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Suppression of Particle Dispersion by Sweeping Effects in Synthetic Turbulence

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Synthetic models of Eulerian turbulence like so called “Kinematic Simulations” (KS) are often used as computational shortcuts for studying Lagrangian properties of turbulence. These models have been criticized by Thomson \& Devenish (2005), who argued on physical grounds that sweeping decorrelation effects suppress pair dispersion in such models. We derive analytical results for Eulerian turbulence modeled by Gaussian random fields, in particular for the case with zero mean velocity. Our starting point is an exact integrodifferential equation for the particle pair separation distribution obtained from the Gaussian integration-by-parts identity. When memory times of particle locations are short, a Markovian approximation leads to a Richardson-type diffusion model. We obtain a time-dependent pair diffusivity tensor of the form $K_{ij}(r,t) = S_{ij}(r)\tau(r,t)$ where $S_{ij}(r)$ is the structure-function tensor and $\tau(r,t)$ is an effective correlation time of velocity increments. Crucially, this is found to be the minimum value of three times: the intrinsic turnover time $\tau_{edd}(r)$ at separation $r$, the overall evolution time $t$, and the sweeping time $t/v_0$ with $v_0$ the rms velocity. We study the diffusion model numerically by a Monte Carlo method. With inertial-ranges like the largest achieved in most current KS (about six decades long), our model is found to reproduce the $t^{3/2}$ power-law for pair dispersion predicted by Thomson \& Devenish and observed in the KS. However, for much longer ranges, our model exhibits three distinct pair-dispersion laws in the inertial-range: a Batchelor $t^2$-regime, followed by a Kraichnan-model-like $t^4$ diffusive regime, and then a $t^6$ regime. Finally, outside the inertial-range, there is another $t^2$ regime with particles undergoing independent Taylor diffusion. These scalings are exactly the same as those predicted by Thomson \& Devenish for KS with large mean velocities, which we argue hold also for KS with zero mean velocity. Our results support the basic conclusion of Thomson \& Devenish (2005) that sweeping effects make Lagrangian properties of KS fundamentally different from hydrodynamic turbulence for very extended inertial-ranges.

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I. INTRODUCTION

How particle pairs separate in a turbulent flow has been a central subject of turbulence research since the classical work of Richardson [1]. Unfortunately, the phenomenon has proved quite difficult to investigate by numerical solution of the fluid equations and by controlled laboratory experiments, especially because of the very large Reynolds numbers required. Many studies have therefore employed “synthetic turbulence” or ensembles of random velocity fields with some of the scaling properties of real turbulent velocities but which can be efficiently sampled even for very long scaling ranges. For example, papers [2–8] have followed this approach and have reported substantial agreement of their numerical simulations with the predictions of Richardson, especially the famous “$t^3$-law” for the growth in time of mean square particle separation distances.

The asymptotic validity of a $t^3$ scaling law for these synthetic turbulence models at very high Reynolds numbers has been called into question, however. Papers [9, 10] for example pointed out that the use of synthetic turbulence to model Eulerian velocity statistics implies sweeping effects of large-scale eddies on particle motions that diverge with the Reynolds number. These authors suggested to employ synthetic ensembles, such as multiplicative cascade models [9] or Gaussian random fields [10], to model instead the turbulent statistics of (quasi)Lagrangian velocities [47]. In a simple one-dimensional Gaussian model of Eulerian velocities, Chaves et al. [10] found analytically that large-scale sweeping effects “localized” particle pairs and prevented them from separating. Subsequently, in a very interesting paper [11], Thomson \& Devenish have proposed an intuitive picture how sweeping affects particle dispersion in synthetic models of Eulerian turbulence. The key point is that large-scale eddies in real turbulence advect both particles and smaller scale eddies, while large-scale eddies in synthetic turbulence advect only particle pairs and not smaller eddies. This fact implies that particle pairs at separations $r$ in synthetic turbulence should experience rapidly changing relative velocities, as they are swept into new, statistically independent eddies. This occurs on a “sweeping” time-scale $\tau_{sw}(r) \sim r/v_0$, where $v_0$ is the rms velocity set by the largest eddies in the synthetic ensemble. Thomson \& Devenish assume a diffusion process of pair separations with an eddy-diffusivity $K(r) \sim \delta u^2(r)\tau_{sw}(r)$ and $\delta u^2(r)$ the mean-square relative velocity at separation $r$. In an ensemble with Kolmogorov scaling $\delta u^2(r) \sim (\varepsilon r)^{2/3}$, this yields

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\[ \frac{dr^2}{dt} \sim K(r) \sim \varepsilon^{2/3} r^{5/3} / v_0 \] and the solution

\[ \langle r^2(t) \rangle \sim \frac{\varepsilon t^6}{v_0^6} \]  

(1)

Note that this implies considerably slower growth than Richardson’s \( t^3 \)-law [48]. Thomson & Devenish argued for the above prediction in the case of a large mean sweeping, with \( v_0 \) replaced by the mean speed \( \overline{v} \). In the case of a zero-mean velocity ensemble, they argued instead for a \( t^{9/2} \)-growth law, intermediate between \( t^3 \) and \( t^6 \) (see our section III below). These predictions were supported in [11] by the numerical technique of “Kinematic Simulations” (KS) [3–5, 7, 12]. The previous contrary numerical results were explained on various grounds, e.g. the use of an adaptive time-stepping scheme in [3] which did not resolve the small sweeping time \( \tau_{sw}(r) \) and its effect on particle dispersion.

The issues raised by the paper of Thomson & Devenish have still not been fully resolved. The numerical simulations in [11] used another form of adaptive time-stepping, which was suggested in [13] to be responsible for the observation of a \( t^{9/2} \) growth. Thomson & Devenish then repeated their simulations with a fixed small time-step and reported the same \( t^{9/2} \) law [14]. The most recent simulations of Nicolleau & Nowakowski [15] for their longest scaling ranges show some evidence of the Thomson-Devenish sweeping effects, but the reported scaling laws are intermediate between those of Richardson and of Thomson-Devenish and agree with neither theory. Thus, there is still considerable uncertainty in the literature regarding the validity of the Thomson-Devenish theory. The question is important, because synthetic turbulence is a useful testing ground for numerical and theoretical methods, and because comparison of particle dispersion in synthetic and real turbulence illuminates the physical mechanisms of the latter.

Because of the disagreement of the numerical simulations of different groups, it is useful to have analytic results. The Thomson-Devenish arguments apply to a wide array of synthetic turbulence models, but Gaussian velocity ensembles are the most mathematically tractable. We therefore consider here the use of Gaussian random fields as models of turbulent Eulerian velocities. More precisely, we take the advecting velocity field \( \mathbf{u}(\mathbf{x}, t) \) to be a Gaussian random field with mean \( \overline{\mathbf{u}}(\mathbf{x}, t) \) and covariance \( C_{ij}(\mathbf{x}, t; y, s) = \langle u'_i(\mathbf{x}, t) u'_j(\mathbf{y}, s) \rangle \) for the fluctuations \( \mathbf{u}' = \mathbf{u} - \overline{\mathbf{u}} \). Specific models of interest are similar to those studied in [10], with \( \overline{\mathbf{u}}(\mathbf{x}, t) = \overline{\mathbf{u}} \) independent of space and time and with covariance defined for \( 0 < \alpha < 2, 0 < \beta < 2 \) by

\[ C_{ij}(\mathbf{x}, t; y, s) = D_2 \int d^d k \ e^{-D_3 k_L^d |t-s|} \frac{e^{i \mathbf{k} \cdot (\mathbf{x} - \mathbf{y})}}{k^{d+\alpha}} P_{ij}(\mathbf{k}) . \]  

(2)

Here \( k_L^2 = k^2 + 1/L^2 \) and \( P_{ij}(\mathbf{k}) = \delta_{ij} - k_i k_j / k^2 \) is the projection onto the subspace of \( \mathbb{R}^d \) orthogonal to \( \mathbf{k} \). The constant \( D_2 \) determines the rms velocity magnitude, while \( D_3 \) determines the intrinsic turnover rate of the synthetic turbulent eddies. The Gaussian random field \( \mathbf{u}(\mathbf{x}, t) \) so defined is statistically homogeneous in space, stationary in time, and solenoidal. The length \( L \) is proportional to the integral length-scale. The scaling properties of the model at scales smaller than \( L \) are similar to those of real turbulence. For example, the single-time covariance for \( r \ll L \) is calculated to be

\[ C_{ij}(\mathbf{x}, t; y, t) \sim D_0 L^\alpha \]  

\[ -D_1 \tau^\alpha [(d+\alpha-1) \delta_{ij} - \alpha \hat{x}_i \hat{x}_j] + O(r^2 / L^2) \]  

(3)

with \( \mathbf{r} = \mathbf{x} - \mathbf{y} \). The constants \( D_0, D_1 \) are both multiples of \( D_2 \), with constants of proportionality involving the parameters \( d \) and \( \alpha \). See [16], p.686. Kolmogorov 1941 dimensional scaling corresponds to the exponents \( \alpha = \beta = 2/3 \). We shall consider also in this paper Gaussian velocity models whose energy spectra coincide with KS models which have been studied numerically [11, 13–15]. The incompressibility of these models will be used at several points, although much of our analysis applies to more general models, e.g. with any degree of compressibility.

The principal results of this paper are as follows. For a general Gaussian model of Eulerian turbulence we carefully derive the diffusion approximation for pair dispersion assumed in the argument of Thomson-Devenish [11], under the assumption of short memory times for particle locations. We furthermore obtain a closed formula, eq.(54), for the 2-particle eddy-diffusivity in a general Gaussian model. For the specific models with covariance (2) we obtained more explicit results, which, under the conditions \( \alpha < 1 \) and either \( \beta < 1 \) or frozen turbulence with \( D_3 = 0 \), verify the Thomson-Devenish argument about sweeping decorrelation effects. In particular, we obtain under these conditions a 2-particle eddy-diffusivity tensor of the form \( K_{ij}(\mathbf{r}, t) = S_{ij}(\mathbf{r}) \tau(\mathbf{r}, t) \), where \( S_{ij}(\mathbf{r}) \) is the structure-function tensor and \( \tau(\mathbf{r}, t) \) is an effective correlation time of velocity increments. Crucially, \( \tau(\mathbf{r}, t) \) is the minimum of the intrinsic turnover time \( \tau_{eddy}(\mathbf{r}) \) at separation \( r \), the overall evolution time \( t \), and the sweeping time \( r / v_0 \). Although this result confirms the sweeping decorrelation effect, we argue that the pair-dispersion law for zero mean-velocity ensembles at high Reynolds numbers is different from the \( t^{9/2} \) suggested by Thomson & Devenish [11]. Instead, we argue that there are distinct ranges of power-laws \( t^2 \), \( t^3 \) and then \( t^4 \) again at successively longer times, exactly as Thomson & Devenish argued for ensembles with large mean velocities. We carry out careful numerical Monte Carlo simulations with our diffusion model which verify these behaviors in the model at very high Reynolds numbers. We also present Monte Carlo results for our diffusion model at the Reynolds numbers employed in current KS work, about six decades of inertial-range, and reproduce then both the “c3-law” and “c9/2-law” that have been reported in KS at comparable Reynolds numbers. We thus argue that the current KS results in the literature are not yet probing asymptotic regimes and the true
scaling in KS at very high Reynolds numbers will be the same as in our diffusion model.

The detailed analytical derivation of diffusion models is presented in section II of the paper, and predictions for their dispersion laws discussed in section III. Our numerical methods are described and validated in section IV, and then used to obtain results for mean-square particle separations and other statistics. A concluding section V briefly discusses the results.

II. DERIVATION OF THE DIFFUSION MODEL

In this section we present the derivations of our main analytical results. A reader who is only interested in physical conclusions and not the detailed justifications may skip to our final formula (54) for the pair-diffusivity and the following discussion.

