A priori and a posteriori analysis of models for Large-Eddy simulation of particle-laden flow

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Abstract

In Large-Eddy simulation of particle-laden flow, the effect of the unresolved scales on the particles needs to be modelled. In this work we analyse three very promising models, namely the approximate deconvolution method (ADM) which was proposed for particle-laden flow independently by Kuerten (Phys. Fluids 18, 2006) and Shotorban and Mashayek (Phys. Fluids 17, 2005) and two stochastic models, proposed by Shotorban and Mashayek (J. Turbul. 7, 2006) and Simonin et al. (Appl. Sci. Res. 51, 1993). We present results from a priori and a posteriori analysis of these models in isotropic turbulence at $Re_\lambda = 52$. This data allows for a direct quantitative comparison of the models. The analysis shows that ADM always leads to improved statistics but that even for high Stokes numbers, the rate of dispersion is not predicted correctly by ADM. Concerning the stochastic models, we found that with the correct choice of model parameters, the models perform well at small Stokes numbers. On the other hand, at high Stokes numbers the stochastic models show significant errors such that it may be recommendable to neglect the small scale effects instead of using one of the stochastic models.

Keywords: Large-Eddy Simulation, particle-laden flow, SGS effects, Approximate Deconvolution, Langevin model

1. Introduction

Large-Eddy Simulation (LES) has become an important tool for the simulation of turbulent flow. State of the art methods provide reliable results and are capable to tackle application relevant challenges. One crucial component for LES is the correct choice of a turbulence model, i.e., a model for the effect of the unresolved subgrid scales (SGS) on the resolved scales. Such models are herein referred to as fluid-LES models.

For Large-Eddy Simulation of particle laden flow, an additional model for the effect of the unresolved scales on the particles is needed, referred to as particle-LES models. The works of Yamamoto et al. (2001); Armenio et al. (1999); Kuerten and Vreman (2005); Fedele and Simonin (2006); Yang et al. (2008); Marchioli et al. (2008) show that neglection of small scale effects is not an option.

Most particle-LES models were developed in a Eulerian-Lagrangian framework, i.e., the carrier fluid flow is computed by solving the Navier–Stokes equations and the particles are
computed by tracing single particles through the domain. Then, modelling reduces to reconstruction of small scale effect on a single particle.

On this basis, a large number of models was proposed. Among these are for example the models of Simonin et al. (1993); Wang and Squires (1996); Shotorban and Mashayek (2006); Kuerten (2006b); Amiri et al. (2006); Gobert et al. (2007); Shotorban et al. (2007); Bini and Jones (2007, 2008); Pozorski and Minier (1998); Pozorski and Apte (2009), just to mention a few. Most of these models are stochastic models, often obtained by extending models which were originally developed for inertia free particles in the context of Reynolds Averaged Navier–Stokes (RANS) simulations, such as the generalised Langevin model by Haworth and Pope (1986).

A deterministic alternative is the approximate deconvolution method (ADM) for particle laden flows (see Kuerten, 2006b; Shotorban et al., 2007; Shotorban and Mashayek, 2005). ADM is based on an approximate inversion of the LES filter and was originally developed in a Eulerian context.

The present study focusses on three very promising particle-LES models, namely the stochastic models proposed by Shotorban and Mashayek (2006) and Simonin et al. (1993) and ADM as proposed by Kuerten (2006b) and Shotorban and Mashayek (2005). For all three models, the respective authors present some results on the accuracy of their models. Their findings are summarised as follows.

Kuerten (2006b) analysed ADM in particle-laden turbulent channel flow at a Reynolds number based on friction velocity of \( Re_\tau = 150 \). He conducted an a posteriori analysis for particles with Stokes numbers of \( St = 1, 5 \) and 25, based on the viscous time scale. His results show that ADM significantly improves rms values of the wall normal component of the particle velocity. The improvement is greater for high Stokes number than for low Stokes number. In addition, Shotorban and Mashayek (2005) found that in a turbulent shear layer, ADM improves particle dispersion.

Shotorban and Mashayek (2005) analysed their Langevin-based model in decaying isotropic turbulence and found that for small Stokes numbers (\( St \leq 2.5 \) based on the Kolmogorov time scale at initialisation) the model leads to correct particle dispersion whereas at higher Stokes number significant deviations can be observed.

Fede et al. (2006) analysed the model of Simonin et al. (1993) in forced isotropic turbulence and found that the model leads to correct kinetic energy for the particles. However, their simulations are restricted to \( St \leq 5 \), based on the Kolmogorov time scale. The present study shows that at higher Stokes numbers the model does not perform very well.

