek: $\int_{\mathbb{R}^d} J_n(x, T) f_n^{\epsilon, \delta}(x) dx$. Nor is any nontrivial estimate needed to find a suitable bound on the difference of the terms in this case. Indeed, the two expressions differ by

$$\epsilon^{d-2} \sum_{i \in I_q(T)} \left(J_n(x_i, T) - \int_{\mathbb{R}^d} J_n(y, T) \delta^{-d} \eta\left(\frac{x_i - y}{\delta}\right) dy \right),$$

which in absolute value is at most $Z\delta \|\nabla J_n\|_\infty \leq C\delta$, since total particle number at time T is less than the initial total $N = Z\epsilon^{2-d}$.

The first and second terms on the right-hand side of (4.1) similarly have counterparts

$$\int_{\mathbb{R}^d \times [0, T]} \frac{\partial J_n}{\partial t}(x, t) \cdot f_n^{\epsilon, \delta}(x) dx dt \quad \text{and} \quad \int_{\mathbb{R}^d \times [0, T]} d(n) \Delta J_n(x, t) \cdot f_n^{\epsilon, \delta}(x) dx dt.$$

Each pair of term and counterpart likewise has a difference which in absolute value is deterministically bounded above by some constant multiple of δ .

4.1.4 Replacing the coagulation term by using the Stosszahlansatz

Given the form of the coagulation term present in (1.9), there is a clear candidate for the form of the term which will form a counterpart to the interaction term appearing in the third line of (4.1): namely, $\int_{\mathbb{R}^d \times [0, T]} J_n(x, t) (Q_1^n(f_n^{\epsilon, \delta}(x)) - Q_2^n(f_n^{\epsilon, \delta}(x))) dx dt$. In stark contrast to the other cases, proving that the replacement of the collision term with this counterpart involves a suitably small error is a major undertaking. We now state the key estimate in this regard, a proposition which we will sometimes call the Stosszahlansatz. Recall that the coefficients $\beta : \mathbb{N}^2 \rightarrow (0, \infty)$ are specified in (1.8).

Proposition 4.1 *For each $n, m \in \mathbb{N}$, we have that*

$$\begin{aligned} & \epsilon^{d-2} \mathbb{E}_N \int_0^T \sum_{i, j \in I_q(t)} \alpha(m_i, m_j) V_\epsilon(x_i - x_j) J_n(x_i, t) \mathbf{1}_{m_i(t)=n, m_j(t)=m} \\ &= \beta(n, m) \int_0^T \int_{\mathbb{R}^d} J_n(x, t) f_n^{\epsilon, \delta}(x, t) f_m^{\epsilon, \delta}(x, t) dx dt + \text{Err}_{n, m}(\epsilon, \delta), \end{aligned}$$

where the error $\text{Err}_{n,m}$ satisfies

$$\lim_{\delta \searrow 0} \limsup_{\epsilon \searrow 0} \sum_{m \in \mathbb{N}} \mathbb{E}_N |\text{Err}_{n,m}(\epsilon, \delta)| = 0.$$

Setting $J_n = 1$ for ease of description, note that the integral on the left-hand side is the cumulative rate at which particle pairs of masses n and m are liable to coagulate during all of $[0, T]$; by the relation (1.6) and the anticipated survival of a positive fraction of particles at any given positive time, we see that the normalization ϵ^{d-2} on the left-hand side is chosen so that the overall expression is of unit order in the high N limit. Proposition 4.1 asserts that this expression is closely approximated by the integral over space-time of the product of the microscopic candidate densities multiplied by the constant coefficient $\beta(n, m)$. As such, this $\beta(n, m)$ is a macroscopic coagulation propensity of pairs of particles of these masses.

Proposition 4.1 is an expression of the type of precollisional particle independence that we discussed for elastic billiards in Subsection 2.3.2; here, this independence is manifested by the presence of the product $f_n^{\epsilon, \delta}(x, t) f_m^{\epsilon, \delta}(x, t)$.

We now explain how Proposition 4.1 may be invoked to show that the coagulation term in (4.1) is suitably approximated by its counterpart. Recall that the instantaneous change $\mathfrak{C}_{x_i, x_j, t, n} J_n$ is comprised of four terms: two gain terms and two loss terms. Consider the third of these terms, which is the first loss term. This term expresses the instantaneous loss of terms $J_n(x_i, t)$ due to the collision at time t of x_i with some other particle x_j . This other particle may have any mass $m_j \in \mathbb{N}$. Writing this term as a sum over that mass, we obtain that the term equals

$$-\epsilon^{d-2} \mathbb{E}_N \int_0^T \sum_{m=1}^{\infty} \sum_{i, j \in I_q(t)} \alpha(m_i, m_j) V_\epsilon(x_i - x_j) J_n(x_i, t) \mathbf{1}_{m_i(t)=n, m_j(t)=m} dt.$$

Applying Proposition 4.1, we find that the term equals

$$-\sum_{m=1}^{\infty} \beta(n, m) \int_{\mathbb{R}^d} J_n(x, t) f_n^{\epsilon, \delta}(x, t) f_m^{\epsilon, \delta}(x, t) dx + \text{Err}_n(\epsilon, \delta).$$

where this error term, after the sum over $m \in \mathbb{N}$, is known to satisfy

$$(4.3) \quad \lim_{\delta \searrow 0} \limsup_{\epsilon \searrow 0} \mathbb{E}_N |\text{Err}_n(\epsilon, \delta)| = 0.$$

Exactly the same considerations apply to the second of the loss terms because the two terms are equal due to the symmetry of the interaction kernel V .

The comparable estimate for each of the gain terms is slightly easier to handle, because a particle of mass n may be produced by only finitely many mass pairs – $(1, n-1), (2, n-2), \dots, (n-1, 1)$ – rather than the infinite number of choices – $(n, 1), (n, 2), \dots$ – which may cause such a particle to disappear. Regarding the first term, an application of Proposition 4.1 yields that

$$\epsilon^{d-2} \mathbb{E}_N \int_0^T \sum_{m=1}^{n-1} \sum_{i, j \in I_q(t)} \alpha(m_i, m_j) V_\epsilon(x_i - x_j) \frac{m}{n} J_n(x_i, t) \mathbf{1}_{m_i(t)=m, m_j(t)=n-m} dt$$

equals

$$\sum_{m=1}^n \beta(m, n-m) \int_{\mathbb{R}^d \times [0, T)} J_n(x, t) \frac{m}{n} f_n^{\epsilon, \delta}(x, t) f_m^{\epsilon, \delta}(x, t) dx dt + \text{Err}_n(\epsilon, \delta),$$

where likewise the error satisfies (4.3). The second gain term differs only in that $\frac{n-m}{n}$ replaces $\frac{m}{n}$; thus, the total gain term satisfies the same statement with this term omitted.

4.1.5 The martingale term is replaced by zero

The martingale term M_T in (4.1) is treated by arguing that it is typically small in absolute value:

Proposition 4.2 *There exists $C > 0$ such that, for each $N \in \mathbb{N}$, $\sup_{T \in (0, \infty)} \mathbb{E}_N M(T)^2 \leq C \epsilon^{d-2}$.*

The martingale term is in a sense much smaller than the collision term treated by the Stosszahlansatz: it vanishes before the low δ limit is even taken. The ideas in the proof of Proposition 4.2 are already in large part seen in the proof of the more substantial Proposition 4.1 and we will not explain their specifics; it is in Section 5 that the martingale term is treated in [10].

4.1.6 The counterpart to the PDE using microscopic candidate densities

By using Propositions 4.1 and 4.2 and the other more trivial estimates, we are able to replace each term in (4.1) with its counterpart, expressed in terms of the microscopic candidate densities, and obtain the following bound on the error in the resulting near identity:

$$\begin{aligned} (4.4) \quad & \int_{\mathbb{R}^d} J_n(x, T) f_n^{\epsilon, \delta}(x) dx - \int_{\mathbb{R}^d} J_n(x, 0) f_n^{\epsilon, \delta}(x) dx \\ &= \int_{\mathbb{R}^d \times [0, T)} \frac{\partial J_n}{\partial t}(x, t) f_n^{\epsilon, \delta}(x) dx dt + \int_{\mathbb{R}^d \times [0, T)} d(n) \Delta J_n(x, t) f_n^{\epsilon, \delta}(x) dx dt \\ & \quad + \sum_{m=1}^n \beta(m, n-m) \int_{\mathbb{R}^d \times [0, T)} J_n(x, t) f_n^{\epsilon, \delta}(x, t) f_m^{\epsilon, \delta}(x, t) dx dt \\ & \quad - 2 \sum_{m=1}^{\infty} \beta(n, m) \int_{\mathbb{R}^d \times [0, T)} J_n(x, t) f_n^{\epsilon, \delta}(x, t) f_m^{\epsilon, \delta}(x, t) dx dt + \text{Err}_n(\epsilon, \delta). \end{aligned}$$

where the error satisfies (4.3) because each of the errors used in the five estimates which we applied does.

4.2 Taking the limit to obtain the Smoluchowski PDE

Our approximate identity (4.4) closely resembles the equation (1.9) satisfied by a weak solution of the Smoluchowski PDE: we simply replace the solution of the latter with the microscopic candidate densities to obtain the former. However, to pass to (1.9) from (4.4) in the limit of low ϵ followed by low δ , we require a little more information about the microscopic models \mathbb{P}_N .

4.2.1 The approximate identity rewritten using empirical measures

Recall from Theorem 1.1 that in fact we express approximation by \mathbb{P}_N for high N of the Smoluchowski PDE by using the empirical measures μ_N^f valued in space-mass-time $\mathbb{R}^d \times \mathbb{N} \times [0, \infty)$. In Section 1.7, we let $\mu_{N,n}^f$ denote the mass n marginal of μ^f , the empirical measure in space-time for particles of mass n , for each $n \in \mathbb{N}$. That is,

$$\mu_{N,n}^f = \epsilon^{d-2} \sum_{i \in I_q(t)} \delta_{(x_i, t)} \mathbf{1}_{m_i=n} dt.$$

On \mathbb{P}_N , the microscopic candidate densities are expressed in terms of the empirical measures by $f_n^{\epsilon, \delta}(x, t) = (\mu_{N,n}^f * \eta^\delta)(x, t)$ for all $(x, t) \in \mathbb{R}^d \times [0, \infty)$, where the convolution is in the space variable.

Suppose that ν is a vague limit point of the sequence $\{\mu_N^f : N \in \mathbb{N}\}$. After extraction of a subsequence if necessary, there exists, by definition, a sequence $\{\nu_N : N \in \mathbb{N}\}$ of measures $\nu_N \in \mathcal{M}$ such that ν_N lies in the support of μ_N^f for each $N \in \mathbb{N}$, and with ν_N converging vaguely to ν . Vague convergence readily implies that $\nu_N * \eta^\delta$ converges pointwise to $\nu * \eta^\delta$. This convergence allows us to take the low ϵ limit of (4.4) to arrive at

$$\begin{aligned} (4.5) \quad & \int_{\mathbb{R}^d} J_n(x, T) \cdot (\nu_n * \eta^\delta)(x) dx - \int_{\mathbb{R}^d} J_n(x, 0) \cdot (\nu_n * \eta^\delta)(x) dx \\ &= \int_{\mathbb{R}^d \times [0, T)} \frac{\partial J_n}{\partial t}(x, t) \cdot (\nu_n * \eta^\delta)(x, t) dx dt + \int_{\mathbb{R}^d \times [0, T)} d(n) \Delta J_n(x, t) \cdot (\nu_n * \eta^\delta)(x, t) dx dt \\ & \quad + \sum_{m=1}^n \beta(m, n-m) \int_{\mathbb{R}^d \times [0, T)} J_n(x, t) \cdot (\nu_n * \eta^\delta)(x, t) \cdot (\nu_m * \eta^\delta)(x, t) dx dt \\ & \quad - 2 \sum_{m=1}^{\infty} \beta(n, m) \int_{\mathbb{R}^d \times [0, T)} J_n(x, t) \cdot (\nu_n * \eta^\delta)(x, t) \cdot (\nu_m * \eta^\delta)(x, t) dx dt + \text{Err}_n(\delta), \end{aligned}$$

where $\lim_{\delta \searrow 0} \text{Err}_n(\delta) = 0$ almost surely.

4.2.2 Preparing for the final step towards the PDE: uniform integrability

In order to conclude the proof of Theorem 1.1, two steps are needed. First,

Proposition 4.3 *Let ν be a vague limit point of $\{\mu_N^f : N \in \mathbb{N}\}$. For each $n \in \mathbb{N}$, the marginal ν_n has a density $f_n : \mathbb{R}^d \times [0, \infty) \rightarrow [0, \infty)$ with respect to Lebesgue measure on space-time $\mathbb{R}^d \times [0, \infty)$.*

Second, we must argue that the collection $\{f_n : n \in \mathbb{N}\}$ solves (1.9). Given the first step, it is the taking of the low δ limit in (4.5) which will yield the second. However, to successfully carry out this limit, an extra piece of information will be needed, namely, for each fixed $n \in \mathbb{N}$, the uniform integrability of the family $f_n * \eta^\delta$ as δ ranges over $(0, 1)$. The next proposition is sufficient in this regard.

Proposition 4.4 *There exists a collection $\{k_n : n \in \mathbb{N}\}$ of positive constants such that, whenever ν is a vague limit point of $\{\mu_N^f : N \in \mathbb{N}\}$, the densities $\{f_n : n \in \mathbb{N}\}$ provided by Proposition 4.3 satisfy $\|f_n\|_{L^\infty(\mathbb{R}^d \times [0, \infty))} \leq k_n$ for each $n \in \mathbb{N}$.*

We now confirm that these two elements applied to (4.5) are enough to yield Theorem 1.1.