A. Gaussian Integration-by-Parts Identity

We show first that the transition probability of particle pairs in Gaussian velocity ensembles obeys an exact evolution equation, as a consequence of the well-known integration-by-parts identity or Donsker-Furutsu-Novikov relation (see [17], section 4.1). Let \( u(x,t) \) be the random turbulent velocity field and let the fluid particle position that satisfies

\[
\frac{d}{dt} x(t) = u(x,t), \quad x(t_0) = a
\]  

be denoted as \( x_u(a,t_0|t) \), or \( x(a,t) \) for short. Define the “fine-grained PDF” of 2-particle positions as

\[
P_{2,u}(x_2,x_1,t|a_2,a_1,t_0) = \frac{2}{n=1}\delta^d(x_n - x_u(a_n,t_0|t)).
\]  

Then the PDF of 2-particle positions is given by

\[
P_2(x_2,x_1,t|a_2,a_1,t_0) = \langle P_{2,u}(x_2,x_1,t|a_2,a_1,t_0) \rangle,
\]

where the average is over the random velocity field \( u \).

Taking the time-derivative of (5) and using (4) it is a calculus exercise to show that

\[
\partial_t P_{2,u}(t) = -\sum_{n=1}^2 \nabla x_n \cdot \left[ \hat{P}(x_n,t) + u'(x_n,t) \right] P_{2,u}(t),
\]

(7)

where the velocity has been decomposed into its mean and fluctuating part \( u(x,t) = \hat{P}(x,t) + u'(x,t) \). The average of the second term on the righthand side can be obtained using Gaussian integration-by-parts [17]

\[
\langle u'_i(x,t) \rangle P_{2,u}(t) = \int dx \int ds C_{ik}(x,t;y,s) \\
\times \left\langle \frac{\delta}{\delta u_k(y,s)} P_{2,u}(t) \right\rangle,
\]

(8)

where \( C_{ij}(x,t;y,s) = \langle u'_i(x,t) u'_j(y,s) \rangle \). To represent the functional derivative we introduce the Lagrangian response function

\[
G_{ij}(a,t;y,s) \equiv \frac{\delta x_i(a,t)}{\delta u_j(y,s)},
\]

(9)

so that

\[
\frac{\delta}{\delta u_k(y,s)} P_{2,u}(t) = \sum_{m=1}^2 -\partial_{x_m} P_{2,u}(t) \cdot G_{jk}(a_m,t;y,s).
\]

(10)

The result of averaging (7) is the drift-diffusion equation

\[
\partial_t P_2(t) = -\sum_{n=1}^2 \nabla x_n \cdot \left[ \hat{P}(x_n,t) P_2(t) \right]
\]

\[
+ \sum_{n,m=1}^2 \partial_{x_n} \partial_{x_m} [D_{ij}(x_n,x_m,t_0) P_2(t)].
\]

(11)

with

\[
\hat{P}(x,t) = \hat{P}(x,t) + \partial_x D_{ij}(x',t_0) |_{x'=x}
\]

the mean velocity plus a fluctuation-induced drift, and with the diffusivity tensor

\[
D_{ij}(x_n,x_m,t_0) = \langle G_{jk}(a_m,t;y,s) \rangle_x (x_2,x_1,t; a_2,a_1,t_0)
\]

(13)

where

\[
\langle G_{jk}(a_m,t;y,s) \rangle_x (x_2,x_1,t; a_2,a_1,t_0) = \frac{\langle G_{jk}(a_m,t;y,s) P_{2,u}(t) \rangle}{P_2(t)}
\]

(14)

is the conditional average of the response function given that the two particles start in locations \( a_2,a_1 \) at time \( t_0 \) and end up at locations \( x_2,x_1 \) at time \( t \).

We now develop a more useful expression for the response function (9). It is straightforward to show by functional differentiation of the equation of motion (4) that

\[
\partial_t G_{ij} = \frac{\partial u_k}{\partial x_k} (x(a,t_0|t), t) G_{kj} + \delta_{ij} \delta^d(y-x(a,t_0|t)) \delta(t-s).
\]

(15)

This equation may be solved as

\[
G_{ij}(a,t;y,s) = \left\{ \begin{array}{ll}
  g_{ij}(y,s) \delta^d(y-x(a,t_0|s)) & t > s > t_0 \\
  0 & \text{o.w.}
\end{array} \right.
\]

(16)

with \( g(y,s) = Texp \left( \int_s^t dr \frac{\partial u_k}{\partial x_k} (x(a,t_0|r), r) \right) \) the time-ordered exponential matrix for the trajectory which satisfies \( x(a,t_0|s) = y \). This notation is made natural by an alternative derivation of (16) based on the flow composition identity

\[
x(a,t_0|t) = x(a,t_0|s), s|t).
\]

(17)
Taking the functional derivative $\delta / \delta u_j(y, s)$ of (17) and using the chain rule gives

$$\frac{\delta x_i(a, t)}{\delta u_j(y, s)} = \frac{\partial x_i}{\partial y_k(y, s | t)} \frac{\delta x_k(a, s)}{\delta u_j(y, s)}. \tag{18}$$

On the other hand, it is readily seen that the functional derivative of the integral form of the particle equation of motion (4), gives

$$\frac{\delta x_k(a, s)}{\delta u_j(y, s)} = \delta_{jk} \delta \delta(y - x(a, t_0 | s)) \theta(s - t_0), \tag{19}$$

with $\theta(t)$ the Heaviside step function. Thus, eq.(16) is rederived with $g_{ij}(y, s | t) = \frac{\partial x_i}{\partial y_j}(y, s | t)$. If (16) is substituted into the formula (13) it yields

$$D_{ij}(x_n, x_m, t, t_0) \equiv \int_{t_0}^{t} ds \int d^d y_m C_{ik}(x_n, s | y_m, s) \times \langle g_{jk}(y_m, s | t) | x_2, x_1, t, y_m, s; a_2, a_1, t_0 \rangle \times P(y_m, s | x_2, x_1, t; y_m, s) \quad \tag{20}$$

where $P(y_m, s | x_2, x_1, t; a_2, a_1, t_0)$ is the conditional probability density of the position of particle $m$ at time $s$ given the positions of both particles at times $t$ and $t_0$. This formula for the diffusivity when substituted into (11),(12) gives the final form of our exact evolution equation for the 2-particle transition probability.

### B. Markovian Approximation

Despite appearances, the evolution in the exact equation (11) is non-Markovian in general. It is clear from formula (20) that the 2-particle diffusion matrix is a function not only of the particle positions $x_1, x_2$ at time $t$, but also of the positions $a_1, a_2$ at time $t_0$. This dependence was suppressed in our notations, but the evolution, in principle, retains a long-time memory of the initial conditions. Only in special cases can the evolution be shown to be Markovian. The famous example is the Gaussian velocity field that is delta-correlated in time, the so-called Krichman model [18, 19], for which

$$C_{ik}(x, t; y, t) = C_{ik}(x, t; y, t) \delta(t - s). \tag{21}$$

Substituting into (20) and using

$$g_{jk}(y_m, t | t) = \delta_{jk} \tag{22}$$

and

$$P(y_m, t | x_2, x_1, t; a_2, a_1, t_0) = \delta(d)(y_m - x_m) \tag{23}$$

gives (with the “1/2 delta-function rule” for the upper limit of integration)

$$D_{ij}(x_n, x_m, t, t_0) = \frac{1}{2} C_{ij}(x_n, x_m, t). \tag{25}$$

Thus, in this case rigorously there is no dependence of the diffusion matrix $D$ on $a_1, a_2$ and the well-known diffusion model is obtained [18, 19]. Another example with Markovian particle evolution is the velocity field obtained as the superposition of Gaussian random wave trains with very high frequencies, so that the group velocity of the waves greatly exceeds the root-mean-square velocity [20]. This example has direct relevance to KS simulations with “eddy-turnover frequency” $\omega_n = \lambda \sqrt{k_n^2 E(k_n)}$ in the limit $\lambda \gg 1$ of large “unsteadiness” parameter.

The description as a diffusion should generally hold reasonably well if the correlation time of the Gaussian velocity field is short enough, since the integrand in (20) then becomes negligible at values of $s < t$ for which there is sizable dependence on $a_1, a_2$. With this motivation, we make the Markovian approximation

$$D_{ij}(x_n, x_m, t, t_0) \equiv \int_{t_0}^{t} ds \int d^d y_m C_{ik}(x_n, s | y_m, s) \times \langle g_{jk}(y_m, s | t) | x_2, x_1, t; y_m, s \rangle P(y_m, s | x_2, x_1, t). \tag{24}$$

The physical assumption is that for times ordered as $t_0 \ll s < t$ the position of the particle at time $s$ is determined mainly by its position at time $t$ and is negligibly dependent on the position at the initial time $t_0$. The worst case for this approximation is clearly the “frozen velocity” model with infinite correlation time, when times $s \geq t_0$ in the integral are not suppressed by decay of correlations. Such $s$ values give an undamped contribution also in general for times $t - t_0$ much smaller than the velocity correlation time. However, it is easy to check that the exact result (13) [or (20)] and the Markovianized result (24) both give

$$\frac{d}{dt} D_{ij}(x_n, x_m, t, t_0) = C_{ij}(x_n, t; x_m, t) + O(t - t_0) \tag{25}$$

so that, for $t - t_0$ much smaller than the correlation time,

$$D_{ij}(x_n, x_m, t, t_0) = C_{ij}(x_n, t_0; x_m, t_0)(t - t_0) + O((t - t_0)^2). \tag{26}$$

Thus the Markovian approximation becomes exact in this limit. We note in passing that the Krichman-Lundgren theory of 2-particle dispersion [21, 22] when applied to the Gaussian velocity ensemble gives a result almost identical to the formula (24) (for more discussion, see [23]).

The formula (24) from the Markovian approximation can be further simplified. It is intuitively clear that conditioning on the location of both particles is superfluous in an average of a random variable that involves only one of these particles. In fact, it can be easily established from the definitions (5),(6) that

$$P(y, s | x, x, t) = \int d^d y' P(y', y, s | x', x, t)$$

$$= \int d^d y' \langle \delta(d)(y' - x(x', t | s)) \delta(d)(y - x(x, t | s)) \rangle$$

$$= \langle \delta(d)(y - x(x, t | s)) \rangle = P(y, s | x, t). \tag{27}$$

A similar argument gives

$$\langle g_{jk}(y, s | t) | x, x, y, s \rangle = \langle g_{jk}(y, s | t) | x, y, s \rangle. \tag{28}$$
More generally, we may define the PDF
\[ P(g,t;\gamma', \gamma, s|x', x, t) = \langle \delta^3 \hat{r}_g (g - g(y, s)) \delta^4 (\gamma' - x'(t|s)) \delta^4 (\gamma - x(t|s)) \rangle \]
and mimic the previous argument to show that
\[ P(g,t;\gamma, s|x', x, t) = P(g,t;\gamma, s|x, t). \]
Then
\[ P(g,t;\gamma, s|x', x, t) = \frac{P(g,t;\gamma, s|x', x, t)}{P(y,s|x', x, t)} = \frac{P(y,s|x', x, t)}{P(y,s|x, t)} = P(g,t;\gamma, s|x, t). \]

It follows from these facts that
\[ D_{ij}(x_n, x, t, 0) = \int_0^t ds \int d^3 y M_{ik}(x_n, t; y, t, s) \times \langle g_{jk}(y, s) \rangle |x_n| \]
which is the final form of the Markovian approximation for the diffusion tensor.