Concluding, all published results were obtained on different configurations and are therefore not comparable. In particular, for the Langevin-based models only data at small Stokes numbers is published. For all models, the available data density over the Stokes number range is not satisfactory. Data rather correspond to probes at specific Stokes numbers but from this data no Stokes number dependent behaviour of the models can be deduced.

The present study aims at a clarification of that issue by providing data which allows a direct comparison of these three particle-LES models on a broad range of Stokes numbers. The data density on the Stokes number range is sufficiently high to allow the deduction of a Stokes number dependence. The testcase is isotropic turbulence at \( Re_\lambda = 52 \). All three models were originally developed such that they should perform well in that testcase but we will show that even by tuning the model constants, the models do not always perform well.
Actually in some cases better results are obtained by neglecting SGS effects than using one of the stochastic models.

This paper is organised as follows. Sections 2 and 3 contain a description of the numerical methods used to compute flow and particle dynamics. Statistics of the single phase simulations are also presented in section 2. In section 4 the three particle-LES models under consideration are presented and section 5 contains results of an a priori and an a posteriori analysis of the models.

2. Numerical Simulation of the carrier flow

In the present work we analyse particle dynamics in forced isotropic turbulence by DNS and LES. For the simulation of the carrier fluid, we use a second order Finite-Volume method together with a third order Runge-Kutta scheme proposed by Williamson (1980) for advancement in time. The conservation of mass is satisfied by solving the Poisson equation for the pressure using an iterative solver proposed by Stone (1968). More details on the flow solver can be found in Manhart (2004).

The flow is driven using a slightly modified version of the deterministic forcing scheme proposed by Sullivan et al. (1994). Sullivan et al. propose a forcing scheme where the energy in the spectral modes below a certain wave number \( \kappa_1 \) is held constant. We additionally imposed a lower bound for the forced wavenumbers, i.e., only the modes in a given range \([\kappa_0, \kappa_1]\) are forced. The Reynolds number in our simulations is always \( Re_\lambda = 52 \), based on the transverse Taylor microscale \( \lambda \) and the rms value of one (arbitrary) component of the fluctuations \( u_{rms} \).

In all computations the flow was solved in a cube on a staggered Cartesian equidistant grid. The size of the computational box and the cell width was chosen such that all scales are resolved, based on the criteria stated by Pope (2000), cf. table 1.

| Simulation parameters and Eulerian statistics from DNS of forced isotropic turbulence. | DNS | 52 |
|---------------------------------|-----|----|
| \( Re_\lambda \)                | 52  |    |
| Number of grid points \( N \)   | 256 |    |
| range of forced wavenumbers \([\kappa_0, \kappa_1]\) | [0.514, 1.54]/\( \lambda \) |    |
| integral length scale \( L_f \) | 2.00 |    |
| time scale of energy containing eddies \( k_f/\epsilon \) | 5.15\( \lambda/u_{rms} \) |    |
| Kolmogorov length scale \( \eta_K \) | 0.070 |    |
| Kolmogorov time scale \( \tau_K \) | 0.248\( \lambda/u_{rms} \) |    |
| length of computational box \( L \) | 11.9\( L_f \) |    |
| cell width \( \Delta x \)       | 1.34\( \eta_K \) |    |
| filter width \( \Delta \)       | 7\( \Delta x \) |    |
| kinetic energy of the filtered field \( \hat{k}_f \) | 0.87/\( k_f \) |    |

The particle-LES models were assessed by a priori and a posteriori analysis. For the a priori analysis, we filtered the DNS field \( \mathbf{u} \) by a box filter \( \mathcal{G} \) with filter width \( \Delta = 7\Delta x \), \( \Delta x \) being the DNS cell width. The filtered DNS was sampled on a correspondingly coarse grid,
resulting in a field $\mathcal{G}\mathbf{u}$ which is comparable to an LES field. The kinetic energy of the filtered field $\hat{k}_f = \langle \mathcal{G}u_i^2 \rangle / 2$ is 87% of the energy of the unfiltered field $k_f = \langle u_i^2 \rangle / 2$, cf. table 1. $\langle \cdot \rangle$ denotes spatial and temporal averaging.

In the a priori analysis, $\mathcal{G}\mathbf{u}$ was used as input for the particle-LES models. Then, the models were assessed with respect to the difference in statistics obtained from unfiltered DNS and filtered DNS with particle-LES model.