4.2.3 Taking the final step: Proof of Theorem 1.1.

By Proposition 4.3, (4.5) holds with f_n in place of ν_n for each $n \in \mathbb{N}$. The uniform boundedness provided by Proposition 4.4 and the Lebesgue differentiation theorem imply that $f_n * \eta^\delta$ converges pointwise to f_n a.e. on $\mathbb{R}^d \times [0, \infty)$ for each $n \in \mathbb{N}$. This same uniform boundedness and the dominated convergence theorem imply that each integral appearing in the identity converges to its counterpart where $f_n * \eta^\delta$ is replaced by f_n . Since the error term vanishes in the low δ limit, we indeed obtain (1.9). \square

5 An outline of the proof of the Stosszahlansatz

Here we explain in outline how we will prove Proposition 4.1.

5.1 Coagulation propensity, and particle pairs at small macroscopic distance

For $z \in \mathbb{R}^d$ and $n, m \in \mathbb{N}$, define under the law \mathbb{P}_N the stochastic process $Q_z = Q_{z,n,m} : [0, \infty) \rightarrow \mathbb{R}$ whose value at time $t \in [0, \infty)$ is given by

$$(5.1) \quad \frac{1}{2} \epsilon^{d-2} \sum_{i,j \in I_q(t)} \alpha(m_i, m_j) V_\epsilon(x_i - x_j + z) J_n(x_i, t) \mathbf{1}_{m_i(t)=n, m_j(t)=m}.$$

In seeking to prove the Stosszahlansatz, then, it is our aim to show that $\mathbb{E}_N \int_0^T Q_0(t) dt$ is close to a $\beta(n, m)$ multiple of the time-integrated product of microscopic candidate densities for particles of mass n and m . Since the microscopic coagulation density is binary in nature, it is unsurprising that $\mathbb{E}_N \int_0^T Q_0(t) dt$, the cumulative rate of coagulation between pairs of particles of such masses (at least if $J_n = 1$), should be approximated by the time integral of such a product of empirically defined densities. We have already explained heuristically in Section 3 why we might expect the macroscopic coagulation propensity $\beta(n, m)$ to have the form (1.8). The challenge now is to find a rigorous means of approximating $\mathbb{E}_N \int_0^T Q_0(t) dt$; as we outline this approach, we will see an alternative explanation for the formula (1.8) emerge.

Consider for a moment the expression Q_z , where $z \in \mathbb{R}^d$ is a small macroscopic quantity, namely, it is fixed at a given small value as we take a high N (or low ϵ) limit. We see that the quantity $\mathbb{E}_N \int_0^T Q_z(t) dt$ is a time-weighted count all instances of pairs of particles, of mass n and mass m , which at some time during $[0, T]$ are such that the mass m particle lies in the tiny ϵ -ball whose centre is displaced from the mass n particle by the small quantity z . Such instances at any given moment of time are weighted by the factor $\alpha(m_i, m_j) V_\epsilon(x_i - x_j + z)$; for later convenience, it is useful to also define \hat{Q}_z , where in the formula for Q_z , we replace $V_\epsilon(x_i - x_j + z)$

by $\hat{V}_\epsilon(x_i - x_j + z)$. Here, $\hat{V}_\epsilon(\cdot) = \epsilon^{-2}\hat{V}(\cdot/\epsilon)$, where $\hat{V} : \mathbb{R}^d \rightarrow [0, \infty)$ is a smooth and compactly supported function. (In fact, we will use two such variants, also writing \bar{Q}_z when \bar{V}_ϵ replaces V_ϵ .) The quantity $\mathbb{E}_N \int_0^T \hat{Q}_z(t) dt$ qualitatively meets the same description as does $\mathbb{E}_N \int_0^T Q_z(t) dt$, a time-weighted count of instances of z -displacements of particle pairs. The next assertion shows that $\mathbb{E}_N \int_0^T Q_0(t) dt$ is well approximated by an appropriately weighted count of particle pairs at small macroscopic distance z :

Proposition 5.1 *For $n, m \in \mathbb{N}$, recall that $u_{n,m} : \mathbb{R}^d \rightarrow [0, 1]$ is specified in (1.7), and let $\bar{V} = V(1 - u_{n,m})$. Then*

$$(5.2) \quad \int_0^T Q_0(t) dt = \int_0^T \bar{Q}_z(t) dt + \text{Err}_{n,m}(\epsilon, z).$$

Regarding the error term: defining Err_δ to be the supremum over $z \in \mathbb{R}^d$ for which $|z| = \delta$ of

$$\limsup_{\epsilon \searrow 0} \sum_{m \in \mathbb{N}} \mathbb{E}_N |\text{Err}_{n,m}(\epsilon, z)|,$$

we have that $\text{Err}_\delta \rightarrow 0$ as $\delta \rightarrow 0$.

From Proposition 5.1, it is simple to derive the Stosszahlansatz Proposition 4.1:

Sketch of proof of Proposition 4.1. We verify the statement only in the case $J_n = 1$; the general case invokes a simple additional estimate.

To derive Proposition 4.1, we begin by averaging the information in Proposition 5.1 over small macroscopic δ . In what follows, the relation $f \simeq_n g$ asserts that $f(n, m, \delta)$ and $g(n, m, \delta)$ are random variables on \mathbb{P}_N such that, for each $n \in \mathbb{N}$,

$$\lim_{\delta \searrow 0} \limsup_{\epsilon \searrow 0} \sum_{m \in \mathbb{N}} \mathbb{E}_N |f_{n,m,\delta} - g_{n,m,\delta}| = 0.$$

For $\delta > 0$, write $\eta_\delta : \mathbb{R}^d \rightarrow [0, \infty)$ for $\eta_\delta = \delta^{-d}\eta(\cdot/\delta)$. Proposition 5.1, the definition of \bar{Q} and substitutions $x_i - z_1 = \omega_1$ and $x_j - z_2 = \omega_2$ imply then that

$$\begin{aligned} & \int_0^T Q_0(t) dt \\ \simeq_n & \int_0^T \int_{\mathbb{R}^d \times \mathbb{R}^d} \bar{Q}_{z_2 - z_1}(t) \eta_\delta(z_1) \eta_\delta(z_2) dz_1 dz_2 dt \\ = & \epsilon^{d-2} \alpha(n, m) \int_0^T \int_{\mathbb{R}^d \times \mathbb{R}^d} \sum_{i,j \in I_q(t)} \bar{V}_\epsilon((x_i - z_1) - (x_j - z_2)) \mathbf{1}_{m_i(t)=n, m_j(t)=m} \eta_\delta(z_1) \eta_\delta(z_2) dz_1 dz_2 dt \\ = & \epsilon^{d-2} \alpha(n, m) \int_0^T \int_{\mathbb{R}^d \times \mathbb{R}^d} \sum_{i,j \in I_q(t)} \bar{V}_\epsilon(\omega_1 - \omega_2) \mathbf{1}_{m_i(t)=n, m_j(t)=m} \eta_\delta(x_i - \omega_1) \eta_\delta(x_j - \omega_2) d\omega_1 d\omega_2 dt. \end{aligned}$$

The virtue of this last expression is that the two particle sums may be passed inside to yield the microscopic candidate densities. Indeed, the expression equals

$$(5.3) \quad \epsilon^{d-2} \alpha(n, m) \int_0^T \int_{\mathbb{R}^d \times \mathbb{R}^d} \bar{V}_\epsilon(\omega_1 - \omega_2) f_n^{\epsilon, \delta}(\omega_1, t) f_m^{\epsilon, \delta}(\omega_2, t) d\omega_1 d\omega_2 dt.$$

Note that the double integral in (ω_1, ω_2) is almost on the diagonal, because \bar{V}_ϵ is supported in the ϵ -ball. When $\omega_1, \omega_2 \in \mathbb{R}^d$ satisfy $\|\omega_1 - \omega_2\| \leq \epsilon$, we have that $|\eta(\frac{x_i - \omega_2}{\delta}) - \eta(\frac{x_i - \omega_1}{\delta})| \leq \epsilon \delta^{-1} \|\nabla \eta\|_\infty$, so that $|f_n^{\epsilon, \delta}(\omega_2, t) - f_n^{\epsilon, \delta}(\omega_1, t)| \leq \epsilon^{d-1} \delta^{-d-1} \|\nabla \eta\|_\infty$.

Thus, at the expense of an error that is small in the sense of the \simeq_n relation, $f_m^{\epsilon, \delta}(\omega_2, t)$ may be replaced by $f_m^{\epsilon, \delta}(\omega_1, t)$ in (5.3); this done, the ω_2 integral may be detached, so that we see that (5.3) satisfies

$$\simeq_n \epsilon^{d-2} \alpha(n, m) \int_{\mathbb{R}^d} \bar{V}_\epsilon(x) dx \int_0^T \int_{\mathbb{R}^d} f_n^{\epsilon, \delta}(\omega, t) f_m^{\epsilon, \delta}(\omega, t) d\omega dt,$$

which since $\bar{V}_\epsilon(\cdot) = \epsilon^{-2} \bar{V}(\cdot/\epsilon)$ equals

$$\epsilon^{2(d-2)} \alpha(n, m) \int_{\mathbb{R}^d} \bar{V}(x) dx \int_0^T \int_{\mathbb{R}^d} f_n^{\epsilon, \delta}(\omega, t) f_m^{\epsilon, \delta}(\omega, t) d\omega dt,$$

Combining the above estimates, we confirm that Proposition 4.1 holds with

$$\beta(n, m) = \alpha(n, m) \int_{\mathbb{R}^d} \bar{V}(x) dx. \quad \square$$

5.2 An outline of the proof of Proposition 4.1

In an attempt to find a convenient representation of the quantity $\int_0^T Q_z(t) dt$, both when $z \in \mathbb{R}^d$ is zero and when it is non-zero and small, we define a z -dependent random variable S_z under \mathbb{P}_N for which the action of the free motion operator \mathfrak{A}_F on S_z produces, among others, the term Q_z . For each pair $(n, m) \in \mathbb{N}^2$, we define $\phi_{n, m}^\epsilon : \mathbb{R}^d \rightarrow (0, \infty)$ so that

$$-\Delta \phi_{n, m}^\epsilon(x) = \frac{\alpha(n, m)}{d(n) + d(m)} \epsilon^{-d} V(z/\epsilon),$$

subject to $\lim_{x \rightarrow \infty} \phi_{n, m}^\epsilon(x) = 0$. We then define a positively valued stochastic process $S_z : [0, \infty) \rightarrow (0, \infty)$ on \mathbb{P}_N : for each $t \geq 0$, we set

$$(5.4) \quad S_z(t)(q) = \epsilon^{2(d-2)} \sum_{i, j \in I_q(t)} \phi_{n, m}^\epsilon(x_i - x_j + z) J_n(x_i, t) \mathbf{1}\{m_i = n, m_j = m\}.$$

The action $\mathfrak{A}_F(S_z - S_0)$ of the free motion operator on $S_z - S_0$ is itself a random variable on \mathbb{P}_N which maps non-negative time $t \in [0, \infty)$ to \mathbb{R} . The term $Q_z(t) - Q_0(t)$ appears in the expression $-\mathfrak{A}_F(S_z - S_0)(t)$, in the case where each of the derivatives in the Laplacian operator falls on $\phi_{n, m}^\epsilon$ rather than on the test function J_n . (Note the minus sign attached to $\mathfrak{A}_F(S_z - S_0)(t)$; it arises from our choice that the functional S_z be positive rather than negative.)

For $T > 0$, consider the \mathbb{P}_N -almost sure identity

$$(5.5) \quad (S_z - S_0)(T) = (S_z - S_0)(0) + \int_0^T \left(\frac{\partial}{\partial t} + \mathfrak{A}_F \right) (S_z - S_0)(t) dt \\ + \int_0^T \mathfrak{A}_C(S_z - S_0)(t) dt + M_T,$$

and note that the process $\{M_T : T \geq 0\}$ is a \mathbb{P}_N -martingale. As we have noted, each of the terms $-\int_0^T Q_z(t)dt$ and $\int_0^T Q_0(t)dt$ appears in the free motion term on the right-hand side. The quantity $\int_0^T Q_0(t)dt$ remains of unit order in the low ϵ limit, as we discussed after the statement of the Stosszahlansatz Proposition 4.1. For similar reasons, $\int_0^T Q_z(t)dt$ may be expected to have this property for any given $z \in \mathbb{R}^d$. Suppose for a moment that it were the case that all the other terms appearing in (5.5) were of smaller order, as a low ϵ and then low z limit is taken. More precisely, suppose that, after the removal of the two terms mentioned above, the remainder satisfies the estimate on the error $\text{Err}_{n,m}$ given in Proposition 5.1. Then in fact Proposition 5.1 would hold, but with the term \bar{V} on the right-hand side of (5.2) replaced by V . Reviewing the proof of Proposition 4.1 from Proposition 5.1, we would find that the Stosszahlansatz indeed holds, but with the formula $\beta = \alpha \int_{\mathbb{R}^d} V(x)dx$ rather than (1.8). In other words, the reasoning that there are no further unit order terms appearing in (5.5) – which the authors of [10] believed when first studying this approach – leads to the fallacious guess $\beta = \alpha \int_{\mathbb{R}^d} V(x)dx$.

The formula's incorrectness means that we should expect some further term in (5.5) to remain of unit order as $\epsilon \rightarrow 0$ and then $z \rightarrow 0$. In Section 3.2, it was explained that this guess is wrong because of an effect of the curtailment of the interaction clock associated with a particle pair at the moment of the concerned particle's collision.