We now consider the special case when the velocity field is statistically homogeneous in space. In that case, the drift velocity in (12) is independent of \( x \) and simplifies to \( \nabla \mathbf{u} (t) = \mathbf{u}(t) \), due to homogeneity and incompressibility [49]. Furthermore, a simplified equation can be derived for the transition probability of the 2-particle separation vector \( \mathbf{r} = x_2 - x_1 \), defined by
\[ P_2(r,t|r_0,0) = \int d^4 a P_2(x+r, x, t|a+r_0, a, t_0) \]
\[ = \int d^4 a P_2(x, x-r, t|a+r_0, a, t_0), \]
which is also independent of \( x \). Since the diffusion tensor \( D_{ij}(x_n, x, t) \) depends only on the difference \( x_n - x_m \) in the homogeneous case, the equation (11) with the substitutions \( \mathbf{r} = x_2 - x_1 \) and
\[ \nabla_{x_2} \longrightarrow \nabla_{\mathbf{r}}, \quad \nabla_{x_1} \longrightarrow -\nabla_{\mathbf{r}}, \]
yields the diffusion equation
\[ \partial_t P_2(r,t|r_0,0) = \partial_r \partial_{\mathbf{r}} [K_{ij}(r,t,0) P_2(r,t|r_0,0)], \]
with the eddy-diffusivity tensor
\[ K_{ij}(r,t,0) = \int_0^t ds \int d^3 y S_{ik}(r,y,t,s) \times \langle g_{jk}(y,s) \rangle |0, t| P(y,s|0, t) \]
and we define the 2nd-order structure function of velocity increments at two points \( \mathbf{0}, \mathbf{y} \) and two times \( t, s \):
\[ S_{ik}(r,y,t,s) = \langle [u_t'(r,t)-u_t'(0,t)][u_t'(y+r,s)-u_t'(y,s)] \rangle. \]

If furthermore the velocity field is assumed to be statistically stationary in time, then we can take \( t-t_0 \longrightarrow t \) and \( t_0 \longrightarrow 0 \), to obtain
\[ \partial_t P_2(r,t|r_0,0) = \partial_r \partial_{\mathbf{r}} [K_{ij}(r,t) P_2(r,t|r_0,0)], \]
with
\[ K_{ij}(r,t) = \int_0^t d\tau \int d^3 y S_{ik}(r,y,0,\tau) \times \langle g_{jk}(y,\tau|0,0; y, \tau) \rangle P(y,\tau|0,0). \]

by the change of variables \( \tau = s - t \).

C. Structure Function and One-Particle Distribution Function

The integral over \( y \) in the above formula (39) converges at large \( y \) because of decay in the two-point structure function and in the 1-particle transition probability. Physically, rapid decay is due to the facts that increments separated by great distances are uncorrelated and particles have low probability to be swept to large distances. Both of these effects can be easily quantified.

To evaluate the two-point structure function, we use a standard identity that expresses it in terms of the single-point 2nd-order structure function ([24], p.102):
\[ S_{ik}(r,y,0,\tau) = \frac{1}{2} [S_{ik}(r+y,0,\tau) + S_{ik}(r-y,0,\tau) - 2S_{ik}(y,0,\tau)]. \]

We first consider the single-time case with \( \tau = 0 \). For the spatial power-law covariance (3) with \( 0 < \alpha < 2 \), the single-point structure function becomes
\[ S_{ij}(r) = 2 \left[ C_{ij}(0,0) - C_{ij}(r,r) \right]_{\tau=0} \sim 2D_{ij} r^\alpha [(d+\alpha - 1)\delta_{ij} - \alpha \hat{r}_{ij}] + O(r^2/L^2) \]
for \( r \ll L \). The formula (40) implies in general that \( S_{ik}(r,y,0,\tau) = S_{ik}(r) \) for \( y \ll r \), whereas in the particular case (41) it gives
\[ S_{ik}(r,0,0,\tau = 0) = O(r^2/y^{2-\alpha}) \]
for \( r \ll y \ll L \). When \( y \gg L \), there is generally exponential or fast power-law decay, depending on the precise assumptions about the fall-off of the spectrum at low \( k \). The 2-time structure function \( S_{ij}(r,0,\tau) \) shows a similar behavior as the single-time structure function, except that there is a new length \( L_\beta(\tau) = (D_\beta |\tau|)^{1/\beta} \) with eddies smaller than this scale decorrelated by time \( |\tau| \). As seen from (2), the decorrelation is associated with an exponential decay of the cospectrum, with \( L_\beta(t) \) acting as
an effective “dissipation scale.” Thus, $S_{ij}(\mathbf{r}; 0, \tau)$ scales $\propto r^2$ for $r \ll L(\tau)$, while formula (41) holds for $L(\tau) \ll r \ll L$. Thus, the decay law (42) is found when $\tau \neq 0$ only for the range of values $\max\{r, L(\tau)\} \ll y \ll L$ and is limited to times $|\tau| < L^2/D_3$. For $y \ll \max\{r, L(\tau)\}$ instead $S_{ij}(\mathbf{r}; y, 0, \tau)$ is independent of $y$ and for $y \gg L$ the decay is again like that for $\tau = 0$.

The 1-particle transition probability should be dominated by large-scale sweeping and thus have the form

$$P(y, \tau|0, 0) = \frac{1}{(2\pi)^{d/2}v_0^d|\tau|} \exp\left(-|y - \mathbf{v}\tau|^2/2v_0^2\tau^2\right)$$

(43)
to a good approximation, with $v_0$ the root mean square velocity. We hereafter consider mainly the case $\mathbf{v} = 0$. For the Gaussian random field with mean zero and covariance (2), $v_0 \propto D_2L^\alpha$. In that case, it has been verified by a formal scaling analysis in [10], section 7, that the leading-order motion of particles for large $L$ is indeed ballistic with a constant, random velocity $\mathbf{v} = y/\tau$ chosen from a Gaussian ensemble with rms value $v_0$. Subleading corrections were also obtained in [10] to account for the effects of the change of the velocity in space and time. Note that (43) decays rapidly for $y \gg v_0|\tau|$.

D. Stability Matrix

The most difficult term to evaluate in (39) is the conditional average of the stability matrix $\mathbf{g}(y, \tau|0)$. Existence of this matrix requires a short-distance cutoff $\eta$ on the “inertial-range” scaling behavior in the model covariance (2) and (3), which otherwise corresponds to velocity fields only H"older continuous and non-differentiable in space. Even with the cutoff, the matrix $\mathbf{g}(y, \tau|0)$ will grow exponentially in $|\tau|$ almost surely, with rate determined by the leading Lyapunov exponent $\lambda \propto (D_1/\eta^{2-\alpha})^{1/2}$. It is thus far from clear a priori that the conditional average even remains finite in the limit $\eta \to 0$.

We begin by evaluating this term for the “frozen” velocity field with infinite correlation time (or $D_3 = 0$ in eq. (2)). A key observation here is that the Gaussian transition probability (43) implies that particles are swept from point $y$ to 0 in time $|\tau|$ along straight lines with a constant speed $v = y/\tau$ generally of order $v_0$. The velocity-gradient field $\nabla u(x, t)$ has a spatial correlation of order $\eta_\delta$, so that the particle trajectories contributing in (39) will see a constant in space but rapidly changing velocity-gradient with a correlation time $\sim \eta/v_0$. Thus, one can expect that the Lagrangian velocity-gradient will be well approximated by the model of a Gaussian field that is delta-correlated in time, for which the statistics of the stability matrix have been much studied.

To make this argument more formally, consider the spatial covariance of the velocity-gradient in the frozen case $C_{ij,mn}(r) = \langle u_{i,m}(r)u_{j,n}^*(0) \rangle$, where $u_{i,m}' = \partial u_i'/\partial x_m$. By twice differentiating (3) and then averaging over the direction of the unit vector $\hat{r}$, it is calculated to be

$$\overline{C}_{ij,mn}(r) = D_i^{\prime}r^{\alpha-2}[(d + 1)\delta_{ij}\delta_{mn} - (\delta_{im}\delta_{jn} + \delta_{in}\delta_{jm})]$$

(44)

with $D_i^{\prime} = D_i^{1/2}\frac{a_3}{\alpha-2}[\frac{a_3}{\alpha-2} + 2 + \frac{d}{\alpha-2}] > 0$ for $d \geq 2$ and $0 < \alpha < 2$. This covariance holds for $r \gg \eta_\delta$, whereas the covariance for $r < \eta \delta$ is essentially constant and can be taken to be given by (44) with $r = \eta_\delta$. A particle swept with velocity $v$ will see a random velocity-gradient with temporal correlation obtained by substituting $r = vt$ in (44). Thus, the (Eulerian) velocity-gradients in a Lagrangian frame can be taken as Gaussian with covariance

$$\langle u_{i,m}(t)u_{j,n}(0) \rangle = D_i^{\prime}t^{\alpha-1}\delta_{ij}\delta_{mn} - (\delta_{im}\delta_{jn} + \delta_{in}\delta_{jm})$$

(45)

with $D_i^{\prime} = 2\left(\frac{2-a_3}{\alpha-2}\right)D_i$ and $\delta(t) = \frac{1}{t_\eta}\Delta(\frac{1}{t_\eta})$, for $t_\eta = \eta/v$ and

$$\Delta(t) = \frac{1 - \alpha}{2(2-\alpha)} \times \left\{ \begin{array}{ll} 1 & \text{for } |t| < 1 \\ \alpha^{-2} & \text{for } |t| > 1 \end{array} \right.$$

(46)

Since $\Delta(t)$ is integrable for $\alpha < 1$ with $\int_{-\infty}^{+\infty} dt \Delta(t) = 1$, one then has $\lim_{t \to 0} \delta(t) = \delta(t)$. It follows from these arguments that the velocity-gradient experienced by the particle should be approximated by a Gaussian matrix-valued process, constant in space and delta-correlated in time, if $\alpha < 1$. This approximation could break down for fixed $\eta \ll L$ if there happens to be a small advection speed $v \ll v_0$.

Now consider the non-frozen velocity field, with covariance given by (2) and (3) for $D_3 \neq 0$. In this case the single-point, 2-time covariance of the velocity-gradient averaged over directions has the form

$$\overline{C}_{ij,mn}(r = 0, \tau) = D_i^{\prime}\eta^{\alpha-2} - D_3\eta^{2-\alpha}\delta_{ij}\delta_{mn} - (\delta_{im}\delta_{jn} + \delta_{in}\delta_{jm})$$

(47)

There is now a short correlation time $t_\eta = \eta^{\beta}/D_3$, which allows us to write

$$\langle u_{i,m}'(t)u_{j,n}^*(0) \rangle = \frac{2D_i^{\prime}}{D_3}t^{\alpha-1}\eta^{\alpha-2} - \delta_{ij}\delta_{mn} - (\delta_{im}\delta_{jn} + \delta_{in}\delta_{jm})$$

(48)

with $\delta(t) = \frac{1}{t_\eta}\Delta(\frac{1}{t_\eta})$ for $\Delta(t) = (1/2)\exp(-|t|)$. Thus, the single-point statistics of the velocity-gradient becomes temporally delta-correlated for vanishing $\eta$. In addition, there is the same decorrelation effect of rapid sweeping through space that occurs in the frozen-field case. The latter will dominate when $\eta/v \ll \eta^{\beta}/D_3$ at small $\eta$ and when the spatial decay of correlations is fast enough, that is, when both $\beta < 1$ and $\alpha < 1$. In any case, we obtain again a model for Lagrangian velocity-gradients that are Gaussian, constant in space and delta-correlated in time. There is here no problem with small
speeds \( v \ll v_0 \), since the correlation time will never be larger than \( \eta^\beta / D_3 \).

The stability matrix has been well-studied for Gaussian velocity-gradient fields, constant in space and white-noise in time. In particular, it has been shown in [25] that the matrix random process \( g(y, \tau|0) \) is a diffusion on the group \( SL(d) \) of \( d \times d \) matrices with determinant 1. We shall use specifically the formula for the transition probability density \( p_T(g) \) of this process starting at the identity, Eq.(7.14) in [25] for \( n = 2 \):

\[
\int_{SL(d)} p_T(g)f(g \cdot r_0)d\mu(g) = \int P_2(r, \tau|r_0, 0)f(r)d^dr
\]

where \( \mu \) is Haar measure on \( SL(d) \) and

\[
\partial_\tau P_2(r, \tau|r_0, 0) = M_2 P_2(r, \tau|r_0, 0)
\]

for

\[
M_2 f(r) = D((d + 1)\delta_{ij} r^2 - 2 r_i r_j) \partial_\tau \partial_{r_i} f(r) = D \partial_\tau \partial_{r_i} \left[ [(d + 1)\delta_{ij} r^2 - 2 r_i r_j] f(r) \right],
\]

where the second line follows by incompressibility. This implies also that the operator is self-adjoint. The operator \( M_2 \) is the generator for the diffusion process of particle pair-separations in the Kraichnan model with a velocity-gradient uniform in space. Pair separations in that model evolve by \( r = r \cdot \nabla u(t) \), which integrates to \( r(t) = g(t) r_0 \). The equation (49) simply expresses this relation. Since \( M_2 f \equiv 0 \) for a general linear function \( f(r) = ar, \) and considering in (49) arbitrary choices of \( a, r_0 \), it follows that

\[
\int_{SL(d)} g p_T(g)d\mu(g) = I,
\]

the identity matrix. This result is due essentially to the fact that a diffusion leaves invariant a linear profile. Finally, we can conclude that

\[
\langle g_{jk}(y, \tau|0), 0; y, \tau \rangle = \delta_{jk}.
\]

This “martingale property”, which asserts that the expected future value of the random matrix is the initial value, holds for the matrix diffusion process in the Kraichnan model but should be valid more generally. The exponential growth of the individual realizations is offset by their rapid rotation in space which leads to large cancellations in the ensemble average. Incompressibility was employed in the argument, but it is not essential to the conclusion (see [26]).