For the a posteriori analysis, $\mathcal{G}\mathbf{u}$ was computed by LES. As fluid-LES model we used the Lagrangian dynamic Smagorinsky model proposed by Meneveau et al. (1996). The forcing parameters for LES were chosen in the same way as for DNS, i.e., the energy contained in the range $[\kappa_0, \kappa_1]$ is equal in LES and DNS. Beyond $\kappa_1$, the energy in LES is lower than in DNS due to the different grids and the fluid-LES model. With our choice of the grid, the kinetic energy resolved by LES $\bar{k}_f$ is approximately equal to the kinetic energy of the filtered DNS field $\hat{k}_f$, cf. table 2. Instantaneous energy spectra $E(\kappa)$ from DNS and LES are plotted in figure 1. In addition, a model spectrum proposed by Pope (2000) is shown. All data were made dimension free by normalising with DNS quantities.

| Table 2: Parameters for LES of forced isotropic turbulence. |
|-----------------|------------------|------------------|
| LES             |                  |                  |
| $Re_\lambda$    | 52               |                  |
| Number of grid points $N$ | 42$^3$          |                  |
| cell width $\Delta x$     | 0.567$\lambda$  |                  |
| time scale of energy containing eddies $\bar{k}_f/\bar{\epsilon}$ | 10.4$\lambda/\bar{u}_{rms}$ |                  |
| resolved kinetic energy $k_f$ | 0.86$k_f$      |                  |

Figure 1: Instantaneous energy spectrum functions together with lines proportional to $\kappa^{-5/3}$ and $\kappa^{-7}$.

3. Discrete particle simulation

In this study we consider dilute suspensions of small particles. Thus, effects of the particles on the fluid and particle-particle interactions are neglected (one way coupling).

The density of the particles was set to $\rho_p = 1800\rho$ where $\rho$ is the density of the fluid. In each simulation the particles were divided in 24 fractions with different diameter $d$. The maximum diameter equals the Kolmogorov length scale. Consequently, the particles can be treated as point particles.
The particle relaxation time
\[ \tau_p = \frac{\rho_p d^2}{\rho 18 \nu} \] (1)
ranges from \( \tau_p = 0.1 \tau_K \) to \( \tau_p = 100 \tau_K \). Corresponding Stokes numbers \( St = \frac{\tau_p}{\tau_K} \) based on the Kolmogorov time scale \( \tau_K \) range from \( St = 0.1 \) to \( St = 100 \).

Based on the works of Armenio and Fiorotto (2001) and Kubik and Kleiser (2004), we assumed that in the given configurations the acceleration of a particle \( \frac{dv}{dt} \) is given by Stokes drag only,
\[ \frac{dv}{dt} = -c_D Re_p \frac{24}{\tau_p} (v - u_{f@p}). \] (2)

Here, \( v(t) \) denotes the particle velocity and \( u_{f@p} \) the fluid velocity at the particle position. The particle Reynolds number \( Re_p \) is based on particle diameter and particle slip velocity \( ||u_{f@p} - v|| \) which leads to a nonlinear term for the Stokes drag. The drag coefficient \( c_D \) was computed in dependence of \( Re_p \) according to the scheme proposed by Clift et al. (1978).

The fluid velocity \( u_{f@p} \) must be evaluated at the particle position \( x_p(t) \), i.e. \( u_{f@p} = u(x_p(t), t) \). Hence, these values must be interpolated. In the present work, a standard fourth order interpolation scheme was implemented, following the recommendations of Yeung and Pope (1988) and Balachandar and Maxey (1989).

In the following, the notation ‘@\( p \)’ is adopted for arbitrary functions \( f(x, t) \), i.e.,
\[ f_{@p}(t) = f(x_p(t), t). \] (3)

For example \( u_f \) refers to the space- and time-dependent solution of the Navier–Stokes equations whereas \( u_{f@p} \) refers to the time-dependent fluid velocity seen by the particle. Correspondingly, \( \mathcal{G} u_f \) refers to the space- and time-dependent solution of the filtered Navier–Stokes equations whereas \( (\mathcal{G} u_f)_{@p} \) refers to the time-dependent filtered fluid velocity seen by the particle.

Equation (2) is a stiff differential equation for small Stokes numbers. The numerical scheme for integrating equation (2) must be capable to handle this. Therefore, equation (2) was solved by a Rosenbrock-Wanner method (see Hairer and Wanner 1990). This method is a fourth order method with adaptive time stepping. The stiff term in equation (2) is linearised in each time step and discretised by an implicit Runge-Kutta scheme.