The further unit order term in (5.5) does indeed exist, and its form reflects this mechanism of curtailment of the interaction clock on particle collision. The term arises from the action of the collision operator \mathfrak{A}_C on S_0 . In the expression $\int_0^T \mathfrak{A}_C S_0(t)dt$, when the form of the collision operator \mathfrak{A}_C and of the functional S_0 is substituted, a sum is obtained. For each summand, four particles are concerned, in two pairs: two particles in the first pair arise from \mathfrak{A}_C , and it is this pair whose infinitesimal interaction rate is being integrated over the time period $[0, T]$, while the difference of the locations of the particles in the second pair form the argument of $\phi_{n,m}^\epsilon$ arising from S_0 . Although each of the two pairs is formed of distinct particles, there may be one or two coincidences between members of the first and of the second pair. When both of these coincidences occur, and the second pair equals the first, the contribution made to $-\int_0^T \mathfrak{A}_C S_0(t)dt$ by such terms is given by

$$(5.6) \quad \epsilon^{2(d-2)} \alpha(n, m) \int_0^T \sum_{k,l \in I_q(t)} V_\epsilon(x_k - x_l) \phi_{n,m}^\epsilon(x_k - x_l) J_n(x_k, t) dt.$$

This term witnesses the abrupt curtailment of the propensity to coagulate of a pair of particles at that moment when they do coagulate. That it is this term which remains of unit order reflects the role of the microscopic repulsion about a given particle in determining the relation (1.8) which is discussed in Section 3.

It turns out that the sum of the remaining terms in (5.5) is indeed negligible in that it satisfies the estimate that $\text{Err}_{n,m}(\epsilon, z)$ does in Proposition 4.1. Only the three unit order terms identified above remain in the limit of low ϵ and then z . That is, we have found that $\int_0^T Q_z(t) dt$ differs from

$$(5.7) \quad \epsilon^{2(d-2)} \alpha(n, m) \int_0^T \sum_{i,j \in I_q(t)} V_\epsilon(x_i - x_j) \\ \times \left[1 + \phi_{n,m}^\epsilon(x_i - x_j) \right] J_n(x_i, t) \mathbf{1}\{m_i = n, m_j = m\} dt.$$

by an error of the form $\text{Err}_{n,m}(\epsilon, z)$ in Proposition 4.1. Note that the ‘1’ that appears in the square bracket corresponds to $Q(0)$, and the other term to the unit order term (5.6). In the language of Proposition 5.1, we have learnt that

$$(5.8) \quad \int_0^T \hat{Q}_0(t) dt = \int_0^T Q_z(t) dt + \text{Err}_{n,m}(\epsilon, z),$$

where $\hat{V} = V(1 + \phi_{n,m})$ and \hat{Q}_0 is defined by the formula (5.1) that specifies Q_0 with \hat{V} replacing V . As a check of working, note that, since $\phi_{n,m} \geq 0$, we are asserting that the positive $\int_0^T Q_z(t) dt$ exceeds the positive $\int_0^T Q_0(t) dt$ by a further positive term of the same order. This is consistent with the explanation offered in Section 3.2: we expect $\int_0^T Q_z(t) dt$ to exceed $\int_0^T Q_0(t) dt$, because the size of the latter term (measuring the cumulative interaction clock of ϵ -displaced particles) is limited by the disappearance of particles on collision, while the former (measuring a comparable quantity for the much more distant z -displaced particles) experiences no such limitation.

Of course, (5.8) is not quite the conclusion we sought: to prove Proposition 5.1, we want to approximate $\int_0^T Q_0(t)$ by $\int_0^T \bar{Q}_z(t)$, so that the modification $Q \rightarrow \bar{Q}$ falls in the z -displaced term; but so far we have obtained such a result where the modification is made to the $z = 0$ term.

In light of this analysis, we may however revisit the approach. Consider a variant X_z of the process S_z : for each $z \in \mathbb{R}^d$, under \mathbb{P}_N , $X_z : [0, \infty) \rightarrow \mathbb{R}$ is given by

$$(5.9) \quad X_z(t)(q) = \epsilon^{2(d-2)} \sum_{i,j \in I_q(t)} u_{n,m}^\epsilon(x_i - x_j + z) J_n(x_i, t) \mathbf{1}\{m_i = n, m_j = m\},$$

where here, for each pair $(n, m) \in \mathbb{N}^2$, we define $u_{n,m}^\epsilon : \mathbb{R}^d \rightarrow (0, \infty)$ so that, for $z \in \mathbb{R}^d$,

$$(5.10) \quad -\Delta u_{n,m}^\epsilon(z) = \frac{\alpha(n,m)}{d(n)+d(m)} \epsilon^{-d} U(z/\epsilon),$$

subject to $\lim_{x \rightarrow \infty} u_{n,m}^\epsilon(x) = 0$. The function $U : \mathbb{R}^d \rightarrow (0, \infty)$ is at yet unspecified; of course the choice $U = V$ would specify the earlier functional S_z . Our aim now is to make a different choice of U , for which the solution of the problem (5.10) exists uniquely, and for which the earlier analysis may be carried out in such a way that its conclusion is not (5.8) but rather the desired

$$(5.11) \quad \int_0^T Q_0(t) dt = \int_0^T \bar{Q}_z(t) dt + \text{Err}_{n,m}(\epsilon, z),$$

for some $\bar{Q} : \mathbb{R}^d \rightarrow [0, \infty)$. Regarding scaling, note that, whenever $U : \mathbb{R}^d \rightarrow (0, \infty)$ is such that (5.10) has a unique solution for some $\epsilon > 0$, and this holds in fact for all $\epsilon > 0$; indeed, writing $u_{n,m} : \mathbb{R}^d \rightarrow (0, \infty)$ for $u_{n,m} = u_{n,m}^1$, we have that, for each $\epsilon > 0$, and for all $z \in \mathbb{R}^d$,

$$(5.12) \quad u_{n,m}^\epsilon(z) = \epsilon^{2-d} u_{n,m}(z/\epsilon).$$

In order to find a candidate for U that may make this plan work, we may hope that, for some suitable class of U , the earlier discussion continues to apply to the extent that the unit order terms that survive in the passage of low ϵ and then low z are the natural counterparts to the three terms identified there.

We want the analogue of the term $\int_0^T \hat{Q}_0(t)$ to be $\int_0^T Q_0(t)$ in the new calculation. Recall that $\hat{Q}_0(t)$ equals (5.7). When we reprise the earlier discussion with $X_z - X_0$ in place of $S_z - S_0$, the counterpart of the expression (5.7) is

$$(5.13) \quad \epsilon^{2(d-2)} \sum_{i,j \in I_q(t)} \left[- (d(n) + d(m)) \Delta u_{n,m}^\epsilon(x_i - x_j) + \alpha(n, m) V_\epsilon(x_i - x_j) u_{n,m}^\epsilon(x_i - x_j) \right] \\ J_n(x_i, t) \mathbf{1}\{m_i = n, m_j = m\}.$$

By (5.10), and the scaling properties $V_\epsilon(\cdot) = \epsilon^{-2} V(\cdot/\epsilon)$ and (5.12) of V and $u_{n,m}$, the quantity in the square brackets above equals

$$\epsilon^{-d} \alpha(n, m) \left(U\left(\frac{x_i - x_j}{\epsilon}\right) + V\left(\frac{x_i - x_j}{\epsilon}\right) u_{n,m}\left(\frac{x_i - x_j}{\epsilon}\right) \right).$$

Our aim is that (5.13) will equal $Q_0(t)$; we see that this demand is equivalent to the identity $U + V u_{n,m} = V$. That is, the function U must be specified by $U = V(1 - u_{n,m})$. Assuming for now that such a choice may be made, consider the term which is analogous to $\int_0^T Q_z(t) dt$ in (5.8) when the earlier analysis is replayed with X_z in place of S_z . This new term equals

$$(5.14) \quad -\epsilon^{2(d-2)} \int_0^T \sum_{i,j \in I_q(t)} (d(n) + d(m)) \Delta u_{n,m}^\epsilon(x_i - x_j + z) J_n(x_i, t) \mathbf{1}\{m_i(t) = n, m_j(t) = m\} dt.$$

Recalling the definition of \bar{Q} from the statement of Proposition 5.1, and noting that $-(d(n) + d(m)) \Delta u_{n,m}^\epsilon(z) = \alpha(n, m) V(z/\epsilon) (1 - u_{n,m}^\epsilon(z))$ for $z \in \mathbb{R}^d$, we see that (5.14) is precisely $\int_0^T \bar{Q}_z(t) dt$.

That is, setting U as described above, our reprisal of the method yields

$$(5.15) \quad \int_0^T Q_0(t) dt = \int_0^T \bar{Q}_z(t) dt + \text{Err}_{n,m}(\epsilon, z),$$

in place of (5.8), which is precisely the form of the statement asserted by Proposition 5.1.

To turn these ideas into a proof of Proposition 5.1, note first that making use of our desired choice of U entails that we argue that the PDE

$$-\Delta u_{n,m} = \frac{\alpha(n,m)}{d(n)+d(m)} V(1 - u_{n,m})$$

has a unique solution $u_{n,m} : \mathbb{R}^d \rightarrow [0, 1)$ satisfying $u_{n,m}(z) \rightarrow 0$ as $z \rightarrow \infty$. This we have already taken care of: see Lemma 1.4.

Our more substantial remaining task is the following. Defining the functional X_z with this choice of $u_{n,m}$, we must argue that the dominant terms in the identity (5.5) (with X_z in place of S_z) are indeed $-\int_0^T (\overline{Q}_z(t) - Q_0(t)) dt$; more precisely, we must show that both the left-hand side of (5.5), and the difference of its right-hand side with $-\int_0^T (\overline{Q}_z(t) - Q_0(t)) dt$, satisfy the demand made of the error $\text{Err}_{n,m}(\epsilon, z)$ in the statement of Proposition 5.1.

6 Proof of Proposition 5.1

We now present the proof of the Stosszahlansatz, or rather, reduce it to two key estimates (and we do so making some simplifications which in no way diminish the essentials of the argument). The job at hand is to carry out the task mentioned in the preceding paragraph. In order to analyse the various error terms, we begin by providing formulas for them.

6.1 The action of the free motion and collision operators on the functional

Recall that, for each $z \in \mathbb{R}^d$, under \mathbb{P}_N , we are defining $X_z : [0, \infty) \rightarrow \mathbb{R}$ by means of

$$(6.1) \quad X_z(t)(q) = \epsilon^{2(d-2)} \sum_{i,j \in I_q(t)} u_{n,m}^\epsilon(x_i - x_j + z) J_n(x_i, t) \mathbf{1}\{m_i = n, m_j = m\},$$

where, for each pair $(n, m) \in \mathbb{N}^2$, we define $u_{n,m}^\epsilon : \mathbb{R}^d \rightarrow (0, \infty)$ so that, for $z \in \mathbb{R}^d$,

$$-\Delta u_{n,m}^\epsilon(z) = \epsilon^{-d} \frac{\alpha(n,m)}{d(n)+d(m)} V(z/\epsilon) (1 - u_{n,m}(z/\epsilon))$$

subject to $\lim_{x \rightarrow \infty} u_{n,m}^\epsilon(x) = 0$.

Our aim is to analyse the high- N behaviour of the terms in the \mathbb{P}_N -almost sure identity

$$(6.2) \quad (X_z - X_0)(T) = (X_z - X_0)(0) + \int_0^T \left(\frac{\partial}{\partial t} + \mathfrak{A}_F \right) (X_z - X_0)(t) dt \\ + \int_0^T \mathfrak{A}_C(X_z - X_0)(t) dt + M_T,$$

where the process $\{M_T : T \geq 0\}$ is a \mathbb{P}_N -martingale.

To simplify our presentation, we will consider only the case that the test function $J : \mathbb{R}^d \times \mathbb{N} \times [0, \infty) \rightarrow [0, \infty)$ takes the form $J(x, m', t) = J_n \mathbf{1}_{m'=n}$, where $J_n \in (0, \infty)$ is a constant; in the general case, $J_n : \mathbb{R}^d \times [0, \infty) \rightarrow [0, \infty)$ is a map on space-time. We also write $\overline{J} : \mathbb{R}^d \times \mathbb{N} \times [0, \infty) \rightarrow [0, \infty)$ for the function $\overline{J}(x, m', t) = \mathbf{1}_{m'=m}$. In this way, we may write

$$(6.3) \quad X_z(t)(q) = \epsilon^{2(d-2)} \sum_{i,j \in I_q(t)} u_{n,m}^\epsilon(x_i - x_j + z) J(x_i, m_i, t) \overline{J}(x_j, m_j, t).$$

We may label terms arising from the action of the free motion operator on the functional as follows:

$$(6.4) \quad \left(\frac{\partial}{\partial t} + \mathfrak{A}_F\right)(X_z - X_0)(t) = H_1 + H_2 + H_3,$$

where

$$\begin{aligned} H_1 &= -\epsilon^{2(d-2)} \sum_{i,j \in I_q} \alpha(m_i, m_j) \left[V^\epsilon(x_i - x_j + z) - V^\epsilon(x_i - x_j) \right] J(x_i, m_i, t) \bar{J}(x_j, m_j, t), \\ H_2 &= -\epsilon^{2(d-2)} \sum_{i,j \in I_q} \alpha(m_i, m_j) V_\epsilon(x_i - x_j) u_{m_i, m_j}^\epsilon(x_i - x_j) J(x_i, m_i, t) \bar{J}(x_j, m_j, t), \\ H_3 &= \epsilon^{2(d-2)} \sum_{i,j \in I_q} \alpha(m_i, m_j) V_\epsilon(x_i - x_j + z) u_{m_i, m_j}^\epsilon(x_i - x_j + z) J(x_i, m_i, t) \bar{J}(x_j, m_j, t). \end{aligned}$$