The result (53) is only strictly known to be valid when the velocity covariance converges to an \( \eta \)-independent result as \( \eta \to 0 \), whereas (45) diverges as \( \sim \eta^{\alpha-1} \) for \( \alpha < 1 \) and (48) diverges as \( \sim \eta^{\alpha+\beta-2} \) for \( \alpha + \beta < 2 \). However, the final result (53) is independent of the amplitude of the covariance (i.e. the value of \( D_1 \)) and thus we conjecture that it extends even to the present cases with diverging covariance. The result yields a simplified formula for the 2-particle eddy-diffusivity:

\[
K_{ij}(r, t) = \int_0^t d\tau \int d^d y S_{ij}(r; y, 0, \tau) P(y, \tau|0, 0),
\]

(54) together with (40),(43). We shall now use this formula to obtain concrete results for the eddy-diffusivity in the Gaussian ensembles whose covariances are given by (2).

### E. The Frozen-in-Time Velocity Field

The simplest case to analyze is the “frozen” field so that

\[
S_{ij}(r; y, 0, \tau) = S_{ij}(r; y).
\]

(55)

Using (43) for \( u = 0 \) and making the change of variables \( u = y^2/2v_0^2 r^2 \),

\[
\int_0^t d\tau P(y, \tau|0, 0) = \frac{1}{\sqrt{8\pi}^{d/2}} \frac{1}{v_0 y^d - 1} \int_0^{\frac{y^2}{2v_0^2 r^2}} dy \Gamma \left( \frac{d - 1}{2}, \frac{y^2}{2v_0^2 r^2} \right)
\]

(56) with the (upper) incomplete gamma function defined by

\[
\Gamma(s, z) = \int_z^{\infty} du u^{s-1} e^{-u}.
\]

Since \( d^d y = y^{d-1} dy d\Omega_y \), with \( d\Omega_y \) the element of \( d \)-dimensional solid angle, we get from (54) that

\[
K_{ij}(r, t) = \frac{1}{\sqrt[2]{2}\Gamma \left( \frac{d}{2} \right)} \int_0^{\infty} dy S_{ij}(r; y) \frac{y^{d-1}}{2v_0^2 r^2} \int_0^{dy} dy \Gamma \left( \frac{d - 1}{2}, \frac{y^2}{2v_0^2 r^2} \right)
\]

(57) where the angle-averaged structure function is defined by

\[
\overline{S}_{ij}(r; y) = \frac{1}{S_d} \int d\Omega_y S_{ij}(r; y)
\]

(58) for \( S_d = 2\pi^{d/2}/\Gamma \left( \frac{d}{2} \right) \) the \( (d - 1) \)-dimensional area of the unit hypersphere in \( d \)-dimensional space.

When the velocity statistics are isotropic, as for the model with zero mean and covariance (2), the eddy-diffusivity tensor can be reduced to two scalar functions \( K_L, K_N \) defined by

\[
K_{ij}(r, t) = K_L(r, t) \hat{r}_i \hat{r}_j + K_N(r, t) (\delta_{ij} - \hat{r}_i \hat{r}_j).
\]

(59)

These two functions are related by incompressibility as \( K_N = K_L + rK'_L/(d - 1) \) and it is convenient to base further analysis on \( K_L \). As is well known, if the separation statistics are also isotropic, then the diffusion equation (35) can be expressed entirely in terms of \( K_L \), as

\[
\partial_r P(r, t) = \frac{1}{r^{d-1}} \partial_r \left( r^{d-1} K_L(r, t) \partial_r P(r, t) \right)
\]

(60) Here the separation PDF satisfies

\[
\int_0^{\infty} P(r, t) r^{d-1} dr = 1.
\]

(61)
The latter integral converges for $\alpha < 1$.

For $D_3 \neq 0$. More generally, consider any velocity field statistically homogeneous in space and stationary in time. Then (32) together with (43) & (53) give

$$D_{ij}(x' - x, t) = \int_0^t ds \int dy C_{ij}(x' - y, t - s) \times \frac{1}{2\pi v_0^2(t - s)^2} exp \left[ -\frac{|y - x - \mathbf{\tau}(t-s)|^2}{2v_0^2(t-s)^2} \right]$$

Since the $y$-integration has the form of a convolution, it is easily evaluated by a Fourier transform:

$$\hat{D}_{ij}(k, t) = \int_0^t ds \hat{C}_{ij}(k, t - s) \times \exp \left[ -\frac{1}{2v_0^2 k^2(t-s)^2} \right]$$

For the model in (2) note $\hat{C}_{ij}(k, t) = \hat{C}_{ij}(k) \exp(-\gamma_k|t|)$ with $\hat{C}_{ij}(k) = D_2 P_{ij}(k)/k^{d-2+\gamma}$ and $\gamma_k = D_3 k^2$. For large $\mathbf{\tau}$, see [11]. Hereafter we take $\mathbf{\tau} = 0$. Then making the change of variables $\sigma = v_0 k(t-s)$, one obtains

$$\hat{D}_{ij}(k, t) = \frac{1}{v_0 k} \int_0^{v_0 k t} d\sigma \hat{C}_{ij}(k) \times \exp \left[ -\gamma_k \frac{\sigma - 1/2\sigma^2}{v_0^2 k^2(t-s)^2} \right]$$

Thus, for $t < 1/v_0 k$,

$$\hat{D}_{ij}(k, t) \sim \hat{C}_{ij}(k) t.$$  

This implies by inverse Fourier transform that

$$K_{ij}(x, t) \sim S_{ij}(x) t, \quad t < r/v_0.$$  

On the other hand, consider fixed $t$ and large $k$. Note that the convection time is smaller than the correlation.
time, or \( v_0 k > \gamma_k \), for \( k > k^* = (D_3/v_0)^{1/3} \) when \( \beta < 1 \).

Thus, for \( k \gg k^* \), (73) gives

\[
\dot{D}_{ij}(k,t) \sim \frac{1}{v_0 k} \mathcal{C}_{ij}(k) \cdot \sqrt{\frac{\pi}{2}} \text{erf} \left( \frac{v_0 k t}{\sqrt{2}} \right). \tag{76}
\]

This formula is exact in the case of frozen turbulence \((D_3 = 0)\) when \( k^* = 0 \). If furthermore \( k \gg 1/v_0 t \), then

\[
\dot{D}_{ij}(k,t) \sim \sqrt{\frac{\pi}{2}} \cdot \frac{1}{v_0 k} \mathcal{C}_{ij}(k) \tag{77}
\]

becomes independent of \( t \) and scales as a power \( k^{-(d+\alpha+1)} \). For \( \alpha < 1 \), we thus obtain by inverse Fourier transform that for \( r \ll \min\{v_0 t, L_*\} \)

\[
K_{ij}(r,t) \sim \sqrt{\frac{\pi}{2}} \frac{D^{(a+1)}_0}{v_0} r^{\alpha+1} \left[ (d+\alpha) \delta_{ij} - (\alpha + 1) \hat{r}_i \hat{r}_j \right]. \tag{78}
\]

It follows that the essential behavior of the frozen field case carries over to the finite time-correlated velocity with \( \alpha < 1 \) and \( \beta < 1 \). Just as for the frozen velocity, \( K_{ij}(r,t) = S_L(r) \tau (r,t) \) and the correlation time satisfies (66) and (68) with \( J_{\infty} = \sqrt{\frac{\pi}{2}} \frac{D^{(a+1)}_0}{D^{(a)}_0} [50]. \)

If, however, \( \beta > 1 \), then the behavior is quite different. Under this assumption \( \gamma_k > v_0 k \) for \( k > k^* \), so that we now make the change of variables \( \sigma = \gamma_k (t - s) \) to obtain

\[
\dot{D}_{ij}(k,t) = \frac{1}{\gamma_k} \int_0^{\gamma_k t} d\sigma \mathcal{C}_{ij}(k) \times \exp \left[ -\frac{1}{2} \left( \frac{v_0 k}{\gamma_k} \right)^2 \sigma^2 \right]. \tag{79}
\]

Equations (74) and (75) again hold, now for \( t \ll 1/\gamma_k \) and \( t \ll 1/\gamma_1 r \), respectively. On the other hand, for fixed \( t \) and \( k \gg k^* = 1/L_* \),

\[
\dot{D}_{ij}(k,t) \sim \frac{1}{\gamma_k} \mathcal{C}_{ij}(k) \left[ 1 - \exp(-\gamma_k t) \right]. \tag{80}
\]

If furthermore \( k \gg 1/L_\beta(t) \), then

\[
\dot{D}_{ij}(k,t) \sim \frac{1}{\gamma_k} \mathcal{C}_{ij}(k) \tag{81}
\]

becomes independent of \( t \) and scales as a power \( k^{-(d+\alpha+\beta)} \). When \( \alpha + \beta < 2 \), we then obtain by inverse Fourier transform that for \( r \ll \min\{L_\beta(t), L_*\} \)

\[
K_{ij}(r,t) \sim \frac{D^{(a+\beta)}_1}{D_3} r^{\alpha+\beta} \left[ (d+\alpha + \beta) \delta_{ij} - (\alpha + \beta + 1) \hat{r}_i \hat{r}_j \right]. \tag{82}
\]

We can again write \( K_{ij}(r,t) = S_L(r) \tau (r,t) \) but now

\[
\tau (r,t) \sim \begin{cases} \frac{t}{(\text{const}) \cdot D_3^{1/3}} & t \ll r^3 / D_3 \\ \frac{t}{D_3} & t \gg r^3 / D_3, r \ll L_* \end{cases}. \tag{83}
\]

Thus, the sweeping decorrelation effect is absent at sufficiently small scales when \( \beta > 1 \) and \( \alpha + \beta < 2 \).

### III. CONSEQUENCES OF DIFFUSION MODEL

In the previous section we have derived a diffusion model which, for homogeneous and isotropic statistics, takes the form (60). For the Gaussian velocity ensemble having covariance (2) with Kolmogorov scaling exponent \( \alpha = 2/3 \), the diffusivity takes the form

\[
K_L(r,t) = \frac{C_L}{v_0} \frac{2/3 \sigma^{5/3}}{D_0} J \left( \frac{v_0 t}{r} \right) \tag{84}
\]

both in the frozen case and in the temporally fluctuating case with \( \beta = 2/3 \). Here \( C_L \) is the Kolmogorov constant in the longitudinal velocity structure function, \( S_L(r) \sim C_L (\epsilon r)^{2/3} \), and \( C'_L = C_{\epsilon} J_{\infty} \). In this section we shall attempt to determine the growth law for the mean-square separation \( \langle r^2 (t) \rangle \) predicted by the model (60,84).