The code was validated via probability density functions (PDFs) for the particle acceleration. To this end, a DNS of forced isotropic turbulence at \( Re_\lambda = 265 \) on 1030\(^3 \) grid points was conducted. This data was then compared to data from a DNS conducted by Biferale et al. (2004) and an experiment conducted by Ayyalasomayajula et al. (2006). Biferale et al. conducted a DNS at \( Re_\lambda = 280 \) and traced inertia free particles (i.e. \( St = 0 \)). Ayyalasomayajula et al.’s experiment was at \( Re_\lambda = 250 \) with particle Stokes numbers \( St = 0.09 \pm 0.03 \). Correspondingly, in the present simulation two particle fractions were traced, one at \( St = 0 \) and another at \( St = 0.1 \). Each fraction consists of 960000 particles. Figure 2 shows that the results from the present simulations agree very well with the referenced data.
Figure 2: Probability density function $P(a)$ of particle acceleration $a$ for validation of the code. X-axis is normalised with respect to the (Stokes number dependent) rms value of $a$. Triangles: reference DNS of $St = 0$ particles conducted by Biferale et al. (2004). Squares: reference experiment of $St = 0.09 \pm 0.03$ particles conducted by Ayyalasomayajula et al. (2006) (renormalised). Continuous and dashed lines: present DNS at $St = 0$ and $St = 0.1$, respectively.

For model assessment, 24 fractions of particles were traced with 80000 particles per fraction. The particles were initialised at random positions (homogeneous distribution) inside the computational box and traced until a statistical steady state was obtained. Then, 1000 time records were taken within a time span of $T = 250\lambda/u$ for computing statistics. The temporal resolution of the statistics equals approximately the Kolmogorov time scale. With this temporal resolution, the Lagrangian correlation functions could be resolved for all Stokes numbers. The time span was large enough to guarantee that averaging in time cancels out oscillations caused by the forcing scheme.

In terms of particle time scales, $T$ is large enough to guarantee reliable statistics. From $\epsilon = 15\nu u_{rms}^2/\lambda^2$ it follows that $T/\tau_p = 250\sqrt{15}/St \approx 968.2/St$. In all simulations, $St \leq 100$, thus $T/\tau_p \geq 9.68$. Hence, statistics were sampled over at least 9.68 times the particle relaxation time.

4. Analysed particle-LES models

In the present section, the three particle-LES models under consideration are presented. They are the approximate deconvolution method which was proposed for particle-laden flow independently by Kuerten (2006b) and Shotorban and Mashayek (2005) and two stochastic models, proposed by Shotorban and Mashayek (2006) and Simonin et al. (1993). In the following, the models are stated and the numerical implementation used in this work is explained.

4.1. Approximate Deconvolution Method (ADM)
Model statement. ADM is well established for incompressible single phase flows (see Stolz and Adams (1999), Schlatter (2004), Stolz, Adams and Kleiser (2001). Kuerten (2006a,b),
Shotorban and Mashayek (2005) and Shotorban et al. (2007) analysed the capabilities of ADM for particle-laden flow. With ADM, the fluid velocity seen by the particle $u_{f_{ADM}}$ is computed from

$$u_{f_{ADM}} = (u_{f_{ADM}})_{@p} = \sum_{n=0}^{N} ((I - G)^{n} G u_{f})_{@p} = (H^{ADM} G u_{f})_{@p}. \quad (4)$$

Here, $I$ stands for identity. $N$ is the number of deconvolution steps. $H^{ADM}$ is called defiltering operator because it is supposed to approximate the inverse of $G$.

Equation (4) is solved once per time step and the particle velocity is computed from

$$\frac{du_{ADM}^{p}}{dt} = \frac{c D R e_{p}}{24 \tau_{p}} (u_{ADM}^{p} - u_{p}^{ADM}). \quad (5)$$

The operator $H = I - G$ can be interpreted as extractor of subgrid scales. With this operator, $u_{f_{ADM}}^{p}$ can be written as

$$u_{f_{ADM}}^{p} = \sum_{n=0}^{N} H^{n} G u_{f} = \sum_{n=0}^{N} H^{n} (I - H) u_{f} = (I - H^{N+1}) u_{f}. \quad (6)$$

For $N \rightarrow \infty$ the transfer function of $H^{N+1}$ equals zero for the resolvable scales ($||k|| < \kappa_{c}$) and one for the unresolvable scales ($||k|| > \kappa_{c}$). This shows that for large $N$, the effect of ADM can be interpreted as improving the LES filter towards a sharp spectral filter.