The terms arising from the collision operator may be labelled

$$\mathfrak{A}_C(X_z - X_0)(t) = G_z(1) + G_z(2) - G_0(1) - G_0(2),$$

where $G_z(1)$ equals

$$(6.5) \quad \begin{aligned} &\frac{1}{2} \sum_{k,l \in I_q} \alpha(m_k, m_l) V_\epsilon(x_k - x_l) \epsilon^{2(d-2)} \sum_{i \in I_q} \\ &\left\{ \frac{m_k}{m_k + m_l} \left[u_{m_k, m_i}^\epsilon(x_k - x_i + z) J(x_k, m_k + m_l, t) \bar{J}(x_i, m_i, t) \right. \right. \\ &\quad \left. \left. + u_{m_i, m_k}^\epsilon(x_i - x_k + z) J(x_i, m_i, t) \bar{J}(x_k, m_k + m_l, t) \right] \right. \\ &+ \frac{m_l}{m_k + m_l} \left[u_{m_l, m_i}^\epsilon(x_l - x_i + z) J(x_l, m_k + m_l, t) \bar{J}(x_i, m_i, t) \right. \\ &\quad \left. \left. + u_{m_i, m_l}^\epsilon(x_i - x_l + z) J(x_i, m_i, t) \bar{J}(x_l, m_k + m_l, t) \right] \right. \\ &- \left[u_{m_k, m_i}^\epsilon(x_k - x_i + z) J(x_k, m_k, t) \bar{J}(x_i, m_i, t) \right. \\ &\quad \left. \left. + u_{m_i, m_k}^\epsilon(x_i - x_k + z) J(x_i, m_i, t) \bar{J}(x_k, m_k, t) \right] \right. \\ &- \left[u_{m_l, m_i}^\epsilon(x_l - x_i + z) J(x_l, m_l, t) \bar{J}(x_i, m_i, t) \right. \\ &\quad \left. \left. + u_{m_i, m_l}^\epsilon(x_i - x_l + z) J(x_i, m_i, t) \bar{J}(x_l, m_l, t) \right] \right\}, \end{aligned}$$

and where

$$(6.6) \quad G_z(2) = -\epsilon^{2(d-2)} \sum_{k,l \in I_q} \alpha(m_k, m_l) V_\epsilon(x_k - x_l) u_{m_k, m_l}^\epsilon(x_k - x_l + z) J(x_k, m_k, t) \bar{J}(x_l, m_l, t).$$

In the triple sum over distinct particle indices (k, l, i) appearing in $G_z(1)$, the particles indexed by k and l are interacting at rate $\alpha(m_k, m_l) V_\epsilon(x_k - x_l)$; when this pair collides, there is an instantaneous change in the value of those terms in $X_z(q)$ that include the location x_i of a given third

particle indexed by i and not involved in the collision. There are two gain terms, associated to the appearance of a new particle at one or other of x_k and x_l , and two loss terms, associated to the disappearance of the particles indexed by k and l .

The term $G_z(2)$ is a double sum over distinct particle indices (k, l) that records the instantaneous change caused by collision of such a particle pair in the value of those terms in $X_z(q)$ expressed in terms only of the elements of that pair. That is, the collision occurs at infinitesimal rate $\alpha(m_k, m_l)V_\epsilon(x_k - x_l)$; when collision happens, the particles with indices k and l disappear, so that the term $\epsilon^{2(d-2)}u_{m_k, m_l}^\epsilon(x_k - x_l + z)J_n \mathbf{1}_{m_j=m}$ no longer appears in $X_z(q)$.

Note that

$$H_2 = G_0(2).$$

We find then that

$$(6.7) \quad \begin{aligned} & \left| \int_0^T H_1(t) dt + \int_0^T H_3(t) dt \right| \\ & \leq |X_z - X_0|(q(T)) + |X_z - X_0|(q(0)) \\ & \quad + \int_0^T |G_z(1) - G_0(1)|(t) dt + \int_0^T |G_z(2)|(t) dt + |M(T)|. \end{aligned}$$

We will now state bounds on these error terms which are sufficient for the purpose of proving Proposition 5.1 (and thus Proposition 4.1 and Theorem 1.1).

Proposition 6.1 *There exists a constant $C > 0$ such that*

1. for all $N \in \mathbb{N}$, $\sum_{m \in \mathbb{N}} \int_0^T \mathbb{E}_N |G_z(1) - G_0(1)|(t) dt \leq CT^{3d/2}|z|(\log 1/|z|)^{3d/2}$;
2. for any $t \geq 0$, $\sum_{m \in \mathbb{N}} \mathbb{E}_N |X_z(t) - X_z(0)| \leq C|z|$;
3. for all $N \in \mathbb{N}$, $\sum_{m \in \mathbb{N}} \int_0^T \mathbb{E}_N |G_z(2)|(t) dt \leq C(\frac{\epsilon}{z})^{d-2}$;
4. and, for each $t \geq 0$, $\sum_{m \in \mathbb{N}} \mathbb{E}_N [M(t)^2] \leq C\epsilon^{d-2}$.

6.2 Proving the error bounds

Two important tools are needed to prove the above bounds. Here, we present these two tools (but do not yet prove the assertions we state about them), and use them to give a proof of Proposition 6.1(1); the three other estimates in this proposition follow in a roughly similar way, and we do not give the proofs of these estimates here.

6.2.1 Particle concentration bounds

The first tool is an assertion that, at any given time, the joint density of any given number of particles is uniformly bounded above. Recall that $\{h_n : n \in \mathbb{N}\}$ is the intensity density profile of particles under \mathbb{P}_N .

Proposition 6.2 For $k \in \mathbb{N}$, let $g_k : \mathbb{R}^{dk} \times [0, \infty) \rightarrow [0, \infty)$ be such that $g_k(y_1, \dots, y_k, t)$ is the density at $(y_1, \dots, y_k) \in (\mathbb{R}^d)^k$ for the ordered presence of the lowest- k indexed particles in \mathbb{P}_N at time t . Then

$$\|g_k\|_{L^\infty(\mathbb{R}^{dk} \times [0, \infty))} \leq \left(\sup_{m \geq 1} m^{-1} d(m)^{-d/2} \right)^k \cdot \sup_{j \in \mathbb{N}} \|h_j\|_{L^\infty(\mathbb{R}^d)}^k.$$

6.2.2 Uniform control on pairwise collision probabilities

Recall from Section 1.11 the notion of Brownian motion on \mathbb{R}^d killed at rate W , where $W : \mathbb{R}^d \rightarrow [0, \infty)$ is a smooth and compactly supported function; recall from there that $u_W : \mathbb{R}^d \rightarrow [0, 1]$ is such that $u_W(x)$ is the probability that Brownian motion killed at rate W and begun at x is killed at some time, and also that we set $u_W^\epsilon(\cdot) = \epsilon^{2-d} u_W(\cdot/\epsilon)$.

Lemma 6.3 There exist a constant $C > 0$ such that, for all continuous $W : \mathbb{R}^d \rightarrow [0, \infty)$ with support in the Euclidean unit ball,

- for $x \in \mathbb{R}^d$, $u_W(x) \leq \frac{1}{\|x\|^{d-2}}$ and $\|\nabla u_W(x)\| \leq \frac{C}{\|x\|^{d-1}}$;

- for $x, z \in \mathbb{R}^d$ and $\epsilon > 0$ such that $\|x\| \geq \max\{2\|z\| + \epsilon, 2\epsilon\}$,

$$(6.8) \quad \left| u_W^\epsilon(x+z) - u_W^\epsilon(x) \right| \leq 2^{3d-6} \frac{\|z\|}{\|x\|^{d-1}}$$

and

$$(6.9) \quad \left| \nabla u_W^\epsilon(x+z) - \nabla u_W^\epsilon(x) \right| \leq 4^d (2^{d-1}d + 1) \frac{\|z\|}{\|x\|^d}.$$

6.2.3 Applying the tools

For the proof of Proposition 6.1(1), we need one further simple estimate, on long-range particle displacement:

Lemma 6.4 Suppose that each element of the initial data $h_n : \mathbb{R}^d \rightarrow [0, \infty)$, $n \in \mathbb{N}$, is supported in a given compact region B , and that $\bar{d} := \sup_{n \in \mathbb{N}} d(n) < \infty$. Then, for some constant $C > 0$ and for all $r > 0$,

$$\mathbb{P}_N \left(|x_1(T)| \geq r \right) \leq C \exp \left\{ -\frac{r^2}{2\bar{d}T} \right\}.$$

Proof. If $R > 0$ is an upper bound on the radius of the region B , then note that $x_1(T)$ under \mathbb{P}_N is stochastically dominated by the maximum modulus during $[0, T]$ of a rate- \bar{d} Brownian motion begun at distance R from the origin. From this and the reflection principle the result follows. \square

Proof of Proposition 6.1(1). Note that

$$\int_0^T \mathbb{E}_N |G_z(1) - G_0(1)|(t) dt \leq \sum_{i=1}^8 D_i,$$

where

$$(6.10) \quad D_1 = \frac{1}{2} \mathbb{E}_N \int_0^T dt \sum_{k,l \in I_q} \alpha(m_k, m_l) V_\epsilon(x_k - x_l) |J(x_k, m_k)| \\ \epsilon^{2(d-2)} \sum_{i \in I_q} |\bar{J}(x_i, m_i)| \cdot \left| u_{m_k, m_i}^\epsilon(x_k - x_i + z) - u_{m_k, m_i}^\epsilon(x_k - x_i) \right|,$$

and the later D_i terms differ from D_1 only in inessential ways. Note that D_1 depends implicitly on n and m , in that the test functions J and \bar{J} have been chosen to charge only particles of mass n and m . The above sum over particle index triples (k, l, i) has at most $N^3 = Z^3 \epsilon^{3(2-d)}$ summands; recalling that $V_\epsilon = \epsilon^{-2} V(\cdot/\epsilon)$ is supported in the ϵ -ball about the origin, we see that $\sum_{m \in \mathbb{N}} D_1$ is at most

$$(6.11) \quad \epsilon^{3(2-d)} \cdot \epsilon^{-2} \cdot \epsilon^{2(d-2)} Z^3 C_1 \mathbb{E}_N \int_0^T \mathbf{1}_{|x_1 - x_2| \leq \epsilon} \cdot \sup_{m \in \mathbb{N}} \left| u_{n,m}^\epsilon(x_3 - x_2 + z) - u_{n,m}^\epsilon(x_3 - x_2) \right| \cdot \mu_t(dx_1, dx_2, dx_3),$$

where we set $C_1 = \|V\|_\infty \cdot \sup_{n,m} \alpha(n, m) \cdot \sup_n J_n \cdot \sup_m \bar{J}_m$, and where the law μ_t is the joint density at time t of the first three indexed particles under \mathbb{P}_N . By Lemma 6.4, and the uniform bound $u^\epsilon \leq 1$, the contribution to the right-hand side made by the integral over $(x_1, x_2, x_3) \in \mathbb{R}^{3d} \setminus [-R, R]^{3d}$ is at most $C \exp\left\{-\frac{R^2}{2\bar{d}T}\right\}$, which equals $|z|$ if we choose $R^2 = 2\bar{d}T \log(C/|z|)$. We bound the integrand on $[-R, R]^{3d}$ by noting that the support of V_ϵ has volume ϵ^d and using Lemma 6.3 to bound the functions $|u_{n,m}^\epsilon(x+z) - u_{n,m}^\epsilon(x)|$ simultaneously. Thus,

$$\sum_{m \in \mathbb{N}} D_1 \leq Z^3 C_1 \int_{[-R, R]^{3d}} \sup_{m \in \mathbb{N}} \left| u_{n,m}^\epsilon(x_3 - x_2 + z) - u_{n,m}^\epsilon(x_3 - x_2) \right| dx + |z| \\ \leq Z^3 \|V\|_\infty \cdot \sup_{n,m} \alpha(n, m) \cdot \sup_n J_n \cdot \sup_m \bar{J}_m \cdot (2\bar{d}T \log(C/|z|))^{3d/2} \cdot C|z| + |z|,$$

where it is (6.8) that determines the value of the constant $C > 0$. This completes the proof of Proposition 6.1(1). \square

6.3 Uniform control on pairwise collision probabilities: proofs

Proof of Lemma 6.3. By the uniqueness element of Lemma 1.6, we have that, for $x \in \mathbb{R}^d$,

$$(6.12) \quad u_W(x) = c_0 \int_{\mathbb{R}^d} W(y) (1 - u_W(y)) \|x - y\|^{2-d} dy,$$

where, for each $d \geq 3$, $c_0^{-1} = d(d-1)\omega_d$, with ω_d equal to the volume of the Euclidean unit ball in d dimensions.

Write $u_\infty : \mathbb{R}^d \rightarrow [0, 1]$ so that for $x \in \mathbb{R}^d$, $u_\infty(x)$ is the probability that Brownian motion begun at $x \in \mathbb{R}^d$ visits the Euclidean unit ball; our notation is used because formally this object coincides with u_W for $W = \infty \mathbf{1}_{\|x\| \leq 1}$. From the interpretation of u_W and u_∞ as killing probabilities, it is

evident that $u_W(x) \leq u_\infty(x)$ for all $x \in \mathbb{R}^d$ and for any continuous W supported in the unit ball. However, for $x \in \mathbb{R}^d$,

$$(6.13) \quad u_\infty(x) = \min \{1, \|x\|^{2-d}\};$$

thus, $u_W(x) \leq \|x\|^{2-d}$ for all $x \in \mathbb{R}^d$ and W as above, as Lemma 6.3 firstly asserts.