Does this model lead to the \( t^{9/2} \)-law of Thomson-Devenish [11]? To answer this question, we must briefly review the argument for the \( 9/2 \)-law. The key idea in [11] is that the mean-square separation pointwise in space depends on the local velocity \( \epsilon \) of the fluctuating velocity. The sweeping effect occurs at points where \( \tau_{sw}(r) = r / \epsilon' \) is smaller than the intrinsic correlation time, \( \tau_{int}(r,t) = t / \epsilon' \). The local correlation time is argued to be the smallest of these:

\[
\tau (r,t) = \min\{\tau_{sw}(r), \tau_{int}(r,t)\}. \tag{85}
\]

Hence, when \( \epsilon' > (\epsilon r)^{1/3} \) (fluctuating) or \( r / \epsilon' \) (frozen), then the mean-square separation conditioned on \( \epsilon' \) is affected by sweeping and shows the slow growth

\[
\langle r^2 (t) \rangle \sim \frac{\varepsilon t^{4/5}}{v_0^5} \tag{86}
\]

but in the opposite case exhibits the faster growth

\[
\langle r^2 (t) \rangle \sim \varepsilon t^3. \tag{87}
\]

Using these growth laws to evaluate \( \tau_{sw} \) and \( \tau_{int} \) in (85), it is easily checked that the \( t^{9/2} \)-law holds for points with \( \epsilon' > (\varepsilon t)^{1/2} \) and the \( t^3 \)-law for points with \( \epsilon' < (\varepsilon t)^{1/2} \). The probability for the latter condition to hold is small but growing in time:

\[
\text{Prob} \left( \epsilon' < (\varepsilon t)^{1/2} \right) \sim \frac{(\varepsilon t)^{3/2}}{v_0^3}. \tag{88}
\]

This formula holds for a Gaussian distribution of 3D velocities \( \epsilon' \), or for any similar distribution \( p(\epsilon') = (1/v_0^3) f(\epsilon' / v_0) \) with variance \( v_0^2 \) and non-vanishing density at the origin. The unconditional mean-square separation can then be estimated from (87) and (88) as

\[
\langle r^2 \rangle \sim \varepsilon t^3 \cdot \frac{(\varepsilon t)^{3/2}}{v_0^3} = \frac{\varepsilon^{5/2} t^{9/2}}{v_0^5}. \tag{89}
\]
The same result can be obtained from the $t^6$ dispersion law (86) by noting that it is a rapidly decreasing function of $v'$, so that the dominant contribution is obtained from the points with $v' \geq (\epsilon t)^{1/2}$ which also occur with probability $\sim (\epsilon t)^{3/2}/v_0^3$. At first sight, it appears that the model (60),(84) may embody these ideas of [11]. The diffusion model implies the exact equation

$$\frac{d}{dt} \langle r^2(t) \rangle = 2 \int K_T(r, t) P(r, t) r^{d-1} dr,$$  \hspace{1cm} (90)

where $K_T = K_L + (d - 1)K_N$ is the trace of the diffusion tensor. The average over $r$ in (90) can thus play the same role as did the average over $v'$ in the argument of [11]. The eddy-diffusivity (84) is equivalent to the correlation time (70). The population of particle pairs with separations $r > v_0 t$ should exhibit a growth law $\langle r^2(t) \rangle \sim \epsilon t^3$, while the pairs with $r < v_0 t$ should exhibit $\langle r^2(t) \rangle \sim \epsilon^{-1} t^3$. It appears possible that averaging over the entire range of pair separations could give rise to the 9/2-law (89) with an intermediate growth rate.

The above reasoning is, however, essentially wrong. The diffusion model (60),(84) does possess a $t^3$ regime, but only in an unphysical way. To see this, note that for both the $t^3$ and the $t^6$ growth laws the condition $r > v_0 t$ is first satisfied only at such long times that $t > v_0^3/\epsilon$. Substituting the standard relation $\epsilon \sim v_0^3/L$ (which follows from the assumed Kolmogorov scaling of the energy spectrum) implies that the $t^6$ law can be self-consistently satisfied only for times greater than a large-eddy turnover time, $t > L/v_0$. In that case, $\langle r^2(t) \rangle$ exceeds $L^2$ and the particle pairs have left the inertial range. As we shall verify below, the model (60),(84) does indeed possess a $t^3$ range when $t > v_0^3/\epsilon$ but this exceeds the validity of the model, which was derived only for the range $r < L$. In the range $r > L$ the two particles should instead execute independent Brownian motions with a constant diffusivity $D_T$ and the mean-square separation grow diffusively as $\langle r^2 \rangle \sim 4D_T t$. Thus, the $t^3$ range is an unphysical artefact of the model (60),(84).

The argument for the asymptotic $t^{9/2}$ law by Thomson & Devenish [11] thus fails for the model (60),(84). Very importantly, however, we shall show below that our diffusion model can produce an apparent $t^{9/2}$ law over a finite range of scales at relatively low Reynolds numbers, for similar choices of parameters with which such growth laws have been observed in kinematic simulations [11, 14, 15]. In this case, the eddy-diffusivity in the equation (60) is not given by the formula (84), which is asymptotically valid only for $L \gg r$, but instead directly from the expression (73), which holds in general. We shall thus suggest that the 9/2 growth law observed in several kinematic simulations is a finite-Reynolds-number effect and does not represent the asymptotic behavior that would be observed with very long inertial ranges.

We argue that the true high-Reynolds-number behavior—both in our diffusion model and in the kinematic simulations—is essentially the same as that found by Thomson & Devenish for the situation of large mean velocity $\overline{\nu}$ ([11], section 3.1). The principal difference is that we obtain also an early-time Batchelor ballistic range [27, 28] with $t^2$ growth. This is followed, as argued in [11], by ranges of diffusive $t^4$ growth, $t^6$ growth and finally by a range of $t^1$ or $t^3$ growth, depending upon whether the correct diffusivity (73) is used for that range or whether the $r \ll L$ approximation (84) is used (inappropriately, since $r \gg L$). We have not been able to find an analytical solution of our model (60),(84) which exhibits all of the above ranges. In this section we shall instead argue using a simple mean-field approximation

$$\frac{d}{dt} r^2 = 2 K_T(r, t), \hspace{1cm} r(0) = r_0$$  \hspace{1cm} (91)

which ignores fluctuations in the random separation $r$. In the following section IV we shall verify our theoretical conclusions by a numerical Monte Carlo solution of the diffusion model.

The Batchelor $t^2$ regime is the only one which we can derive directly from our diffusion model (60) without any approximation. We take as our initial condition for the diffusion equation the spherical delta function

$$P_0(r) = \frac{\delta(r - r_0)}{r_0^{d-1}}$$  \hspace{1cm} (92)

with all pairs initially at separation $r_0$. If this is substituted into the exact equation (90), it yields

$$\left. \frac{d}{dt} \langle r^2(t) \rangle \right|_{t=0} = 0, \hspace{1cm} \left. \frac{d^2}{dt^2} \langle r^2(t) \rangle \right|_{t=0} = 2S_T(r_0)$$  \hspace{1cm} (93)

where the trace of the short-time result (75) was used,

$$K_T(r, t) \sim 2S_T(r) t,$$  \hspace{1cm} (94)

for $t \ll r_0/v_0$. Taylor series expansion then gives

$$\langle r^2(t) \rangle - r_0^2 = S_T(r_0) t^2 + O(t^3),$$  \hspace{1cm} (95)

which is the well-known result of Batchelor [27, 28]. The mean-field approximation (91) is exact in this regime, since sufficient time has not passed to change $r$ substantially from its initial (deterministic) value $r_0$.

As noted in [11], there is an interval of times $t > r_0/v_0$ when $r$ has still not changed substantially from its initial value $r_0$. For $r \approx r_0$ but $t \gg r_0/v_0$, the result (94) is replaced with

$$K_T(r_0, t) \sim K_T(r_0, \infty) = C'_T \frac{v_0^{2/3}}{r_0^{1/3}}$$  \hspace{1cm} (96)

where $C'_T = \frac{1}{\epsilon^3} C'_T$ as a consequence of incompressibility. The growth law then becomes diffusive

$$\langle r^2(t) \rangle - r_0^2 \sim 2K_T(r_0, \infty)t,$$  \hspace{1cm} (97)

this period lasting until the “takeoff time” $t_{to}$ when $K_T(r_0, \infty)t_{to} \sim r_0^2$, or

$$t_{to} \sim \frac{r_0 v_0^{1/3}}{\epsilon^{2/3}}.$$  \hspace{1cm} (98)
See [11]. Together with the previous Batchelor regime, this diffusive range is obtained from the mean-field model (91) simplified to \( dr^2/dt = 2K_T(r_0, t) \). It is interesting that the diffusive behavior (97) at early times is the analogue in the Kraichnan white-noise advection model [18, 19] of the Batchelor ballistic range (e.g. see [29], section II.B). This is not an accident. The large-scale sweeping of particle pairs through stationary eddies produces an effective small correlation time \( r_0/v_0 \) which makes the velocity field appear to be temporally white-noise for times \( t \gg r_0/v_0 \). This is closely connected with previous attempts to simulate the Kraichnan white-noise ensemble by sweeping fixed large-scale velocity fields rapidly across the computational domain [30, 31].

For times greater than the “takeoff time” \( t_{to} \) but smaller than the “end-of-sweeping time” \( t_{es} = v_0^2/\varepsilon \), one must solve the mean-field equation (91) with

\[
dr^2/dt = 2C_T \varepsilon^{2/3} v_0^{5/3}/v_0,
\]

which leads to the \( \ell^6 \)-law (1). Instead for \( t > t_{es} \) one must solve

\[
dr^2/dt = 2C_T(\varepsilon r)^2/3t, \quad C_T = \frac{11}{3} C_L
\]

at least for the model (84). As previously discussed, this leads to the Richardson \( \ell^6 \)-law but in an unphysical way, since \( r > L \) lies outside the validity of the model (84). For \( t > t_{es} \) and \( r > L \) in reality \( K_T(r, t) \sim 2D_T \), where \( D_T \) is the 1-particle diffusivity of Taylor [32]. Thus, one must solve

\[
dr^2/dt = 4D_T
\]

which yields the very long-time diffusive range.

Our picture of particle dispersion in the zero-mean synthetic turbulence ensembles is thus very close to that in the large mean-velocity ensembles. This is in contrast to Thomson & Devenish [11], who argue for a distinct behavior of particle dispersion in the two cases. To understand better why we reach a different conclusion, it is useful to rederive our results for the eddy-diffusivity in a slightly different way. For convenience we consider only the case of frozen velocity fields. Taking the longitudinal component of the formula (54) yields

\[
K_L(r, t) = \int_{-\infty}^{0} dr \int d^d y S_L(r, y) P (y, \tau | 0, 0).
\]

As discussed in section II C the factor \( P (y, \tau | 0, 0) \) arises as the density of the Gaussian large-scale velocity \( v = y/\tau \). Changing this variable in the above integral yields

\[
K_L(r, t) = \int K_L(r, t | v) \exp \left( -\frac{v^2}{2v_0^2} \right) \frac{dv}{(2\pi v_0^2)^{d/2}}
\]

with

\[
K_L(r, t | v) = S_L(r) \tau (r, t | v)
\]

and

\[
\tau (r, t | v) = \int_{-\infty}^{0} d\tau F \left( \frac{|v|}{\tau} \right) \sim \begin{cases} t & t \ll r/v \\ \infty & t \gg r/v \end{cases}
\]

for \( \infty = \int_{-\infty}^{0} dw F (w) \). Since \( F (|v|/\tau) \) is the correlation coefficient of increments \( \delta u (r) \) at distance \( v/\tau \) apart, \( K_L(r, t | v) \) and \( \tau (r, t | v) \) can be interpreted as pair diffusivity and correlation time for given large-scale velocity magnitude \( v \). It is easy to average these quantities over \( v \) and recover the previous results for \( K_L(r, t) \) and \( \tau (r, t) \), in particular formula (70), and our predictions in this section for \( \langle r^2 (t) \rangle \). It was already observed in [11] (section 3.2, p. 292) that averaging the pair-diffusivity over the large-scale sweeping velocity would lead to the \( \ell^6 \)-law also for the zero-mean velocity ensembles. Thomson & Devenish argued, however, that correct results should be obtained by averaging \( \langle r^2 (t) \rangle \), rather than by averaging \( K_L(r, t | v) \). We find that the opposite is true. The exact integration-by-parts identity for Gaussian velocity fields leads to our formula (103) in which the effective diffusivity is indeed averaged over large-scale sweeping velocity.