**Implementation of the model in this work.** In the present work, the ADM defiltering operator $H^{ADM}$ was computed in three different ways. First, it was computed as proposed by Kuerten (2006b). Second, it was computed making use of the DNS spectrum and third, a model spectrum was used.

If a dynamic Smagorinsky model is used as fluid-LES model, then Kuerten (2006b) proposes to compute $H^{ADM}$ as approximate inverse of the corresponding test filter. In his work and in the present work, this is a box filter. Kuerten (2006b) approximates its inverse by a second-order Taylor expansion in the filter width. The transfer function of this filter is shown in figure 3. In the following this approach is referred to as ADM$^{Kuerten}$.

However, it is not clear whether an inverted box filter gives highest accuracy. Therefore ADM was tested by two more approaches. In both approaches, the ADM filter is constructed such that the product of filter transfer function and LES spectrum is as close as possible to a target spectrum under the constraint that the filter stencil covers up to $5^3$ LES cells. The target spectrum is either the DNS spectrum or the model spectrum proposed by Pope (2000). The results from the corresponding defiltering operators are referred to as ADM$^{DNS}$ and ADM$^{mod}$, respectively. The corresponding transfer functions are also shown in figure 3. Evidently ADM$^{mod}$ leads to a very much stronger amplification around $\kappa_{c}$ than ADM$^{DNS}$. This was to be expected because around $\kappa_{c}$ the model spectrum is higher than the DNS spectrum, cf. figure 4.
ADM does not take explicitly into account that the model itself affects the particle path. More precisely, the model inherently assumes that the resolved spectrum seen by the particle is not modified by the model itself. In order to differentiate between this model assumption and other approximation errors of the model, the a priori analysis was conducted such that the model does not affect the particle path.

More precisely, in the a priori analysis for ADM, for each particle two different values for the particle velocity were computed simultaneously. One value, referred to as DNS particle velocity, is the velocity obtained from the DNS flow field. The second value, referred to as modelled particle velocity, is the velocity obtained from a filtered DNS field and ADM. The particles were tracked with the DNS particle velocity and statistical samples were taken from the modelled velocity. This approach basically tests whether ADM is capable to do what it is supposed to do, neglecting the effect of ADM on the particle path.

4.2. Langevin-based models proposed by Shotorban and Mashayek and Simonin et al.

ADM cannot reconstruct scales smaller than the LES grid. In order to circumvent this, Shotorban and Mashayek (2006) and Simonin et al. (1993) propose stochastic models based on a Langevin equation for the fluid velocity seen by a particle. Such models were originally developed for inertia free particles by Pope (1983), Heinz (2003) and Gicquel et al. (2002), referred to as generalised Langevin models.

Statement of the model proposed by Shotorban and Mashayek (2006). Shotorban and Mashayek adopted generalised Langevin models for inert particles. They propose to compute the fluid velocity seen by the particles $\mathbf{u}_{f\partial p}^{Sho}$ from the stochastic differential equation (Langevin equation)

$$
\frac{d\mathbf{u}_{f\partial p,i}^{Sho}}{dt} = \left( \mathcal{G} \left( \frac{\partial u_{f,j}}{\partial t} + u_{f,j} \frac{\partial u_{f,i}}{\partial x_j} \right) \right)_{\partial p} - \frac{\mathbf{u}_{f\partial p,i}^{Sho} - (\mathcal{G} u_{f,i})_{\partial p}}{T_L} dt + \sqrt{C_0 \epsilon} \, dW_i \quad (7)
$$
and the particle velocity from

\[ \frac{du^\text{Sho}_p}{\text{d}t} = \frac{c_D \Re_p}{24\tau_p} (u^\text{Sho}_f @p - u^\text{Sho}_p) \, \text{d}t. \]  
\[ (8) \]

The reader is reminded that ‘@p’ denotes ‘at the particle position’, cf. equation (3). The first term on the right hand side of equation (7) is the filtered material derivative of the fluid velocity and can be computed from the right hand side of the filtered Navier–Stokes equation. The second term is a drift term for the random variable \( u^\text{Sho}_f @p \), leading to a relaxation of \( u^\text{Sho}_f @p \) against \( (\mathcal{G}u_f) @p \). The last term is a diffusion term for \( u^\text{Sho}_f @p \). \( W \) denotes a Wiener process and \( \epsilon \) is the (modelled) dispersion of subgrid scale kinetic energy. The model parameters \( T_L \) and \( C_0 \) are specified below.