By using this monotonicity to compare the formulas (6.12) and (6.13) along a sequence of $x \in \mathbb{R}^d$ for which $x \rightarrow \infty$, we obtain that, for all such potentials W ,

$$(6.14) \quad \int_{\mathbb{R}^d} W(y)(1 - u_W(y)) \, dy \leq c_0^{-1}.$$

Note that

$$\nabla u_W(x) = c_0 \int_{\mathbb{R}^d} W(y) \frac{x-y}{\|x-y\|^d} (1 - u_W(y)) \, dy.$$

From (6.14), we see that $\|\nabla u_W(x)\| \leq c_0^{-1} \int_{\|y\| \leq 1} \|x-y\|^{1-d} \, dy \leq C \|x\|^{1-d}$ whenever $\|x\| \geq 2$, as we also asserted. On the other hand, that $\|\nabla u_W(x)\| \leq C$ when $\|x\| \leq 2$ is straightforward. We have obtained Lemma 6.3's second assertion.

Turning to derive (6.8), note that, for $x \in \mathbb{R}^d$,

$$u_W^\epsilon(x) = c_0 \epsilon^{2-d} \int_{\mathbb{R}^d} W_\epsilon(y)(1 - u_{W_\epsilon}(y)) \cdot \|x-y\|^{2-d} \, dy.$$

(For the rest of the proof, we will denote the Euclidean norm on \mathbb{R}^d by $|\cdot|$ rather than by $\|\cdot\|$)
Thus, for $x, z \in \mathbb{R}^d$,

$$\begin{aligned} & |u_W^\epsilon(x+z) - u_W^\epsilon(x)| = \epsilon^{2-d} |u_W(\frac{x+z}{\epsilon}) - u_W(\frac{x}{\epsilon})| \\ & \leq c_0 \int_{\mathbb{R}^d} W(y)(1 - u_W(y)) \left| |x+z-\epsilon y|^{2-d} - |x-\epsilon y|^{2-d} \right| \, dy \\ & \leq \sup_{|y| \leq \epsilon} \left| |x+z-y|^{2-d} - |x-y|^{2-d} \right| \\ & = \sup_{|y| \leq \epsilon} \frac{|x-y|^{d-2} - |x+z-y|^{d-2}}{|x+z-y|^{d-2} |x-y|^{d-2}}, \end{aligned}$$

the second inequality by (6.14).

Note that

$$(6.15) \quad |x+z-y|^{d-2} - |x-y|^{d-2} \leq \left(|x-y| + |z|\right)^{d-2} - |x-y|^{d-2}$$

and that

$$(6.16) \quad |x-y|^{d-2} - |x+z-y|^{d-2} \leq |x-y|^{d-2} - \left(|x-y| - |z|\right)^{d-2}.$$

The right-hand sides of (6.15) and (6.16) each take the form $\alpha^{d-3}|z|$, for some $\alpha \in [|x-y|-|z|, |x-y|+|z|]$. Note that if $|y| \leq \epsilon$, then

$$|x-y|-|z| \geq |x|-\epsilon-|z| \geq 0,$$

since $|x| \geq |z| + \epsilon$. As a result, we have that $|x-y|+|z| \leq 2|x-y|$, so that

$$\left| |x-y|^{d-2} - |x+z-y|^{d-2} \right| \leq 2^{d-3}|z||x-y|^{d-3}.$$

Hence,

$$\left| u_W^\epsilon(x+z) - u_W^\epsilon(x) \right| \leq 2^{d-3}|z| \sup_{|y| \leq \epsilon} |x-y|^{-1} |x+z-y|^{2-d}.$$

From $|x| \geq \max\{2|z| + \epsilon, 2\epsilon\}$ and $|y| \leq \epsilon$ follows $|x+z-y| \geq |x-y|/2$ and $|x-y| \geq |x|/2$; thus, the above supremum is at most $2^{2d-3}|x|^{1-d}$. We obtain (6.8).

In seeking to prove (6.9), note that

$$\frac{x+z-y}{|x+z-y|^d} - \frac{x-y}{|x-y|^d} = \frac{(x+z-y)|x-y|^d - (x-y)|x+z-y|^d}{|x+z-y|^d|x-y|^d}.$$

Note that, for any $a \in \mathbb{R}^d$,

$$(6.17) \quad \left| (a+z)|a|^d - |a+z|^d a \right| \leq |a| \left| |a+z|^d - |a|^d \right| + |z||a|^d \leq (2^{d-1}d+1)|z||a|^d,$$

as long as $|z| \leq |a|$. Given that

$$\nabla u_W^\epsilon(x) = -(d-2)c_0 \int_{\mathbb{R}^d} W(y)(1-u_W(y)) \frac{x-y}{|x-\epsilon y|^d} dy,$$

we may apply (6.17) with the choice $a = x - \epsilon y$ and then use (6.14) to obtain

$$\begin{aligned} & \left| \nabla u_W^\epsilon(x+z) - \nabla u_W^\epsilon(x) \right| \\ & \leq (2^{d-1}d+1)|z|c_0 \int_{\mathbb{R}^d} W(y)(1-u_W(y)) |x+z-\epsilon y|^{-d} dy \\ & \leq (2^{d-1}d+1)|z| \sup_{|y| \leq \epsilon} |x+z-y|^{-d}. \end{aligned}$$

From $|x| \geq \max\{2|z| + \epsilon, 2\epsilon\}$ and $|y| \leq \epsilon$, we see that $|x+z-y| \geq |x-y|/2$ and $|x-y| \geq |x|/2$. We conclude that

$$\left| \nabla u^\epsilon(x+z) - \nabla u^\epsilon(x) \right| \leq 4^d (2^{d-1}d+1) \frac{|z|}{|x|^d},$$

as required. □

7 Particle concentration bounds and uniform integrability

The aim of this section is to prove the particle concentration upper bound, Proposition 6.2, and the closely related uniform integrability assertions, Propositions 4.3 and 4.4.

Proposition 6.2 asserts that, if $d : \mathbb{N} \rightarrow (0, \infty)$ decreases, but not too rapidly, then supremum norm bounds enjoyed initially by the particle profile propagate to all later times up to factors determined by the diffusion rates d . Such particle concentration results play an essential role in our derivation of the Smoluchowski PDE; unlike in our preceding work, here we present a proof using probabilistic techniques. The proof will occupy several pages and invokes not unrestrictive hypotheses on $d(\cdot)$. We make use of this approach because of the attractive probabilistic perspective that it offers one of the more technical aspects of our kinetic limit derivation of the Smoluchowski PDE; and because the uniform integrability Proposition 4.4 is an immediate corollary.

It is much quicker to describe the supremum norm propagation effect in terms of solutions to the PDE. In order to illustrate the effect succinctly to begin with, and perhaps also for the benefit of analytically minded readers who may wish to skip some details in the upcoming proof of Proposition 6.2, we first present the statement and proof of [11, Lemma 4.1]. Such a reader may also wish to consult [31], where an analogous particle distribution result, Theorem 3.1, is proved by means not unlike, but more analytic than, our approach to establishing Proposition 6.2.

7.1 An analytic bound on particle concentration

Our analytic lemma concerns a weak solution $\{f_n : n \in \mathbb{N}\}$ of the system (1.9).

Lemma 7.1 *Assume $d(\cdot)$ is non-increasing. Then, for all $x \in \mathbb{R}^d$ and $t \geq 0$,*

$$(7.1) \quad \sum_{n=1}^{\infty} n d(n)^{d/2} f_n(x, t) \leq d(1)^{d/2} u(x, t).$$

where u is the unique solution to $u_t = d(1)\Delta u$ subject to the initial condition $u(x, 0) = \sum_{n=1}^{\infty} n f_n(x, 0)$.

Proof. For $D > 0$, let $\{S_t^D : t \geq 0\}$ denote the diffusion rate D heat semigroup. That is, for any continuous function $f : \mathbb{R}^d \rightarrow \mathbb{R}$, $S_t^D f : \mathbb{R}^d \rightarrow \mathbb{R}$ is given by

$$S_t^D f(x) = \int_{\mathbb{R}^d} f(x - y) \cdot \frac{1}{(2\pi Dt)^{d/2}} e^{-\frac{y^2}{2Dt}} dy.$$

The heat semigroup satisfies the property that, if $D_1 \geq D_2$ and $g \geq 0$, then

$$(7.2) \quad D_1^{d/2} S_t^{D_1} g \geq D_2^{d/2} S_t^{D_2} g;$$

this is a consequence of an elementary bound on the normal density, which in fact we will shortly state as (7.10).

Using the shorthand $Q_n(x, s) = Q_n(f)(x, s)$, $s \in [0, \infty)$, for the collision operator in (1.2), Duhamel's principle implies the basic relation that, for each $n \in \mathbb{N}$, and for all $(x, t) \in \mathbb{R}^d \times [0, \infty)$,

$$(7.3) \quad f_n(x, t) = S_t^{d(n)} h_n(x) + \int_0^t S_{t-s}^{d(n)} Q_n(x, s) ds,$$

where recall that $h_n = f_n(\cdot, 0)$ denotes the initial condition.

We will argue that, for each $\ell \in \mathbb{N}$, and for all $(x, t) \in \mathbb{R}^d \times [0, \infty)$,

$$(7.4) \quad \sum_1^\ell nd(n)^{d/2} f_n(x, t) \leq d(1)^{d/2} S_t^{d(1)} \left(\sum_1^\ell nh_n \right) (x) + d(\ell)^{d/2} \int_0^t S_{t-s}^{d(\ell)} \left(\sum_1^\ell nQ_n(x, s) \right) ds;$$

first, let us show that this claim proves the lemma.

Consider the expression $\sum_1^\ell nQ_n(x, t)$ for any $(x, t) \in \mathbb{R}^d \times [0, \infty)$. Coagulations at (x, t) between pairs of particles whose combined mass is at most ℓ , or each of whose masses are at least $\ell + 1$, do not contribute to the expression, while the remaining coagulations contribute negatively. Thus, $\sum_1^\ell nQ_n(x, t) \leq 0$. We find then from (7.4) that

$$(7.5) \quad \sum_1^\ell nd(n)^{d/2} f_n \leq d(1)^{d/2} S_t^{d(1)} \left(\sum_1^\ell nf_n^0 \right).$$

In this way, we see that, to prove the lemma, it suffices to derive (7.4).

We will establish this bound by induction on $\ell \in \mathbb{N}$. When $\ell = 1$, the bound holds as an equality due to (7.3).

Supposing that (7.4) is valid at index ℓ , we now derive it at index $\ell + 1$. By (7.2) and (7.4), we learn that, for all $(x, t) \in \mathbb{R}^d \times [0, \infty)$,

$$(7.6) \quad \sum_1^\ell nd(n)^{d/2} f_n(x, t) \leq d(1)^{d/2} S_t^{d(1)} \left(\sum_1^\ell nh_n \right) (x) + d(\ell + 1)^{d/2} \int_0^t S_{t-s}^{d(\ell+1)} \left(\sum_1^\ell nQ_n(x, s) \right) ds$$

because $d(\ell) \geq d(\ell + 1)$ and $\sum_1^\ell nQ_n(x, t) \leq 0$.

Applying (7.2) to (7.3) with $n = \ell + 1$ yields

$$(7.7) \quad f_{\ell+1}(x, t) \leq \left(\frac{d(1)}{d(\ell+1)} \right)^{d/2} S_t^{d(1)} h_{\ell+1}(x) + \int_0^t S_{t-s}^{d(\ell+1)} Q_{\ell+1}(x, s) ds.$$

We multiply both sides of (7.7) by $(\ell + 1)d(\ell + 1)^{d/2}$ and add the result to (7.6). The outcome is

$$\sum_1^{\ell+1} nd(n)^{d/2} f_n(x, t) \leq d(1)^{d/2} S_t^{d(1)} \left(\sum_1^{\ell+1} nh_n \right) (x) + d(\ell + 1)^{d/2} \int_0^t S_{t-s}^{d(\ell+1)} \left(\sum_1^{\ell+1} nQ_n(x, s) \right) ds.$$

This completes the proof. □

7.2 Proof of Proposition 6.2

We begin the proof by reformulating the proposition as Proposition 7.3 and proving this. First we record some notation concerning particle locations.

Definition 7.2 For $t \geq 0$, write $\mathcal{X}_t = \{x_i : i \in I_{q(t)}\}$ for the set of particle locations at time t . For $m \in \mathbb{N}$, further write $\mathcal{X}_t^m = \{x_i : i \in I_{q(t)}, m_i = m\}$ for the set of mass m particle locations at time t .

Recall that μ denotes Lebesgue measure on \mathbb{R}^d .

Proposition 7.3 Suppose that $d : \mathbb{N} \rightarrow (0, \infty)$ is non-increasing. For $m \in \mathbb{N}$, let $\ell_m = \|f_m(\cdot, 0)\|_{L^\infty(\mathbb{R}^d)}$. For any $T \geq 0$ and $A \subseteq \mathbb{R}^d$ open,

$$\mathbb{P}_N(|\mathcal{X}_T \cap A| = k) \leq N^k K^k \mu(A)^k.$$

where $K > 0$ denotes the constant $K = Z^{-1} \sum_{m_0 \geq 1} \ell_{m_0} m_0 d(m_0)^{d/2} \sup_{m \geq m_0} m^{-1} d(m)^{-d/2}$.

We begin by proving Proposition 7.3 in the case of one particle.

Lemma 7.4 Suppose that $d : \mathbb{N} \rightarrow (0, \infty)$ is non-increasing. Let $N \geq 1$. For any $T \geq 0$ and $A \subseteq \mathbb{R}^d$ open,

$$\mathbb{P}_N(\mathcal{X}_T \cap A \neq \emptyset) \leq NK\mu(A),$$

where, ℓ_{m_0} and K are specified in Proposition 7.3.