IV. NUMERICAL SIMULATIONS

We now present numerical results for the diffusion models derived in the previous sections, both to confirm our theoretical predictions of their behavior and to obtain new conclusions where no analytical results are available.

A. Methods and Tests

As in [11], we shall solve the diffusion equation (38) using a Monte Carlo method for the equivalent (Ito) stochastic differential equation

\[
dr_i = b_{ij}(r, t) dW_j(t), \quad i, j = 1, ..., d
\]

where Einstein summation convention is used, \( W_j(t) \) is a vector Wiener process and \( 2K_{ij} = b_{ik}b_{jk} \), with lower-triangular square-root \( b_{ij} \) calculated by Cholesky decomposition. We can integrate the stochastic equations (106) using the standard Euler-Maruyama scheme:

\[
r_i (t_k) = r_i (t_{k-1}) + b_{ij}(r, t_{k-1}) \sqrt{\Delta t} N_{k,j} \quad i, j = 1, ..., d
\]

\[
t_k = t_{k-1} + \Delta t
\]

where \( N_{k,j} \) for \( j = 1, ..., d, k = 1, 2, 3, ... \) is an independent, identically distributed sequence of standard normal random variables. The normal random variables are obtained from uniform pseudorandom numbers generated by the Mersenne Twister algorithm [33] which are then transformed to normal by the Box-Muller method [34].

Unfortunately, the range of time that we must cover are so large that it is completely impossible for us to use a constant timestep \( \Delta t \). Instead we use an adaptive scheme
similar to that of [11]. The stepsize is determined over geometric intervals $T(m) < t < T(m + 1)$ with 

$$T(m) = A \cdot 10^{Bm} \quad \text{for} \ m = 0, 1, 2, \ldots, M.$$  \hspace{1cm} (108)

The choice of the constants $A$, $B$ and $M$ is explained in Appendix A. In each such interval, we take 

$$\Delta t = C_\Delta \min \left( \frac{r^2}{K_T(r, t)}, \Delta T \right)$$ \hspace{1cm} (109)

where $\Delta T = T(m + 1) - T(m)$ and $K_T$ is the trace of $K_T$. $C_\Delta$ is the main parameter controlling the size of the time step within the intervals $T(m) < t < T(m + 1)$. We performed convergence tests for values of $C_\Delta$ ranging from 1 to $10^{-6}$ to evaluate the quality of the timestepping. The results of these tests are presented in Appendix B. A large number $S$ of independent samples of the process (106) are generated with initial separations $r(t = 0) = r_0$ uniformly distributed over a sphere of radius $|r_0|$, and statistics obtained by averaging over realizations. Most of the results presented below used $S = 10^4$.

There is considerable debate in the literature, however, whether such adaptive time-stepping schemes lead to converged, unbiased results for the statistics [11, 13–15]. To test our numerical methods, we found it useful to consider somewhat simpler diffusion models where exact analytical results are available for comparison. The models with a power-law diffusivity 

$$K_L(r) = Dr^\zeta, \quad 0 < \zeta < 2$$ \hspace{1cm} (110)

have been very well studied. It has been shown that the long-time evolution is self-similar, with a dispersion law 

$$\langle r^2(t) \rangle \sim g(Dt)^{2/\zeta}, \quad g = \frac{\gamma^{4/\zeta} \Gamma(d + 2/\zeta)}{\Gamma(d/\zeta)}$$ \hspace{1cm} (111)

and a stretched-exponential PDF 

$$P(r, t) = \frac{1}{\langle r^2(t) \rangle^{d/2}} \exp \left[ -\alpha \left( \frac{r}{\langle r^2(t) \rangle^{1/2}} \right)^\gamma + \beta \right]$$ \hspace{1cm} (112)

where $\gamma = 2 - \zeta$, 

$$\alpha = \left[ \frac{\Gamma((d + 2)/\gamma)}{\Gamma(d/\gamma)} \right]^{\gamma/2},$$ \hspace{1cm} (113)

$$\beta = \ln \left[ \frac{\gamma^{(d + 2)/\gamma}}{\Gamma((d/\gamma) + (d + 2)/2)} \right],$$ \hspace{1cm} (114)

with the normalization condition $\int_0^\infty r^{d-1} P(r, t) dr = 1$. See [35], eqs.(3.14),(3.22) and the general, self-similar solutions found in [16] for the case of finite Corrsin invariant [51]. Incidentally, note that the mean-field equation (91) leads to power-law growth with the same exponent $2/\gamma$ as in (111) but with a different prefactor $g^{MF} = (\gamma(d + \zeta))^{2/\gamma}$ than $g$. It is not hard to show that $g^{MF} > g$, with $g^{MF} \to g$ as $d \to \infty$ from Stirling’s approximation.

Notice that the inertial-range model (84) reduces to the time-independent diffusivity 

$$K_L(r, \infty) = C_L^r \varepsilon^{2/3} v_0^5/\nu_0,$$ \hspace{1cm} (115)

as long as $r \ll v_0 t$. This is a special case of the power-law diffusivity (110) with $\zeta = 5/3$, or $\gamma = 1/3$, so that the mean-square separation grows as $t^6$. This case is thus most suitable to test our numerical methods. For the purposes of comparison in the next section with the more complex model (84), we take $D = C_L^r \varepsilon^{2/3}/v_0$ with $C_L^r = 1.256$ and $d = 3$ so that 

$$\langle r^2(t) \rangle \sim g_6^{PD} \varepsilon^{4t^6}/v_0^6$$ \hspace{1cm} (116)

with the power-law diffusion model predicting $g_6^{PD} = 15.968$. This model also has the self-similar PDF of form (112) with $d = 3, \gamma = \frac{1}{3}$, so $\alpha = 11.3714, \beta = 10.1767$.

We now employ the numerical scheme discussed earlier to see which of these exact results we can successfully reproduce. As we see in Fig. 1, long ranges of perfect $t^6$ power-laws can be obtained in log-log plots. Furthermore, Fig. 2 is a semilog plot of the dispersion

![FIG. 1: Monte Carlo results for $\langle r^2(t) \rangle$ in the power-law diffusion model (115) with $S = 10^4$ and $C_\Delta = 1$.](image-url)
14 different times in the long \( t^6 \)-range. Self-similarity is well confirmed by the collapse of rescaled curves for different times, and the analytical result (112) is accurately reproduced up to almost 16 rms values of the separation.

Our conclusion from these exercises is that the adaptive time-stepping scheme should be adequate for exponents of dispersion power-laws, and even for prefactors and PDF’s with good accuracy. Since the primary issue in this work is the exponents, we shall employ the adaptive schemes when necessary to cover extensive ranges where constant time-steps are unfeasible. As additional checks on our numerical results for exponents from adaptive schemes, we test for convergence using constants \( C_\Delta \) ranging from 1 to \( 10^{-6} \). We also compare our Monte Carlo results for the diffusion equation with a separate numerical solution of the mean-field equation (91), integrated with a Fortran 90 implementation of the Watt and Shampine RKF45 ODE solver [36, 37]. This standard ODE integration method is also adaptive, but with variable time-step determined by preselected error tolerances. We therefore can have confidence that the numerical results for the mean-field theory are well converged.

B. The Inertial-Range Model

We consider first the model (84) obtained for Kolmogorov scaling exponents in the limit \( L \gg 1 \) and thus physically applicable only for separations \( r \) in the inertial range of scales. This diffusion model applies for both the frozen velocity case and the finite-time correlated case (since \( \beta = 2/3 < 1 \)). For the purpose of simplifying the numerical work, we opted not to use the exact scaling function \( J(x) \) given by integral (67), which in three dimensions yields a complicated expression in terms of generalized hypergeometric functions. Instead, we built a function with the same asymptotic behaviors (68) as the true \( J(x) \). We took

\[
J(x) = J_\infty \operatorname{erf}(\lambda x) = \begin{cases} x & x \ll 1 \\ J_\infty & x \gg 1 \end{cases}
\]  

(117)

with \( \lambda = \frac{\sqrt{\pi}}{2} \) and \( J_\infty = \frac{D^{1/3}}{D^{1/3} + 1} \approx 0.6396 \). Our expectation was that only these general features should be sufficient to observe the scaling regimes predicted in the previous section. This idea was borne out by the numerical results. In Fig. 4 we plot \( \langle |r(t)|^2 \rangle \) for the inertial-range diffusion model with \( r_0 = 10^{-20} \). On the same graph we plot for comparison the numerical solution \( r^2(t) - r_0^2 \) of the mean-field equation (91). The two agree very well, and clearly exhibit the four predicted regimes with power-laws \( \propto t^2, t^3, t^6 \) and \( t^6 \), successively. A convergence analysis of our adaptive scheme for these results is presented in Appendix B.

To further test the theoretical predictions, we investigate the crossover times between the different regimes and the prefactors of the scaling laws. We use \( C_\Delta = 0.1 \), the largest value that gave accurate results for prefactors in the test problem. In Fig. 5(a), for example, we show for various values of \( r_0/L \) the quantity \( \langle |r(t)|^2 \rangle \) compensated by the Batchelor-range prediction \( \frac{1}{3} C_L (cr_0)^{2/3} t^{2/3} \) plotted versus the time \( t/\tau_{sw} \).
rescaled with the sweeping time $\tau_{sw} = r_0/v_0$. The Batchelor prediction fits the Monte Carlo data to within 0.13\% relative error and the end of this regime is very close to $t/\tau_{sw} = 1$. We similarly show in Fig. 5(b) for the same choices of $r_0/L$ the mean-square separation $\langle |r(t) - r_0|^2 \rangle$ compensated by the Kraichnan-like “diffusive-range” prediction $12\epsilon^2 t^{2/3}\sim 5/3 \epsilon t$ plotted versus $t/t_{to}$ with the “takeoff time” $t_{to}$ given by equation (98). The diffusive-range prediction is verified with a 1.8\% error and the end of this regime quite convincing scales as $\sim 10^{-2}t_{to}$. In Fig. 5(c) we show the corresponding plot of mean-square separation compensated by $\epsilon^2 t^6/v_0^6$ versus $t/t_{to}$. We see that a $t^6$ range begins at time $\sim 10^2 t_{to}$ and extends to the end-of-sweeping time $t_{es} = v_0^2/\epsilon$ with a prefactor $g_6^{MC} \approx 15.97$ of the $t^6$-law. This Monte Carlo value is equal within numerical errors to the exact value $g_6^{PD}$ for the power-law diffusion model (115). It is interesting that the transition between the $t^1$ and $t^6$ scaling ranges is quite broad, covering about four decades. We show finally in Fig. 5(d) the mean-square separation compensated by the Richardson prediction $\epsilon t^3$ plotted versus $t/t_{es}$. For $t > t_{es}$ there is a clear $t^3$ regime with Richardson constant $g_3^{MC} \approx 9.00$. This value agrees very well with the exact constant for a self-similar diffusion model with diffusivity $K_L(r,t) = C_L(c\tau)^{2/3} t$ and $C_L = 1.9636$. Of course, as emphasized earlier, this entire regime of the inertial-range diffusion model is unphysical and will not be observed in KS model simulations.

FIG. 4: Numerical results for $\langle |r(t) - r_0|^2 \rangle$ in the inertial-range model (84): Monte Carlo solution of the diffusion equation (8) with $C_\Delta = 1, S = 10^4$ and mean-field approximation (9).