**Statement of the model proposed by Simonin et al. (1993).** Simonin et al. (1993) also propose to model the fluid velocity seen by the particles by a stochastic process. Fede et al. (2006) presented in detail how to deduct Simonin et al.’s model for particle-laden flow starting from the Navier–Stokes equations. This results in a different Langevin equation than the equation proposed by Shotorban and Mashayek.

In contrast to Shotorban and Mashayek, Simonin et al. propose to transport the resolved scales by particle velocity (and not by fluid velocity). The model can be formulated via a Langevin equation for the unresolved scales

\[ du^\text{Sim}'_f @p,j = \left( -u^\text{Sim}'_f @p,i \left( \frac{\partial \mathcal{G} u_f,i}{\partial x_j} \right) @p + \left( \frac{\partial \tau_{i,j}}{\partial x_j} \right) @p + \Gamma_{ij} u^\text{Sim}'_f @p,j \right) \, \text{d}t + \sqrt{C_0 \epsilon} \, dW. \]  
\[ (9) \]

\( \tau_{ij} = \mathcal{G} \left( u_i u_j \right) - \mathcal{G} u_i \mathcal{G} u_j \) is the SGS stress tensor. The model constant \( C_0 \) is equivalent to \( C_0 \) of Shotorban and Mashayek’s model. For isotropic turbulence, Fede et al. (2006) recommend

\[ \Gamma_{ij} = -\frac{\left( \frac{1}{2} + \frac{3}{4} C_0 \right)}{k_{\text{sgs}}} \epsilon \delta_{ij} = -\frac{1}{T_L} \delta_{ij}, \]  
\[ (10) \]

\( \delta_{ij} \) denoting the Kronecker delta function. This form of \( \Gamma \) was adopted in the present work.

The fluid velocity seen by the particles is then computed from \( u^\text{Sim}_f @p = (\mathcal{G} u_f) @p + u^\text{Sim}'_f @p \) and the particle velocity is computed from

\[ \frac{du^\text{Sim}_p}{\text{d}t} = \frac{c_D \Re_p}{24\tau_p} (u^\text{Sim}_f @p - u^\text{Sim}_p) \, \text{d}t. \]  
\[ (11) \]

**Implementation of the model in this work.** For model closure, two parameters need to be specified, namely the time scale \( T_L \) and the Kolmogorov constant \( C_0 \). Based on the recommendation of the model’s authors, in the present work \( T_L \) was set to

\[ T_L = \frac{k_{\text{sgs}}}{\left( \frac{1}{2} + \frac{3}{4} C_0 \right) \epsilon}, \quad \epsilon = C_\epsilon \frac{k^{3/2}}{\Delta}. \]  
\[ (12) \]

\( k_{\text{sgs}} \) denotes the subgrid kinetic energy and was computed from the DNS data. \( \epsilon \) denotes the SGS rate of dispersion. The model constants \( C_\epsilon \) and \( C_0 \) were set to \( C_\epsilon = 1 \) and \( C_0 = 2.1 \),
following Berrouk et al. (2007) and Gicquel et al. (2002). For the model of Simonin et al. (1993), the SGS stress tensor $\tau$ was computed in accordance with the fluid-LES model, i.e., using an eddy viscosity hypothesis.

As mentioned above, for the a priori analysis of ADM, the particles were traced along the path computed from DNS. For the stochastic models, this would be in contradiction to the model assumptions because the model takes explicitly into account that the particle path depends on the modelled small scale fluctuations (see Shotorban and Mashayek, 2005; Fede et al., 2006). Therefore here the particle paths were computed from the modelled fluid velocity.

The stochastic differential equations (7) and (9) were solved by an Euler-Maruyama scheme (see e.g. Kloeden and Platen, 2010). The stiff terms $-u_{f,\text{Sho}}^{\text{Sho}}/T_L$ and $\Gamma_{ij}u_{f,\text{Sim}}^{\text{Sim}}$ were discretised implicitly. Shotorban and Mashayek (2005) and Fede et al. (2006) used an explicit Euler-Maruyama scheme. These authors focussed on small Stokes numbers. In the present simulations no significant differences between explicit and implicit discretisation was found at small Stokes numbers. At high Stokes numbers, the explicit scheme was found to produce significantly worse results. In particular, the kinetic energy seen by the particles explodes at high Stokes numbers when using an explicit scheme. It should be noted that the terms under consideration are linear and therefore implicit schemes do not produce any computational overhead. In the following, ‘Sho’ denotes results from the model proposed by Shotorban and Mashayek (2005) and ‘Sim’ denotes results for the model proposed by Simonin et al. (1993).