The proofs of Proposition 7.3 and Lemma 7.4 rely on the specific details of pairwise collision seen in the dynamics under \mathbb{P}_N which we defined in Section 1.3; the reader may wish to recall these now by consulting the paragraphs in Section 1.3 under the heading “the precise mechanism of collision”.

7.2.1 The method of proof: the tracer particle

The tool with which we will prove Proposition 7.3 and Lemma 7.4 is the *tracer particle*, which is simply a particle selected uniformly at random at time zero. This notion has been introduced already, *a propos* of the translation invariant model on a torus seen in Section 3. Note however that, in that case, particle symmetry meant that we might call a particle of any given index the tracer particle.

Definition 7.5 Under the law \mathbb{P}_N , let U denote a random variable having the uniform law on $\{1, \dots, N\}$, independently of other randomness. The tracer particle is the particle whose index is U . The location and mass of the tracer particle will be denoted by $x_{\text{tp}} : [0, \infty) \rightarrow \mathbb{R}^d \cup \{c\}$ and $m_{\text{tp}} : [0, \infty) \rightarrow \mathbb{N} \cup \{c\}$.

The basic relation between tracer particle location and the distribution of particles at any given time $T \geq 0$ is now stated.

Lemma 7.6 Fix $T \geq 0$. For each $A \subseteq \mathbb{R}^d$ open,

$$\mathbb{E}_N |\mathcal{X}_T \cap A| = N \mathbb{P}_N(x_{\text{tp}}(T) \in A),$$

Proof. Note that $\mathbb{E} |\mathcal{X}_T \cap A| = \mathbb{E} \sum_{i=1}^N \mathbf{1}_{x_i(T) \in A} = N \mathbb{P}_N(x_U(T) \in A)$. □

7.2.2 A stronger inductive hypothesis

We will prove Lemma 7.4 by formulating a stronger inductive hypothesis where the parameter for the induction is the initial total particle number N . As a shorthand, we write

$$(7.8) \quad \nu_{x,s}(\mathrm{d}y) = (2\pi s)^{-d/2} \exp\left\{-\frac{\|x-y\|^2}{2s}\right\} \mathrm{d}y$$

for the law of a normal random variable of mean $x \in \mathbb{R}^d$ and variance $s \geq 0$.

Lemma 7.7 *Let $N \geq 1$. For given $x \in \mathbb{R}^d$, $m_0 \in \mathbb{N}$ and $\chi \in (\mathbb{R}^d \times \mathbb{N})^{N-1}$, let $\mathbb{P}_N^{x,m_0,\chi}$ denote the law \mathbb{P}_N conditionally on $x_{\mathrm{tp}}(0) = x$, $m_{\mathrm{tp}}(0) = m_0$ and on the other $N-1$ particles at time zero having locations and masses given by χ . Then, for all such (m_0, x, χ) , and for any $T \geq 0$ and $A \subseteq \mathbb{R}^d$ open,*

$$(7.9) \quad \mathbb{P}_N^{x,m_0,\chi}(x_{\mathrm{tp}}(T) \in A) \leq \sup_{m \geq m_0} \frac{m_0}{m} \left(\frac{d(m_0)}{d(m)}\right)^{d/2} \cdot \nu_{x,2d(m_0)T}(A).$$

Before beginning this lemma's proof, we give an overview of the argument, which is an induction on $N \geq 1$. The case $N = 1$ may seem to be a triviality. Collision being impossible for a single particle, $x_{\mathrm{tp}}(T)$ is normally distributed with mean x and variance $2d(m_0)T$, while $m_{\mathrm{tp}}(T)$ equals m_0 almost surely; from which (7.9) follows. One might object however that conditioning on non-collision on the part of the tracer particle will bias the law of its trajectory; formally, we will treat the case $N = 1$ as an instance of the generic step of the induction.

In the case of several particles, a key role is played by the following uniform bound on normal densities: for all $x, y \in \mathbb{R}^d$ and $s \geq 0$, and for all $m, m' \in \mathbb{N}$ such that $m' > m$,

$$(7.10) \quad \frac{\mathrm{d}\nu_{x,2d(m')s}}{\mathrm{d}\nu_{x,2d(m)s}}(y) \leq \left(\frac{d(m)}{d(m')}\right)^{d/2}.$$

This bound is a consequence of $d : \mathbb{N} \rightarrow (0, \infty)$ being non-increasing.

We make some informal comments about the case $N = 2$ in order to illustrate the idea of the proof of Lemma 7.7. The tracer particle may make only one collision in this case. Consider the tracer particle dynamics until the first collision time for this particle at some time t , which we may assume to be less than T (for the other case is in effect the $N = 1$ case). Suppose that the tracer particle has mass m_0 before time t and collides with a mass m_1 particle at that time. Compare the subsequent dynamics to an altered one in which the two particles in the model do not interact. In the ordinary dynamics, the tracer particle survives the collision with probability $\frac{m_0}{m_0+m_1}$, and then pursues a Brownian trajectory of diffusion rate $2d(m_0+m_1)$; thus, the conditional probability given the tracer particle trajectory until first collision at time t that $x_{\mathrm{tp}}(T) \in A$ equals $\frac{m_0}{m_0+m_1} \nu_{x_{\mathrm{tp}}(t),2d(m_0+m_1)(T-t)}(A)$. On the other hand, in the altered dynamics, the tracer particle remains of mass m_0 at time t , and thus has conditional probability $\nu_{x_{\mathrm{tp}}(t),2d(m_0)(T-t)}(A)$ of achieving $x_{\mathrm{tp}}(T) \in A$. The uniform bound (7.10) on normal densities implies that the conditional probability of $x_{\mathrm{tp}}(T) \in A$ for the ordinary dynamics exceeds that for the altered dynamics by a factor of at most $\frac{m_0}{m_0+m_1} \left(\frac{d(m_0)}{d(m_0+m_1)}\right)^{d/2}$. Noting that the tracer particle in the altered dynamics is simply a Brownian particle of diffusion rate $2d(m_0)$ for all time, we may average over the tracer particle trajectory

until first collision, and, in doing so, we see that $\mathbb{P}_N(x_{\text{tp}}(T) \in A)$ is at most a $\frac{m_0}{m_0+m_1} \left(\frac{d(m_0)}{d(m_0+m_1)} \right)^{d/2}$ -multiple of the probability $\mu_{x,2d(m_0)T}(A)$ that the altered dynamics tracer particle reaches A at time T . Taking a supremum in m_1 heuristically explains (7.9) when $N = 2$. When $N > 2$, the tracer particle may collide several times, with particles of successive masses m_1, m_2, \dots, m_k , say. In essence, the same line of argument works, with a comparison factor of $\frac{\sum_{j=0}^{i-1} m_j}{\sum_{j=0}^i m_j} \left(\frac{d(\sum_{j=0}^{i-1} m_j)}{d(\sum_{j=0}^i m_j)} \right)^{d/2}$ being associated to the i -th collision. The product of these telescoping factors, $\frac{m_0}{\sum_{j=0}^k m_j} \left(\frac{d(m_0)}{d(\sum_{j=0}^k m_j)} \right)^{d/2}$, is then an upper bound on the ratio of the probabilities of $x_{\text{tp}}(T) \in A$ in the interacting model \mathbb{P}_N and in the model formed from \mathbb{P}_N by the suppression of all collisions on the part of the tracer particle. Heuristically this also explains the form (7.9) when $N \geq 2$.

7.2.3 Proof of Lemma 7.7

We turn to the rigorous argument establishing the general inductive step. Let $N \geq 1$ be given. Assume then that the statement of Lemma 7.7 is known for values of the inductive parameter strictly less than N . We will analyse $\mathbb{P}_N^{x,m_0,\chi}(x_{\text{tp}}(T) \in A)$ as an average of the conditional probability of $x_{\text{tp}}(T) \in A$ given the tracer particle trajectory until immediately before the first collision. To this end, for $t \geq 0$, we write \mathcal{F}_t for the σ -algebra generated by the \mathbb{P}_N -random variables $x_{\text{tp}} : [0, t] \rightarrow \mathbb{R}^d \cup \{c\}$ and $m_{\text{tp}} : [0, t] \rightarrow \mathbb{N} \cup \{c\}$, so that the information available in \mathcal{F}_t is the data given by monitoring the tracer particle during $[0, t]$. We also write \mathcal{F}_t^- for the sigma-algebra generated by $\{\mathcal{F}_s : 0 \leq s < t\}$, representing the information concerning the tracer particle's history which is available immediately before time t . Let $\sigma_1 \geq 0$ denote the time of the tracer particle's first collision. We also set $\sigma_1^T = \sigma_1 \wedge T$.

Note then that we may express

$$(7.11) \quad \mathbb{P}_N^{x,m_0,\chi}(x_{\text{tp}}(T) \in A) = \mathbb{E} \mathbb{P}_N^{x,m_0,\chi}(x_{\text{tp}}(T) \in A \mid \mathcal{F}_{\sigma_1^T}^-).$$

Note that in the right-hand side, we are writing the probability that $x_{\text{tp}}(T) \in A$ as an average over tracer particle histories up to, but not including, the first collision time. Given any instance of data in $\mathcal{F}_{\sigma_1^T}^-$ for which $\sigma_1 < T$, it is known that the tracer particle is about to experience a collision at time σ_1 , even though the mass of the second particle participating in the collision and the collision's outcome – the survival or perishing of the tracer particle – remain random events. Consider the instantaneous future of the tracer particle trajectory under the law $\mathbb{P}_N^{x,m_0,\chi}(\cdot \mid \mathcal{F}_{\sigma_1^T}^-)$. We may denote by m_1 the mass of the particle with which the tracer particle collides at time σ_1 ; note that under the conditional law, m_1 is a random variable. Let \mathbf{S} denote the event that the tracer particle survives this collision, so that the conditional probability of \mathbf{S} under $\mathbb{P}_N^{x,m_0,\chi}(\cdot \mid \mathcal{F}_{\sigma_1^T}^-, m_1)$ equals $\frac{m_0}{m_0+m_1}$. If \mathbf{S}^c occurs, then x_{tp} immediately arrives in the cemetery state c , so that, if $\sigma_1 < T$, there is no possibility that $x_{\text{tp}}(T) \in A$ in this event. On the other hand, should \mathbf{S} occur, $m_{\text{tp}}(\sigma_1) = m_0 + m_1$, so that the conditional probability that $x_{\text{tp}}(T) \in A$ occurs is given as follows: $\mathbb{P}_N^{x,m_0,\chi}(\cdot \mid \sigma_1 < T)$ -almost surely,

$$(7.12) \quad \mathbb{P}_N^{x,m_0,\chi}(x_{\text{tp}}(T) \in A \mid \mathcal{F}_{\sigma_1^T}^-, m_1, \mathbf{S}) = \int \mathbb{P}_n^{x_{\text{tp}}(\sigma_1), m_0+m_1, \phi}(x_{\text{tp}}(T - \sigma_1) \in A) d\mu_{(n,\phi)}.$$

$$\begin{aligned}
(7.13) \quad \mathbb{P}_N^{x,m_0,\chi}(x_{\text{tp}}(T) \in A) &= \mathbb{E}_N^{x,m_0,\chi} \left[\mathbb{E} \left[\mathbf{1}_{x_{\text{tp}}(T) \in A} \mid \mathcal{F}_{\sigma_1^T}^- \right] \right] \\
&= \mathbb{E}_N^{x,m_0,\chi} \left[\mathbb{E} \left[(\mathbf{1}_{\sigma_1 < T} + \mathbf{1}_{\sigma_1 \geq T}) \mathbf{1}_{x_{\text{tp}}(T) \in A} \mid \mathcal{F}_{\sigma_1^T}^- \right] \right].
\end{aligned}$$

Note that

$$\begin{aligned}
(7.14) \quad \mathbb{E} \left[\mathbf{1}_{\sigma_1 < T} \mathbf{1}_{x_{\text{tp}}(T) \in A} \mid \mathcal{F}_{\sigma_1^T}^- \right] &= \mathbf{1}_{\sigma_1 < T} \cdot \mathbb{E} \left[\mathbf{1}_{x_{\text{tp}}(T) \in A} \mathbf{1}_S \mid \mathcal{F}_{\sigma_1^T}^- \right] \\
&= \mathbf{1}_{\sigma_1 < T} \int \frac{m_0}{m_0+m_1} \mathbb{P}_n^{x_{\text{tp}}(\sigma_1), m_0+m_1, \phi} (x_{\text{tp}}(T - \sigma_1) \in A) d\mu_{(n, m_1, \phi)},
\end{aligned}$$

where $\mu_{(n, m_1, \phi)}$ is the $\mathcal{F}_{\sigma_1^T}^-$ -measurable random measure that specifies the following data:

- n , the number of surviving particles immediately after the collision at time σ_1 ;
- m_1 , the mass of the particle with which x_{tp} collides at time σ_1 ;
- and $\phi \in (\mathbb{R}^d \times \mathbb{N})^{n-1}$, the vector of locations and masses of the particles other than the tracer particle at this time;

and $\mu_{(n, m_1, \phi)}$ denotes the $\mathcal{F}_{\sigma_1^T}^-$ -measurable random measure that specifies the conditional distribution of (n, m_1, ϕ) under the law $\mathbb{P}_N^{x, m_0, \chi}(\cdot \mid \mathcal{F}_{\sigma_1^T}^-)$.