FIG. 5: Monte Carlo results for $\langle |r(t) - r_0|^2 \rangle$. $C_\Delta = 0.1$, $S = 10^4$. Each panel shows the same curves with different scalings. (a) Batchelor regime. (b) Kraichnan regime. (c) $t^6$ regime. (d) Richardson regime. The initial separations are $r_0/L = 10^{-5}(\circ), 10^{-8}(\bigcirc), 10^{-11}(\triangle), 10^{-14}(\triangledown), 10^{-17}(\triangledown), 10^{-20}(\bigtriangledown)$. 
C. Comparison with KS Models

Our derivation of diffusion model approximations was sufficiently general that we can consider cases of more direct relevance for KS simulations, with any energy spectrum and without the approximation of large \( L \). Using the formula (76), which is exact for frozen turbulence, one obtains by inverse Fourier transform in 3D that

\[
D_{ij}(r,t) = \sqrt{\frac{\pi}{2}} \int \frac{C_{ij}(k)}{v_0k} \text{erf} \left( \frac{v_0kt}{\sqrt{2}} \right) e^{ikr}d^3k. \tag{118}
\]

It is convenient to assume statistical isotropy, so that

\[
\hat{C}_{ij}(k) = \frac{E(k)}{4\pi k^2} P_{ij}(k), \tag{119}
\]

where \( P_{ij}(k) \) is the projection operator onto the subspace orthogonal to \( k \). The trace of the diffusivity tensor becomes

\[
D_T(r,t) = \frac{\sqrt{2\pi}}{v_0} \int_0^\infty \frac{dk}{k} E(k) \text{erf} \left( \frac{v_0kt}{\sqrt{2}} \right) \frac{\sin(kr)}{kr} \tag{120}
\]

and \( D_L(r,t) \) can be recovered from

\[
D_L(r,t) = \frac{1}{v_0} \int_0^r D_T(\rho,t) \rho^{d-1}d\rho.
\]

Finally, the diffusivity that appears in equation (60) is

\[
K_L(r,t) = 2(D_L(0,t) - D_L(r,t)).
\]

To apply these results to the KS models \([11, 13–15]\), let us recall that those models have a discrete set of wavenumbers distributed as

\[
k_n = k_1 \left( \frac{k_N}{k_1} \right)^{\frac{n-1}{3}}, \tag{121}
\]

for \( n = 1, \ldots, N \) where \( k_1 = 2\pi/L \), \( k_N = 2\pi/\eta \) and \( \eta \) is the analogue of the Kolmogorov dissipation length. The energy spectrum generally adopted in these models is

\[
E(k) = C_K \varepsilon^{2/3} \sum_{n=1}^N k_n^{-5/3} \delta(k - k_n)\Delta k_n \tag{122}
\]

where \( \Delta k_n = (k_{n+1} - k_{n-1})/2 \) and \( C_K = 1.5 \) is the Kolmogorov constant, so that \( C_L \approx 1.9636 \) \([52]\). Here \( \varepsilon \) is a constant with dimensions of energy dissipation per mass chosen to prescribe values of the rms velocity:

\[
v_0 = \sqrt{\frac{2}{3}} \int_{k_1}^{k_N} E(k)dk. \tag{123}
\]

The formula (120) with the KS spectrum (122) yields

\[
D_T(r,t) = \frac{C_K \varepsilon^{2/3}}{v_0}\sqrt{\frac{2\pi}{k_n^{11/3}}} \sum_{n=1}^N \frac{\text{erf} \left( \frac{k_n r}{\sqrt{2}} \right)}{k_n^{11/3}} \sin(k_n r) \Delta k_n \tag{124}
\]

The assumption of isotropy in this formula is only approximately valid for KS simulations. It would be possible to use the general result (118), without assuming isotropy, which would lead to a discrete sum over wavevectors rather than wavevector magnitudes. However, this would make numerical implementation a bit more difficult, without essentially different physics.

We now present simulation results for diffusion models based on Gaussian velocity fields with the spectra of KS models, or, to be brief, “KS diffusion models”. The same Monte Carlo method was employed as for the inertial-range model. In all of our numerical studies we take \( v_0 = L = 1 \). We tried various values for the number of modes \( N \) and we found that the numerical results on dispersion laws in log-log plots for \( N \gtrsim 100 \) are not significantly different (see Appendix C). All of our presented results are for \( N = 500 \), a comparable number to that in the KS studies \([11, 13–15]\). We have also followed the practice in the KS literature of choosing the smallest length-scale \( \eta = r_0/10 \), for initial separation \( r_0 \). We have checked that taking \( \eta < r_0/10 \) leads to identical results. We have done no systematic study of the opposite case \( \eta > r_0/10 \), with particles starting in the “dissipation range”, when an initial regime of exponentially rapid separation would be expected.

Our first set of numerical experiments investigated whether these more realistic models would exhibit the power-law scaling ranges predicted in section III, with a \( k_N/k_1 \) sufficiently large. In Fig. 6 we plot the numerical results for the mean-square separation \( \langle |r(t) - r_0|^2 \rangle \) obtained from the KS diffusion model with \( r_0 = 10^{-20} \). We observe very clearly the predicted ranges with power-laws \( t^2, t^4, t^6 \) and, lastly, the diffusive \( t^4 \) range at long times expected for a model with finite \( L \). For comparison, we also plot numerical solutions of the mean-field equation (91) using the diffusivity (124). As before, the mean-field theory predictions are quite close to the full Monte Carlo solution of the diffusion model. Lastly, we plot the solution of the mean-field equation for the inertial-range large-\( L \) diffusivity, with the same choice of constants \( L, v_0 \) and \( \varepsilon \). As expected, the dispersion law from this approximation agrees quite well with that of the KS diffusion model for \( r < L \), but predicts a spurious \( t^3 \) power-law range for \( r > L \). The good agreement justifies \emph{a posteriori} our simplification of the scaling function \( J(x) \) in section IV B. Our most important general conclusion from this set of experiments is that the KS diffusion models and, we believe, the KS models themselves should exhibit the above four scaling ranges with successive power-laws \( t^2, t^4, t^6 \) and then \( t^4 \) again, whenever the scale ratio \( k_N/k_1 \) is sufficiently large.

In order to discriminate between various alternative theories, it is useful to compare predictions not only for mean-square separations but also for the full probability density \( P(r,t) \). In Fig. 7 we plot the Monte Carlo probability distribution calculated for 23 different times spread within the \( t^6 \) range. These are rescaled to test for self-similarity and collapse quite well. It should be empha-
sized that the overall evolution of our IR and KS diffusion models is not self-similar, globally in time. This can be seen most clearly in the existence of time ranges with distinct power-law growth laws, whereas a truly self-similar evolution should have just one power-law. In a sufficiently long $t^6$ range, however, one should expect a self-similar evolution. For example, the inertial-range model (84) in the $t^6$ range reduces to the time-independent diffusivity $K_L(r, \infty) = C^Z \varepsilon^{2/3}/\nu_0$, except for $r \gg \nu_0 t$. Since $r \sim \nu_0 t$ is nearly the maximum particle separation that can be achieved in the time $t$, only a very tiny large-$r$ tail will experience a different eddy-diffusivity than this. In Fig. 7 we also compare the Monte Carlo results for the KS diffusion model with exact parameter-free predictions (113), (114) of the power-law diffusion model (110) for $d = 3$ and $\zeta = 5/3$. The agreement is reasonably good. The curves collapse to a straight line with the stretching exponent 1/3 but with a slightly steeper slope than the power-law diffusion model. This suggests that if our Monte Carlo could be carried out for a sufficiently long $t^6$-range, then the PDF would approach the exact self-similar form of the power-law model.

We have conjectured that the growth laws of the KS models themselves, asymptotically for $k_N/k_1 \gg 1$, are the $t^2$, $t^4$, $t^6$ and $t^3$ powers that we have found in the KS diffusion models. How can this be reconciled with the $t^{9/2}$ law predicted in [11] and verified to greater or lesser extent in subsequent KS simulations [11, 14, 15]? We argue that the observed $t^{9/2}$ is an artefact of the modest $k_N/k_1$ ratios achieved in these simulations relative to ours, which tends to “blend” the distinct scaling ranges, in particular the early-time $t^1$ and $t^6$ ranges, between which lies a broad transition zone. In support of this argument, we have performed a sequence of Monte Carlo simulations of the KS diffusion model with scale ratios $k_N/k_1 = 10^3, 10^4, 10^5, 10^6$. The last ratio is chosen to correspond roughly to that employed in the previous KS simulations [11, 14, 15]. Because the range of time-scales is not so great, we have been able to carry out the time-integration not only with the adaptive algorithm employed up until now, but also with a constant time step $\Delta t = 0.1 \frac{\nu_0}{\nu_0}$ which resolves the effects of even the smallest eddies, equivalent to that used in recent KS simulations [14, 15]. The results of the two time-advancement schemes for the dispersion curves are identical when plotted in log-log. As illustrated in Fig. 8, a $t^{9/2}$ regime seems to appear as we increase the ratio $k_N/k_1$. This figure should be compared with Fig. 2 of [14] and Fig. 1 of [15], which it matches very closely. Although we see a similar “$t^{9/2}$-range” at the values of $k_N/k_1$ used in previous KS simulations, covering 1-2 decades in time, it is clear from our results in Fig. 6 that this is only a transitional regime of the KS diffusion model. In fact, for the case $k_N/k_1 = 10^6$ which shows the long “$t^{9/2}$-range” we find $t_{10^6} \approx 10^{-2}$ and thus the broad transition zone between the $t^1$ and $t^6$ laws covers the interval from $10^{-4}$ to $10^0$. This includes all of the apparent “$t^{9/2}$-range”. If we go to $k_N/k_1 = 10^8$, the power-law steepens into a

FIG. 6: Numerical results for $\langle |r(t) - r_0|^2 \rangle$ in the KS diffusion model (124), $k_N/k_1 = 10^{21}$. Monte Carlo solution of the diffusion equation (86) with $C_{\Delta} = 1, S = 10^4$ and mean-field ($\cdots$). Also MC results ($\odot$) for inertial-range model (84).

FIG. 7: Logarithm of the rescaled PDF of pair-separations at 23 different times in the $t^6$ range, for KS diffusion model with $k_N/k_1 = 10^{21}$. Monte Carlo results for $S = 10^4$ and $C_{\Delta} = 1$. The straight line marked with circles ($\odot$) is the analytical result (112) for $d = 3$ and $\gamma = 1/3$. 

\[ \frac{\ln P(r,t)/\langle r^2(t) \rangle^{3/2}}{(r/\sqrt{\langle r^2(t) \rangle})^{1/3}} \]
\(t^3\)-law. At still larger values of \(k_N/k_1\) four asymptotic scaling ranges emerge, with distinct power-law scalings of \(t^2\), \(t^3\), \(t^6\) and \(t^1\). We expect that the same is true of the KS models themselves at sufficiently large \(k_N/k_1\).

in KS there are only a single pair of discrete modes for each wavenumber magnitude and isotropy is only approximate at high wavenumbers. Perhaps the most significant difference in the model assumptions is that our diffusion theory assumes an exactly Gaussian random velocity field. The KS velocity fields instead converge to Gaussian in physical space [54] only for the limit \(N \to \infty\). Since KS calculations in the cited works use \(N\) only a few thousand at most, the Gaussian assumption is only approximately valid in physical space and not at all in Fourier space. We have employed exactly Gaussian velocity fields in our theory because this permits analytical calculations and, most importantly, because the Thomson-Devenish arguments on sweeping effects are so general that they apply to a wide range of models of Eulerian turbulence, including Gaussian velocity fields. The details of the transition to the asymptotic high Reynolds-number scalings could, however, depend upon the precise model adopted for the advecting velocity field.