5. A priori and a posteriori analysis of particle-LES models

The present section contains results from a priori and a posteriori analysis for assessment of the three models presented in section 4. The analysis comprises the kinetic energy seen by the particles, particle kinetic energy and rate of dispersion.

5.1. Assessment of ADM

Figures 4 to 6 show the kinetic energy seen by the particles, particle kinetic energy and the rate of dispersion for the three ADM implementations under consideration. In addition, results from filtered DNS and LES without particle-LES model are shown. In order to obtain comparable results, the presented data from filtered DNS without model corresponds to the data which ADM receives, i.e., in particular in the filtered DNS the particles were traced along unfiltered paths.

The kinetic energy seen by the particles (figure 4) shows a Stokes number dependence due to particle clustering. The Stokes number dependence can be observed in all simulations although the results from filtered DNS and LES are shifted towards higher Stokes numbers in comparison to the unfiltered DNS. This shift is not corrected by ADM.

As expected, the kinetic energy seen by the particles is lower in LES or filtered DNS than in unfiltered DNS. ADM leads to a clear improvement of the kinetic energy seen by the particles although a clear gap between results from ADM and DNS persists. This gap was to be expected because ADM does not reconstruct the smallest scales.

Among the three ADM approaches, Kuerten’s model shows least improvement. This is not surprising because Kuerten’s approach corresponds to a single defiltering step, $N = 1$. The other two approaches must perform better here because the wider stencils allow for
larger values of $N$, leading to higher kinetic energy. On the other hand it must be mentioned that $\text{ADM}^{\text{Kuerten}}$ is computationally less expensive than $\text{ADM}^{\text{DNS}}$ or $\text{ADM}^{\text{mod}}$ and somewhat more generalistic in the sense that $\text{ADM}^{\text{Kuerten}}$ only assumes a specific filter whereas $\text{ADM}^{\text{DNS}}$ and $\text{ADM}^{\text{mod}}$ are based on the spectra of this specific flow.

On first sight, the comparison of $\text{ADM}^{\text{DNS}}$ against $\text{ADM}^{\text{mod}}$ is surprising. $\text{ADM}^{\text{mod}}$ shows closer resemblance to the unfiltered result than $\text{ADM}^{\text{DNS}}$ although $\text{ADM}^{\text{DNS}}$ is based on the unfiltered field. Actually this is an effect of two errors cancelling out each other. Around the cutoff wavenumber the model spectrum is higher than the DNS spectrum but interpolation of the fluid velocity on the particle position leads to strong damping around the cutoff wavenumber. In other words, for $\text{ADM}^{\text{mod}}$ the damping properties of the interpolation scheme brings the spectrum seen by the particles closer to the spectrum of the DNS flow field.

Figure 5 shows $k_u$, the kinetic energy of the particles themselves. As expected from the observations on $k_{u\text{ap}}$, $k_u$ is underestimated by ADM. With $St \to \infty$ the error vanishes.

Most interesting is the rate of dispersion, shown in figure 6. It was computed from the product of particle kinetic energy and integral time scale. The results from a priori and a posteriori analysis differ qualitatively. In the a priori analysis, ADM leads to an underprediction of the rate of dispersion for all Stokes numbers. The reader is reminded that in the a priori analysis of ADM, the particles were traced along DNS particle paths.

The a posteriori analysis shows that actually in LES the rate of dispersion is overestimated by ADM as a result of too high integral time scales. In the a posteriori analysis of ADM, the particles were traced with the modelled particle velocity. Thus, the qualitative difference between a priori analysis and a posteriori analysis could be twofold; either due to the difference in particle paths or due to a defect in the fluid-LES model. Additional simulations showed that the principal reason is that the fluid-LES model leads to too high life times for the large eddies and thus to an overprediction of particle dispersion. Consequently,
it depends on the fluid-LES model whether the rate of dispersion is under- or overestimated by LES.

However, a priori and a posteriori analysis show that even for high Stokes numbers where small scale effects should be negligible (see Yamamoto et al. [2001]), the rate of dispersion is not predicted correctly by ADM.