Since a collision occurs at time σ_1 , n is necessarily at most $N - 1$, so that the inductive hypothesis may be applied to bound above the term $\mathbb{P}_n^{x_{\text{tp}}(\sigma_1), m_0+m_1, \phi}(x_{\text{tp}}(T - \sigma_1) \in A)$ appearing in the integrand above. We find that, for each $m \geq 1$,

$$\begin{aligned}
&\sup_{1 \leq k \leq N-1, \phi \in (\mathbb{R}^d \times \mathbb{N})^k} \mathbb{P}_k^{x_{\text{tp}}(\sigma_1), m_0+m, \phi} (x_{\text{tp}}(T - \sigma_1) \in A) \\
&\leq \sup_{m' \geq m_0+m} \frac{m_0+m}{m'} \left(\frac{d(m_0+m)}{d(m')} \right)^{d/2} \cdot \nu_{x_{\text{tp}}(\sigma_1), 2d(m_0+m_1)(T-\sigma_1)}(A) \\
&\leq \sup_{m' \geq m_0+m} \frac{m_0}{m'} \left(\frac{d(m_0+m)}{d(m')} \right)^{d/2} \cdot \nu_{x_{\text{tp}}(\sigma_1), 2d(m_0)(T-\sigma_1)}(A),
\end{aligned}$$

where the latter inequality invoked the uniform bound (7.10) in the guise $\frac{d\nu_{x, 2d(m')}(y)}{d\nu_{x, 2d(m_0)}} \leq \left(\frac{d(m_0+m)}{d(m')} \right)^{d/2}$ for all $x, y \in \mathbb{R}^d$.

Returning to (7.14), we obtain

$$\mathbb{E} \left[\mathbf{1}_{\sigma_1 < T} \mathbf{1}_{x_{\text{tp}}(T) \in A} \mid \mathcal{F}_{\sigma_1^T}^- \right] \leq \mathbf{1}_{\sigma_1 < T} \cdot \sup_{m' \geq m_0+1} \frac{m_0}{m'} \left(\frac{d(m_0)}{d(m')} \right)^{d/2} \cdot \nu_{x_{\text{tp}}(\sigma_1), 2d(m_0)(T-\sigma_1)}(A).$$

We also have

$$\mathbb{E} \left[\mathbf{1}_{\sigma_1 \geq T} \mathbf{1}_{x_{\text{tp}}(T) \in A} \mid \mathcal{F}_{\sigma_1^T}^- \right] = \mathbf{1}_{\sigma_1 \geq T} \mathbf{1}_{x_{\text{tp}}(T) \in A}.$$

Applying the two preceding equations to (7.13), we find that

$$(7.15) \quad \begin{aligned} & \mathbb{P}_N^{x, m_0, \chi}(x_{\text{tp}}(T) \in A) \\ & \leq \mathbb{E}_N^{x, m_0, \chi} \left[\mathbf{1}_{\sigma_1 < T} \cdot \sup_{m' \geq m_0 + 1} \frac{m_0}{m'} \left(\frac{d(m_0)}{d(m')} \right)^{d/2} \cdot \nu_{x_{\text{tp}}(\sigma_1), 2d(m_0)(T-\sigma_1)}(A) + \mathbf{1}_{\sigma_1 \geq T} \mathbf{1}_{x_{\text{tp}}(T) \in A} \right]. \end{aligned}$$

To bound above the right-hand side, note first that

$$\mathbb{E}_N^{x, m_0, \chi} \left[\mathbf{1}_{\sigma_1 < T} \cdot \nu_{x_{\text{tp}}(\sigma_1), 2d(m_0)(T-\sigma_1)}(A) \right] \leq \nu_{x, 2d(m_0)T}(A).$$

To see this, recall that, when $\sigma_1 < T$, $\nu_{x_{\text{tp}}(\sigma_1), 2d(m_0)(T-\sigma_1)}(A)$ is the probability that an independent Brownian motion of diffusion rate $2d(m_0)$ beginning from $x_{\text{tp}}(\sigma_1)$ at time σ_1 visits A at time T ; in the above left-hand side, the mean is taken over trajectories of a Brownian motion $x_{\text{tp}} : [0, \sigma_1] \rightarrow \mathbb{R}^d$ of diffusion rate $2d(m_0)$ with initial condition $x_{\text{tp}}(0) = x$. That is, the value of this left-hand side is given by the probability that a Brownian motion of diffusion rate $2d(m_0)$ which at time zero is at x visits A at time T (and that an auxiliary stopping time, σ_1 , occurs before T).

Similarly, we have that

$$\mathbb{E}_N^{x, m_0, \chi} \left[\mathbf{1}_{\sigma_1 \geq T} \mathbf{1}_{x_{\text{tp}}(T) \in A} \right] \leq \nu_{x, 2d(m_0)T}(A).$$

Indeed, the subprobability measure $\mathbb{P}_N^{x, m_0, \chi}(\cdot \cap \{\sigma_1 \geq T\})$ is stochastically dominated by Brownian motion begun at x of diffusion rate $2d(m_0)$, since the trajectory of x_{tp} suffers no collision on $[0, T]$ under this defective law.

The last inferences when allied with (7.15) yield

$$\mathbb{P}_N^{x, m_0, \chi}(x_{\text{tp}}(T) \in A) \leq \sup_{m' \geq m_0} \frac{m_0}{m'} \left(\frac{d(m_0)}{d(m')} \right)^{d/2} \cdot \nu_{x, 2d(m_0)T}(A);$$

this is (7.9) for index N , so that the inductive proof of Lemma 7.7 is complete. \square

7.2.4 Proof of Lemma 7.4.

Using Lemma 7.6, and the fact that the tracer particle has mass $m \in \mathbb{N}$ at time zero with probability $Z^{-1} \|h_m\|_{L^1(\mathbb{R}^d)}$, Lemma 7.7 yields the deduction that

$$\mathbb{P}_N(\chi_T \cap A \neq \emptyset) \leq NZ^{-1} \sum_{m_0 \geq 1} \|h_{m_0}\|_{L^1(\mathbb{R}^d)} \cdot m_0 d(m_0)^{d/2} \sup_{m \geq m_0} m^{-1} d(m)^{-d/2} \cdot \mathbb{E}(\nu_{X_{m_0}, 2d(m_0)T}(A)),$$

where the mean in the final term is over X_{m_0} , a random variable having the law of $x_{\text{tp}}(0)$ given that $m_{\text{tp}}(0) = m_0$, so that X_{m_0} has density $\frac{h_{m_0}(\cdot)}{\|h_{m_0}\|_{L^1(\mathbb{R}^d)}}$ on \mathbb{R}^d . The heat equation decreases the L^∞ -norm, so that $\mathbb{E}(\nu_{X_{m_0}, 2d(m_0)T}(A)) \leq \frac{1}{\|h_{m_0}\|_{L^1(\mathbb{R}^d)}} \ell_{m_0} \cdot \mu(A)$, where recall that $\ell_{m_0} = \|h_{m_0}\|_{L^\infty(\mathbb{R}^d)}$. In this way, we obtain Lemma 7.4. \square

7.2.5 Monitoring several tracer particles at once

The proof of Proposition 7.3 is derived similarly as was Lemma 7.4. In asking about the distribution at time T of collections of particles of size k , rather than about single particles, we make use of a natural generalization of the concept of the tracer particle.

Definition 7.8 Fix $k \in \mathbb{N}$. Under the law \mathbb{P}_N , we define an independent random variable $(U_1, \dots, U_k) \in \{1, \dots, N\}^k$ having the uniform distribution on k -sized subsets of $\{1, \dots, N\}$ (so that there are no coincidences among the U_i). For $1 \leq i \leq k$, the i -th tracer particle is the particle whose index is U_i . The location and mass of the i -th tracer particle will be denoted by $x_{\text{tp}}^i : [0, \infty) \rightarrow \mathbb{R}^d \cup \{c\}$ and $m_{\text{tp}}^i : [0, \infty) \rightarrow \mathbb{N} \cup \{c\}$.

Here is the analogue of Lemma 7.6.

Lemma 7.9 Let $k \in \mathbb{N}$ and let $\{A_i : 1 \leq i \leq k\}$ be a collection of open balls in \mathbb{R}^d . Then for each $T \geq 0$,

$$\mathbb{P}\left(\mathcal{X}_T \cap A_i \neq \emptyset \text{ for each } 1 \leq i \leq k\right) = k! \cdot N^k \cdot \mathbb{P}\left(x_{\text{tp}}^i(T) \in A_i \text{ for } 1 \leq i \leq k\right).$$

As Lemma 7.4 was, Proposition 7.3 will follow from a stronger assertion which will be established by induction on N .

Lemma 7.10 Given $\bar{x} = (x_1, \dots, x_k) \in \mathbb{R}^d$, $\bar{m} = (m_1, \dots, m_k) \in \mathbb{N}^k$ and $\chi \in (\mathbb{R}^d \times \mathbb{N})^{N-k}$, we denote by $\mathbb{P}_N^{\bar{x}, \bar{m}, \chi}$ the law \mathbb{P}_N , conditionally on $x_{\text{tp}}^i(0) = x_i$, $m_{\text{tp}}^i(0) = m_i$ for $1 \leq i \leq k$, and on the other $N - k$ particles at time zero having locations and masses given by χ . Then, for all such data, and for any $T \geq 0$ and open $A_i \subseteq \mathbb{R}^d$, $1 \leq i \leq k$,

$$(7.16) \quad \mathbb{P}_N^{\bar{x}, \bar{m}, \chi}\left(\bigcap_{1 \leq i \leq k} \{x_{\text{tp}}^i(T) \in A_i\}\right) \leq \prod_{i=1}^k \sup_{m \geq m_i} \frac{m_i}{m} \left(\frac{d(m_i)}{d(m)}\right)^{d/2} \nu_{x_i, 2d(m_i)T}(A_i).$$

Proof. The proof of the lemma is in essence the same as that of Lemma 7.7. It works by induction on N , beginning with $N = k$. Under the law \mathbb{P}_N , let $\sigma \geq 0$ denote the time of the earliest collision experienced by one of the k tracer particles. Set $\sigma' = \min\{\sigma, T\}$. If $\sigma' < T$ then either two tracer particles collide at time σ' , rendering the event that each such particle reaches its target set A_i at time T impossible, or one of the tracer particles collides with one of the $N - k$ non-tracer particles. The latter case is the non-trivial one, and it may be analysed exactly as in the proof of Lemma 7.7: if the tracer particle in question perishes on collision, it is assigned to the state c and there is nothing to prove; otherwise, it survives, assumes some added mass, among at most $N - 1$ other particles. The inductive hypothesis is applied to the new post-collision scenario. \square

Proof of Proposition 7.3. Applying Lemma 7.10 to the random initial condition $(x_{\text{tp}}, m_{\text{tp}})$ arising from the law \mathbb{P}_N , the worst case of the resulting bound occurs when $m_{\text{tp}}(0) = 1$. Using this bound, and that the density of $\nu_{x, 2d(1)T}$ is uniformly at most $(2d(1)T)^{-d/2}$, we obtain the result. \square

7.3 Uniform integrability

We now use Proposition 7.3 to prove Propositions 4.3 and 4.4.

Proof of Propositions 4.3 and 4.4. We seek a contradiction by assuming that Proposition 4.4 is false. Thus, for some $m \in \mathbb{N}$ and for all $K > 0$, there exists $\epsilon > 0$, $T > 0$ and $A \subseteq \mathbb{R}^d$ open such that, for some subsequence $\{N_i : i \in \mathbb{N}\} \subseteq \mathbb{N}$,

$$\mathbb{P}_{N_i}(|\chi_T^m \cap A| \geq KN_i \mu(A)) \geq \epsilon.$$

This assertion is also true for i sufficiently high if we demand that $A = A_{n_i, c}$ for some open set $A_{n_i, c}$ such that $N_i \mu(A_{n_i, c}) = c$, where $c > 0$ is fixed.

However, Proposition 7.3 implies that, for each $k \in \mathbb{N}$,

$$\mathbb{P}_{N_i}(|\chi_T^m \cap A| \geq k) \leq (CN\mu(A))^k,$$

where $C = Z^{-1} \sum_{m_0=1}^m \ell_{m_0} m_0 d(m_0)^{d/2} \sup_{m \geq m' \geq m_0} (m')^{-1} d(m')^{-d/2}$. For our choice $A = A_{n_i, c}$, we thus have that, for i sufficiently high,

$$\mathbb{P}_{N_i}(|\chi_T^m \cap A_{n_i, c}| \geq Kc) \leq (Cc)^{Kc},$$

Thus, $\epsilon \leq (Cc)^{Kc}$ for all $K > 0$. We arrive at a contradiction by choosing $c \in (0, C^{-1})$. \square

8 Review and summary

In this section, we list those instances where our derivation of Theorem 1.1 was incomplete, explaining where the proof is furnished in [10]; discuss an imprecision in the proof of [10] which we have sought to clarify by our presentation in this survey; and make a limited comparison between our present method of proof of key estimates by means of particle concentration bounds with the approach adopted in [10].

8.1 The list of shortcuts in the survey's proofs

In our proof of Theorem 1.1, several steps are omitted. Beyond the absence of a proof of the classical Feynman-Kac formula, in the guise that $v(x, t)$ in Lemma 1.6 satisfies the PDE (1.14), and the full justification for the uniqueness assertion in Lemma 1.6 (for which the reader may see the analytic treatment given in [10, Section 6]), the missing steps are:

- In the reduction of Proposition 4.1 to Proposition 5.1 undertaken in Section 5.1, the test function $J_n(x, t)$ was chosen to be equal to be identically one. The more general case requires only a few further lines of argument: see [10, Section 3.5].
- Proposition 4.2, showing smallness of the martingale $M(T)$ in (4.1), is not proved. See [10, Section 5].