Finally, we note that for \(k_N/k_1 < 10^4\), the short range of superdiffusive growth of dispersion in our KS diffusion model approximates a \(t^3\)-law. This agrees with the observations of [13, 15] for KS models. Note, however, that the physics is completely different from turbulent Richardson diffusion, which would allow \(t^3\) ranges of arbitrary extent. In fact, the narrow range of such a power-law in our KS diffusion model arises only because of the “merging” of many distinct ranges. In particular, the exponent of the apparent power-law must decrease with decreasing \(k_N/k_1\) to match the \(t^3\)-law starting at \(r = L\), until finally the superdiffusive range disappears entirely when \(k_N/k_1 \approx 1\).

\[\langle |r(t)|^2 \rangle / L^2 = \begin{cases} k_N/k_1 = 10^3 \\ k_N/k_1 = 10^4 \\ k_N/k_1 = 10^5 \\ k_N/k_1 = 10^6 \end{cases}\]

![Fig. 8: Monte Carlo results for \(\langle |r(t)|^2 \rangle / L^2\) in KS diffusion model for various values of \(k_N/k_1\), with \(\Delta t = 0.1 v_0/\nu, S = 10^4\).](image)

A pertinent question is where the predicted transition away from a \(t^{9/2}\) power should begin. The numerical results for our KS diffusion model imply that this transition should occur around \(k_N/k_1 = 10^{7−8}\). This does not agree with the KS results of [11], who found that the \(t^{9/2}\) scaling remains valid up to \(k_N/k_1 = 10^{10}\) (see Fig.12 and especially Fig.14b in [11]). It is possible that those results are vitiated by the adaptive time-integration scheme employed in that work. On the other hand, [14] using a fixed, small time-step that resolved sweeping effects found a \(t^{9/2}\) power-law in KS with \(k_N/k_1 = 10^8\), where our diffusion model produces a distinctly different \(t^5\) power. It is therefore worth considering possible sources of discrepancy between our KS diffusion model and KS itself. In the first place, it must be admitted that the Markov approximation we invoked in section II.B is a plausible but uncontrolled approximation. It is implicit also in the theoretical discussion of [11], but it may not be well-satisfied in KS, particularly for the frozen velocity case [53]. In addition, there are important differences in the underlying model assumptions of our work and the KS models of [11, 14]. Our KS diffusion model in (124) assumes a discrete set of wavenumbers with magnitudes in a geometric sequence, just like KS. However, our KS diffusion model is exactly isotropic, with the energy distributed evenly over the spheres at these radii, whereas

V. CONCLUSIONS

We have derived in this paper a diffusion equation for particle-pair dispersion in synthetic Eulerian turbulence modelled by Gaussian velocity ensembles. The main analytical result is the formula (54) for the 2-particle diffusivity and its special cases, the associated time-scale (65) for frozen velocities and the diffusivity (76) for finite time-correlated velocities. Although the description of pair-dispersion as a diffusion process is not exact (except in certain limiting cases), it arises from a well-motivated set of analytical approximations. Our results confirm the physical argument of Thomson & Devenish [11] that pair-dispersion in such models is fundamentally altered by sweeping decorrelation effects, not experienced by particle pairs in hydrodynamic turbulence. Thus, the \(t^3\)-law observed in previous simulations with synthetic turbulence is quite likely an artefact either of the numerical approximations employed [3] or of the shortness of the inertial ranges [2, 4–8]. However, we argue as well for a similar origin of the \(t^{9/2}\)-law proposed by Thomson & Devenish [11] for synthetic turbulence ensembles with zero mean velocities. Solutions of our diffusion model for such ensembles at Reynolds numbers comparable to those employed in KS simulations that show
a $t^{3/2}$-law range reproduce that finding, but our model yields instead distinct $t^2$, $t^1$, $t^6$ and $t^1$-ranges at higher Reynolds numbers. We thus argue that the asymptotic high Reynolds-number behavior of particle dispersion in synthetic Eulerian turbulence with zero mean-velocities is the same as that predicted by Thomson & Devenish [11] for ensembles with large mean velocities.

Synthetic models of turbulence such as Kinematic Simulations have been used to investigate turbulent transport of passive objects (particles, lines, etc.) in such varied problems as environmental flow, aeroacoustics, kinematic magnetic dynamo, and superfluids [38, 39]. However, such numerical studies must clearly be employed with utmost caution, especially to derive conclusions about turbulent transport at very high Reynolds numbers. Furthermore, the difference in sweeping effects in synthetic Eulerian turbulence and in real hydrodynamic turbulence could imply substantially different physics even at relatively low Reynolds numbers. It is noteworthy that several numerical simulations of pair dispersion in hydrodynamic turbulence have observed not only the $t^3$-law but also the stretched-exponential probability density function (PDF) of separations or “distance-neighbor graph” predicted by Richardson, already at modestly high Reynolds numbers [29, 40, 41]. None of the KS simulations which have observed a $t^3$-law of which we are aware [2–8] have also reported observing Richardson’s prediction for the self-similar PDF. The paper [6] observed stretched-exponential PDFs of pair separation distances in two-dimensional KS that are in good agreement with low Reynolds-number experiments in quasi-2D stratified layers [42, 43], but which do not have the 2/3 stretching exponent in Richardson’s PDF as observed in the higher Reynolds-number simulations [29, 40, 41]. This is another indication that the “$t^3$-law” observed in those KS is a non-asymptotic and non-Richardson effect arising from rather different physics. For engineering or environmental modeling purposes it may frequently not matter what is the precise origin of the $t^3$-law. However, in cases where knowledge of extreme or tail events is important (e.g. the range of dispersal of trace amounts of a dangerous contaminant), KS could lead to erroneous predictions. Discrepancies in particle dispersion properties between KS and hydrodynamic turbulence will grow even larger as the Reynolds numbers increase. There are flows with very extended inertial ranges in astrophysics (e.g. over eleven decades of Kolmogorov spectrum in the interstellar medium [44]), where actual particle dispersion properties must be quite different than what is predicted by KS at the corresponding Reynolds numbers.

Acknowledgments

We thank two anonymous referees for their careful reviews and suggestions to improve the paper. We are especially grateful to David J. Thomson for his useful comments on our paper and, in particular, for the important observation in footnote [53]. The work of GE was partially supported by the NSF Grant CMMI-0941530 at Johns Hopkins University.

Appendix A: Time-Discretization Scheme for the Monte Carlo Integrations

To construct our time-discretization scheme, we began by choosing the sequence of times

$$T(m) = A \cdot 10^{Bm} \quad \text{for } m = 0, 1, 2, ..., M.$$  \hspace{1cm} (A1)

at which we wish to record our numerical observations. To guarantee that we cover the relevant range of times to observe all of the distinct regimes, we choose $T(M) \gg t_{ex} = v_0^3/\varepsilon = 0.292$ and $T(0) \ll t_{sw} = r_0/v_0 = r_0$ (substituting our parameter values $v_0 = 1$ and $\varepsilon = 3.42$). We took $T(M) = 10^4$ and $T(0) = 10^{-5} r_0$. We chose $B = \frac{1}{\Delta v}$ so that there are 250 recorded points per decade, remaining the same as we vary $r_0$. Thus

$$A = 10^{-5} r_0, \quad M = 250 \log_{10} \left( \frac{10^9}{r_0} \right).$$

Since we chose $r_0$ to be negative integer powers of 10, $M$ defined as above is also an integer.

In each interval $T(m) < t < T(m + 1)$ we took the time-step for the numerical solution of the stochastic differential equation to be

$$\Delta t = C_\Delta \min \left( \frac{r^2}{K_T(r,t)}, \Delta T \right)$$  \hspace{1cm} (A2)

with $\Delta T = T(m + 1) - T(m)$ and $C_\Delta$ ranging from 1 to $10^{-6}$. This choice guarantees two properties. First, the square separation $r^2$ never changes by more than a fraction $C_\Delta$ of its current value in one step. Second, the time-step never exceeds $\Delta T$, so that every interval contains at least one point. Since the time series of separations $r(t_k)$ generated numerically do not hit the recording times exactly, we calculate $r(T(m))$ by linear interpolation using the nearest times before and after $T(m)$. 

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Appendix B: Convergence in Time-Step

We tested the dependence of the log-log plots of dispersion on the value of $C_\Delta$. We plot in Fig. 9 the Monte Carlo results for values of $C_\Delta$ ranging from $1$ to $10^{-6}$. There is no observable change in the behavior.

FIG. 9: Monte Carlo results for $\langle |r(t) - r_0|^2 \rangle$ in the inertial-range diffusion model calculated with $S = 10^2$ samples and varying $C_\Delta = 1$ to $C_\Delta = 10^{-6}$.

Appendix C: Convergence in Fourier Modes

We also tested the dependence of our dispersion results for the KS diffusion models on the number of Fourier modes $N$. We show in Fig. 10 log-log plots of the dispersion curves for different values of $N$, obtained from Monte Carlo calculations with $C_\Delta = 1$ and $S = 10^4$. The results are nearly indistinguishable for $N \gtrsim 100$. All of our simulations in the text used $N = 500$.

FIG. 10: Monte Carlo results for $\langle |r(t) - r_0|^2 \rangle$ in the KS diffusion model calculated with $C_\Delta = 1$, $S = 10^4$ samples, varying number of Fourier modes from $N = 10$ to $N = 10^4$.

& Majda [3] employed a homogeneous Gaussian random velocity field with zero mean which was swept by a spatially uniform “large-scale” velocity. However, this setup is statistically equivalent to a homogeneous Gaussian field with a nonzero mean and does not remove sweeping effects, in particular those of large, non-uniform eddies.

[48] The two laws can be written as $<r^2(t)>_{Rich} \sim L^2(t/t_L)^3$ and $<r^2(t)>_{TD} \sim L^2(t/t_L)^6$ in terms of the velocity integral scale $L$ and the large-eddy turnover time $t_L \sim L/v_0$, by using $\varepsilon \sim v_0^3/L$. In their regime of validity $t < t_L$, one has $<r^2(t)>_{TD} \ll <r^2(t)>_{Rich}$.

[49] This is the first point where we have invoked incompressibility.

[50] Here the superscript in $D_i^{(\alpha)}$ is used to indicate the spatial scaling exponent $\alpha$ for which the constant in (3) is calculated. Using eqs. (2.14) and (2.16) in [16] to calculate $D_i^{(\alpha)}$ gives

$$J_\infty = \sqrt{\frac{\pi}{8}} \frac{\alpha}{\sigma + 1} \Gamma \left( \frac{\sigma + 2}{2} \right) \frac{\Gamma \left( \frac{\sigma + 3}{2} \right)}{\Gamma \left( \frac{\sigma + 5}{2} \right)}$$

[51] The equations of [35] are unfortunately marred by several misprints.

[52] The standard formula $C_L = \frac{2\pi}{(3+\alpha)(2+\alpha)} \sin(\pi\alpha/2) C'K$ relates the constants $C_L, C'_K$ for a $k^{-(1+\alpha)}$ power-law spectrum; e.g. see [24], eq. (13.100). This leads to $C_L \approx 1.9727$ and $C'_K = C_L J_\infty \approx 1.2618$. However, the standard formula assumes a continuous spectrum of infinite extent, not the discrete spectrum in the KS model with finite extent. Direct numerical evaluation yielded the slightly different values $C_L \approx 1.9636$ and $C'_K \approx 1.2559$, which are thus the choice of the previous two subsections.

[53] After initial submission of this paper, it was pointed out to us by David J. Thomson (private communication) that the Markovian approximation is questionable for the conditional probability $P(y_m,s|x_0,x_1,t_0,a_0,a_i,t_0)$ in the exact diffusivity formula (20). Lagrangian particles move, to leading order, along straight lines at constant vector velocity and conditioning on the locations $x_m, a_m$ of particle $m$ at times $t, t_0$ fixes its velocity as $v_m = (x_m - a_m)/(t - t_0)$. Thus, a good approximation to $P(y_m,s|x_0,x_1,t_0,a_0,a_i,t_0)$ is the delta function $\delta(y_m - x_m - v_m(s-t))$ and not the Markov approximation $P(y_m,s|x_0,t) = \frac{\exp(-|y_m - x_m|^2/2\sigma_0^2(s-t)^2)}{(2\pi\sigma_0^2(s-t)^2)^{d/2}}$ in which $v_m$ is taken effectively to be a normal random variable with variance $\sigma_0^2$. When velocity correlation times are not very short, the approximation proposed by Thomson implies a non-Markovian evolution of the transition probability in which particles located at position $x$ at time $t$ will experience different effective diffusivities depending upon their initial location $a$ at time $t_0$, or, equivalently, depending upon their initial velocities $v(a,t_0)$. It is plausible that using this better approximation in our formula (20) will lead to the Thomson-Devenish $t^{d/2-1}$-law rather than the $t^d$-law obtained from the Markovian approximation employed in this paper.

[54] Technically, convergence of the random velocity field is proved to be weak in the topology of uniform convergence for continuous functions on physical space. See [45, 46].