![Figure 5: A priori (left) and a posteriori (right) analysis of ADM, particle kinetic energy.](image)

![Figure 6: A priori (left) and a posteriori (right) analysis of ADM, rate of dispersion. Result from LES without particle-LES model (not shown for reasons of clarity) is almost identical to the results from LES with ADM.](image)
5.2. Assessment of the Langevin-based models

The testcase for the Langevin-based models is again isotropic turbulence at \( Re_\lambda = 52 \). Figure 7 shows the kinetic energy of the fluid seen by the particles. In particular the results from the a priori analysis are very disappointing. The model of Shotorban and Mashayek (2006) shows too little \( k_{\text{up}} \) for \( St = 0.1 \) although \( k_{\text{sgs}} \) was computed from the DNS data. We found that this is an effect of the convective term in the model, \( \left( G \left( \frac{\partial u_i}{\partial t} + u_j \frac{\partial u_i}{\partial x_j} \right) \right)_{\text{up}} \). If this term is neglected, then \( k_{\text{up}} \) is correctly predicted by the model.

Concerning the a priori analysis, for low Stokes numbers the model of Simonin et al. (1993) leads to a more accurate prediction of \( k_{\text{up}} \) than the model of Shotorban and Mashayek (2006) but for high Stokes numbers one can observe the inverse. In the a posteriori analysis, the model of Simonin et al. (1993) gives a satisfactory prediction for \( k_{\text{up}} \) whereas the model of Shotorban and Mashayek (2006) leads to an overestimation.

The error in the kinetic energy of the particles is simply a consequence of the error in the kinetic energy seen by the particles, cf. figure 8. In the a priori analysis of the model of Simonin et al. (1993), the excess in \( k_{\text{up}} \) for high \( St \) is not apparent in \( k_u \). On the other hand, in the a posteriori analysis of the model of Shotorban and Mashayek (2005), the excess in \( k_{\text{up}} \) for high \( St \) is also visible in \( k_u \). This is probably due to the different time stepping. In LES, the time step size is higher than in DNS. For the particles \( St = 100 \), the ratio between particle relaxation time and time step size is about 1000 for DNS and about 200 for LES. Thus, the \( St = 100 \)-particles can rather follow the modelled fluctuations in LES than in DNS. Therefore, the excess in \( k_{\text{up}} \) at high Stokes number is rather reflected in the a posteriori analysis than in the a priori analysis.

Concerning the rate of dispersion, figure 9, the result from the a priori analysis is very discouraging, the a posteriori results are somewhat better concerning accuracy of the models. Nevertheless, these tests show that both models do not necessarily improve the result of the LES in comparison to an LES without particle-LES model or LES with ADM. In particular at high Stokes numbers it might be recommendable to use no model instead of one of the
stochastic models, in accordance with the numerical results of Shotorban and Mashayek (2005).

![Figure 8: A priori (left) and a posteriori (right) analysis of the Langevin-based models, particle kinetic energy.](image)

![Figure 9: A priori (left) and a posteriori (right) analysis of the Langevin-based models, rate of dispersion.](image)

6. Conclusions

We have presented data from DNS, filtered DNS and LES of particle-laden isotropic turbulence with three different models for the effect of the unresolved scales on the particles (particle-LES models). The models under consideration are approximate deconvolution (ADM) as proposed by Kuerten (2006b) and two stochastic models, based on the works of

![Figure 8: A priori (left) and a posteriori (right) analysis of the Langevin-based models, particle kinetic energy.](image)
Shotorban and Mashayek (2006) and Simonin et al. (1993). The present work allows for the first time a direct comparison of these models because all models were assessed on the same testcase.

The models were assessed by a priori and a posteriori analysis. The analyses comprise the kinetic energy seen by the particles, particle kinetic energy and rate of dispersion.

ADM was implemented in three different ways, one following Kuerten (2006b) and two approaches via optimisation against spectra. With all three approaches a defect in the rate of dispersion could be observed. ADM was found to underestimate the rate of dispersion if applied on the filtered DNS field. In LES however, ADM was found to overpredict the rate of dispersion. This discrepancy was explained as an effect of approximation errors in the fluid-LES model.

The stochastic models showed very poor performance in the numerical simulations. In particular for high Stokes numbers, LES or filtered DNS with stochastic models showed larger difference to DNS results than LES or filtered DNS without particle-LES model.

According to these results, the stochastic models are not recommendable because, in dependence of the configuration, LES without particle-LES model can perform better, sometimes even tremendously better, than LES with a stochastic particle-LES model. On the other hand, results from ADM are quite promising. At least ADM was found to lead to an improvement for all Stokes numbers.

On the other hand, ADM only enhances the resolved scales but does not actually model scales which cannot be represented on the grid. At high Reynolds numbers, where the LES grid is very coarse due to computational limitations, ADM can be expected to perform worse than at low Reynolds numbers. Therefore a new particle-LES model is needed for LES at high Reynolds numbers.

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