- In the proof of Proposition 5.1 in Section 6, the form of the test functions $J(x, n, t)$ and $\bar{J}(x, m, t)$ was simplified so that they have no space-time dependence. When this simplification is omitted, the action of the free motion operator in (6.4) generates some extra terms, where one or both of the derivatives in the Laplacian fall on the test functions. The resulting terms tend to be smoother than the existing terms, and the methods of treating them are the same as for their rougher counterparts. See the start of [10, Section 3] for the full scale version of (6.4) and [10, Section 3.3] for bounds on the terms appearing in that version.
- We have offered no proof of Proposition 6.1(2), (3) and (4). Similarly to the previous point, this omission is a simplification of the presentation of the proof of the Stosszahlansatz. The key tools – uniform killing probability bounds and particle concentration results – apply equally to prove these statements as we saw that they did to prove Proposition 6.1(1). See Sections 3.3 and 3.4 of [10] for the relevant bounds, valid under the original assumptions.

8.2 A momentary spotlight on an obscurity

In the opening paragraph of [10, Section 4], the step counterpart to that of the present Section 4.2 is discussed: the low ϵ limit is taken of the approximate identity that is (4.1) in this survey. However, in the replacement of the collision term by its counterpart expressed using microscopic candidate densities, which happens by means of the Stosszahlansatz, [10] neglects to clarify that this replacement must be made simultaneously over the infinitely many mass pairs $\{(n, m) : m \in \mathbb{N}\}$, rather than merely being made for one such mass pair. What permits this simultaneous replacement is that the bound satisfied by the error $\text{Err}_{n,m}$ in Proposition 4.1 contains a sum over $m \in \mathbb{N}$. As we have seen, the reason why we are able to prove the Stosszahlansatz with such an error bound relies on the uniform control on the killing probabilities u_W that we saw in Subsection 6.2.2, specifically that Lemma 6.3 is valid as the kernel W varies over choices having given compact support. Of course, in the present survey, we have not presented a proof of all of the required estimates for Theorem 1.1. A complete proof of the result, under the original assumptions made in [10], is formed by rendering [10, Lemma 3.2] uniform over $(n, m) \in \mathbb{N}^2$; the changes needed to do this are contained in the proof of the present Lemma 6.3.

8.3 Comparison with later kinetic limit derivations of the Smoluchowski PDE

Several ramifications of the statement and technique of proof of [10] have been explored by Fraydoun Rezakhanlou, sometimes in collaboration with the author and others. We end by summarising the results so obtained and comparing the approaches to proof in these further articles, both with the original one in [10] and with that expounded here.

8.3.1 The particle concentration bound: its robustness and limitations

The principal technical novelty presented in this survey is the use of the particle concentration bound Proposition 6.2 to yield the key error bounds Proposition 6.1: the technique used in [10] was quite different. Our present technique requires stronger hypotheses, but when it may be applied, it

yields strong conclusions about diverse aspects of particle dynamics. We now explain this summary by drawing a contrast with the method used in [10].

First, to expand, Proposition 6.2 offers strong conclusions about the lack of build-up in particle concentration at positive times in the models \mathbb{P}_N . However, it has content only under the fairly restrictive hypothesis that $d(m)$ decays no faster than $m^{-2/d}$, and, regarded as a tool to prove Proposition 6.1, its use must be accompanied by the assumption that $\alpha(\cdot, \cdot)$ is bounded above uniformly. Here we make a comment about the one simple aspect of the quite different approach that was adopted in [10] to prove the key estimates that correspond to the present Proposition 6.1. We now state a result giving an upper bound on the duration $[0, T]$ total mean collision rate in the models \mathbb{P}_N . It is [10, Lemma 3.1], which we call the “bound on the collision”.

Lemma 8.1 *For any $N \in \mathbb{N}$, and for $T > 0$,*

$$\epsilon^{d-2} \mathbb{E}_N \int_0^T ds \sum_{i,j \in I_q} \alpha(m_i, m_j) V_\epsilon(x_i - x_j) \leq Z$$

Proof. Let $X(t)$ denote the number of surviving particles in \mathbb{P}_N at time t . Consider the variant of (4.1) in which the term J_n is replaced by $\mathbf{1}$ and the result summed over $n \in \mathbb{N}$, so that, for example, the first sum is a total particle count at time T . Taking expectations, we find that, for any $T > 0$,

$$(8.1) \quad \mathbb{E}_N X(T) = \mathbb{E}_N X(0) + \int_0^T \mathbb{E}_N \mathfrak{A}_F X(t) dt + \int_0^T \mathbb{E}_N \mathfrak{A}_C X(t) dt.$$

Particle count is conserved by free motion, so that $\mathbb{E}_N \mathfrak{A}_F X(t) = 0$. On the other hand, the integrated mean collision rate $\int_0^T \mathbb{E}_N \mathfrak{A}_C X(t) dt$ is equal to $\mathbb{E}_N \int_0^T ds \sum_{i,j \in I_q} \alpha(m_i, m_j) V_\epsilon(x_i - x_j)$. By (8.1), this quantity equals $\mathbb{E}_N X(0) - \mathbb{E}_N X(T)$ which is at most $\mathbb{E}_N X(0) = N$. Since $N = Z\epsilon^{2-d}$ by (1.6), we obtain the result. \square

Of course, the proof is almost a triviality. However, it already highlights differences with, and the limitations of, the particle concentration bound Proposition 6.2. Let us try to emulate this unprepossessing lemma’s conclusion by using Proposition 6.2. By merely applying this result without using further tricks, the best we can do is the following.

Claim 8.2 *For any $N \in \mathbb{N}$ and $T > 0$, we have that*

$$\begin{aligned} & \epsilon^{d-2} \mathbb{E}_N \int_0^T ds \sum_{i,j \in I_q} \alpha(m_i, m_j) V_\epsilon(x_i - x_j) \\ & \leq TZ^2 \cdot \|V\|_\infty \cdot \sup_{n,m} \alpha(n, m) \cdot \left(\sup_{m \geq 1} m^{-1} d(m)^{-d/2} \right)^2 \cdot \sup_{n \in \mathbb{N}} \|h_n\|_\infty. \end{aligned}$$

Proof. The relation $N = Z\epsilon^{2-d}$ and Proposition 6.2 applied with $k = 2$ show that, for any time

$t \in (0, \infty)$,

$$\begin{aligned} \sum_{i,j \in [1,N]} \mathbb{P}_N(\|x_j(t) - x_i(t)\| \leq \epsilon) &\leq N^2 \mathbb{P}_N(\|x_2(t) - x_1(t)\| \leq \epsilon) \\ &\leq Z^2 \epsilon^{2(2-d)} \cdot \epsilon^d \cdot \left(\sup_{m \geq 1} m^{-1} d(m)^{-d/2} \right)^2 \cdot \sup_{n \in \mathbb{N}} \|h_n\|_\infty. \end{aligned}$$

Note also that

$$\sum_{i,j \in [1,N]} \mathbb{E}_N \left[\alpha(m_i(t), m_j(t)) V\left(\frac{x_i(t) - x_j(t)}{\epsilon}\right) \right] \leq \sup_{n,m} \alpha(n, m) \cdot \|V\|_\infty \cdot \sum_{i,j \in [1,N]} \mathbb{P}_N(\|x_j - x_i\| \leq \epsilon).$$

If we multiply the above left-hand side by ϵ^{d-2} , and then further by ϵ^{-2} – which we do because $V_\epsilon(\cdot) = \epsilon^{-2} V(\cdot/\epsilon)$ – and integrate over $[0, T]$, then we obtain the left-hand side in the statement of the claim. Thus, the claim follows from the above two inequalities. \square

Claim 8.2 falls short of Lemma 8.1 in a number of ways, including the appearance of the factor $\sup_{n,m} \alpha(n, m)$ on its right-hand side. The reason that such a factor appears is because particle pair presence at distance of order ϵ at positive times is penalized due to the microscopic repulsion phenomenon which has been central to this survey, and which in particular we discussed in Section 3. Crudely, if $\alpha(n, m)$ is high, then the density for such presence is not of order ϵ^{d-2} but of order $\alpha(n, m)^{-1} \epsilon^{d-2}$. Proposition 6.2 is not built to acknowledge this microscopic repulsion effect and the unwanted α goes uncanceled when the proposition is applied.

Lemma 8.1 experiences no such limitation. But of course its description of positive time particle distribution is limited to a very specific aspect of the overall dynamics; in comparison, Proposition 6.2 is a robust tool, that will say something meaningful about any such aspect, when its hypotheses are satisfied.

Lemmas 3.2 and 3.3 of [10] form the counterpart to the present Proposition 6.2 in the sense that they are tools used to prove the present Proposition 6.1. The two lemmas treat several aspects of particle dynamics other than the bound on the collision given in Lemma 8.1. The proofs of these results generalize that of Lemma 8.1 in the sense that the mean value of variants of (4.1) are considered and their terms bounded. In the case of [10, Lemma 3.3], some of these terms involve sums over triples of particle indices, with one of the indices having no restriction on the mass parameter. In such cases, (1.10) in the original assumptions is invoked during a proof by induction in order to find suitable bounds.

8.3.2 The planar case: the route to the PDE

In Section 2.5, we mentioned that, in [9], the kinetic limit derivation counterpart to that of [10] was undertaken for dimension $d = 2$; we also reviewed the main changes to Theorem 1.1’s statement in this case. The technique of proof is the same as in the original work, with the principal technical change concerning the particle distribution result [10, Lemma 3.2], where the proof may not be directly utilized because the non-negativity of the solution H of Poisson’s equation $-\Delta H = J$ (for a given non-negative J) is enjoyed in dimension $d \geq 3$ but not $d = 2$. We refer the reader to [9]

for further discussion of this technicality. However, we note that the robustness of the particle concentration bound Proposition 6.2 has the virtue of permitting the extension of the proof of Theorem 1.1 developed in this survey to the case $d = 2$, under the survey assumptions, without any comparable technical difficulty arising.

The kinetic limit derivations of [10] and [9] were extended to cases of variable radial dependence for particles in [23]; see the present Subsection 2.4.2. The technique of proof, including the treatment of particle distribution bounds, is similar to that of the earlier works.

8.3.3 The case of continuous mass and a different approach to particle concentration

As we mentioned after the equations in Section 1.2, the Smoluchowski coagulation-diffusion PDE has a continuous counterpart, where the mass parameter is now a non-negative real. The kinetic limit derivation of the PDE is revisited in [31] in the case that $d \geq 3$ for the PDE with continuous mass parameter. The principal innovation of the article is [31, Theorem 3.1], a tool for proving particle concentration which is novel in comparison with that of [10] and [9]. This technique has distinct similarities with our tracer particle proof of Proposition 6.2: the comparison of diffusion rate dependent terms in [31, (3.6)] is a rough counterpart to (7.10).

8.3.4 The planar case with fragmentation: equilibrium fluctuations

The Smoluchowski PDE may be modified to include interaction terms corresponding to pairwise particle fragmentation. In the models \mathbb{P}_N , we may model this fragmentation effect by declaring that a particle of given mass is subject to fragment at the ring times of a Poisson clock that ticks at a mass-dependent rate. On fragmenting, the particle splits in two. The detailed rule for this splitting may be chosen to be a “dual” of the rule specified for coagulation under the heading “the precise mechanism for collision” in Section 1.3: the fragmenting particle retains its location, and some random proportion of its precollisional mass, while a new particle, bearing the residue of that mass, appears in a randomly selected microscopic vicinity of the fragmenting particle’s location.

Of course, one may attempt to carry out a kinetic limit derivation of the Smoluchowski coagulation-fragmentation-diffusion PDE from microscopic models \mathbb{P}_N that have been altered in this manner; no such derivation has yet been carried out. In [22], Ranjbar and Rezakhanlou studied a different aspect of particle dynamics in the case that dimension d equals two. The assertion that macroscopic particle densities adhere to a solution of the Smoluchowski PDE is in a sense a weak law of large numbers. What of the analogue of the central limit theorem, a result describing the typical fluctuation of particle density statistics in high indexed \mathbb{P}_N from the density profile offered by the PDE solution? In [22], the authors define empirical fluctuation fields under \mathbb{P}_N , modelling the discrepancy as a function of space-time of the empirical density of particles from the prediction made by the PDE solution, normalized by a square root of total particle number, in the style of the central limit theorem. In [22, Conjecture 2.1], it is conjectured that, in a high N limit, the fluctuation field converges to a random limit that solves an Ornstein-Uhlenbeck equation under which the density profile diffuses freely and is subject to coagulative and fragmenting forces specified by a linearization of those present in the PDE, as well as to a space-time dependent white noise stimulus determined by the PDE solution.

To endeavour to prove this conjecture may seem a fanciful task, given that the kinetic limit derivation of an analogue of Theorem 1.1 is yet to be attempted. However, the authors advance a case for the conjecture by rigorously analysing the system at equilibrium. The modified microscopic models \mathbb{P}_N have mechanisms for both coagulation of pairs of particles, and fragmentation of particles into pairs; and these mechanisms have been chosen so that the film of coagulation event when played in reverse shows the fragmentation event. Thus, the equilibrium measures of the laws \mathbb{P}_N may be explicitly identified: under them, the distribution of particles of any given mass is simply a Poisson process (of some constant intensity determined by the mass), with the particle cloud of particles of differing mass being independent. In [22, Theorem 3.1], the conjecture of convergence to the Ornstein-Uhlenbeck process mentioned above is proved for the system at these equilibria. The proof requires an understanding of the relation between microscopic and macroscopic interaction propensities which extends (2.4) to treat fragmentation but also involves an unexpected interaction with the free motion dynamics. Note, however, that one subtlety which we have emphasised is absent: the microscopic repulsion effect, elucidated in Section 3 and reflected in the formula (1.8), is no longer present, because the microscopic system at equilibrium is a union of independent Poisson processes.

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