

Investigations on one-way coupling effects of particle-laden decaying isotropic turbulent flows

Julian Stemmermann, Steffen Trienekens, Christian Soika

Aachen, 3rd of July 2017

Contents

1	Nomenclature	3
2	Introduction	5
3	Mathematical models	6
3.1	Single-phase flow	6
3.1.1	The Navier-Stokes equations	6
3.2	Particle dynamics	8
4	Numerical methods	9
4.1	Direct numerical simulation	9
4.2	Particle clustering	9
5	Results	11
5.1	Boundary conditions and simulation properties	11
6	Conclusion	12
7	References	13

1 Nomenclature

E	Specific inner energy
Pr	Prandtl number
R	Universal gas constant
Re	Reynolds number
S	Sutherland temperature
T	Temperature
η	Kolomogorov scale
\boldsymbol{I}	Identity tensor
\boldsymbol{S}	Rate-of-strain-tensor
$\boldsymbol{\tau}$	Stress tensor
μ	Dynamic viscosity
ρ	Density
ρ_p	Particle density
$\tau_{m\text{athrmp}}$	Particle response time
\boldsymbol{H}^i	Stores the inviscid variables in the flux-vector included in the Navier-Stokes equations
\boldsymbol{H}^v	Stores the viscous variables in the flux-vector included in the Navier-Stokes equations
\boldsymbol{H}	Container for fluctuating variables in the Navier-Stokes equations
\boldsymbol{Q}	Container for conserved variables in the Navier-Stokes equations
∇	Nabla-operator
\boldsymbol{q}	Heat conduction
\boldsymbol{u}	Three-dimensional velocity
\boldsymbol{v}_p	Particle velocity
\boldsymbol{x}_p	Particle position
c_p	Specific isobaric heat capacity
c_v	Specific isochoric heat capacity
e	Specific internal energy
f_D	Drag correction

k_t	thermal conductivity
m_c	number of clustered particles
p	Pressure
r_p	Particle radius

2 Introduction

in Computational Methods for Multiphase Flow ist auf den Seiten 3-9 ein interessantes Beispiel.

3 Mathematical models

Sollen wir hier noch isotrope Turbulenz etc. erklären?

3.1 Single-phase flow

In this section the mathematical basics for understanding and simulating turbulent flows are discussed. However, it should be pointed out that this is no complete treatise of the mathematical and physical basics. The reader can achieve further insight on this topic by looking at different books and papers, e.g. [7].

3.1.1 The Navier-Stokes equations

The Navier-Stokes-Equations are of great importance for understanding turbulent phenomena. This set of equations exists in forms for compressible and incompressible fluids. For an infinitesimal small volume element $d\tau$ and using the cartesian coordinate system, they can be written in the so-called 'divergence form':

$$\frac{\partial \mathbf{Q}}{\partial t} + \nabla \mathbf{H} = 0 \quad (3.1)$$

It should be noted by the reader that this work only contains investigations about chemically inert fluids and particles, and that the simulation results only fit under this condition. The vector \mathbf{Q} contains all the variables which are conserved, i.e. the density ρ , the velocity \mathbf{u} and the specific inner energy E :

$$\mathbf{Q} = \begin{pmatrix} \rho \\ \rho \mathbf{u} \\ \rho E \end{pmatrix} \quad (3.2)$$

\mathbf{H} is the flux vector which stores all the floating variables and may be split up into two parts:

$$\mathbf{H} = \mathbf{H}^i + \mathbf{H}^v \quad (3.3)$$

The contents of the two vectors are displayed below:

$$\mathbf{H}^i = \begin{pmatrix} \rho \mathbf{u} \\ \rho \mathbf{u} \mathbf{u} + p \\ \mathbf{u}(\rho E + p) \end{pmatrix} \quad (3.4)$$

$$\mathbf{H}^v = -\frac{1}{Re} \begin{pmatrix} 0 \\ \boldsymbol{\tau} \\ \boldsymbol{\tau} \mathbf{u} + \mathbf{q} \end{pmatrix} \quad (3.5)$$

\mathbf{H}^i is called inviscid flux and contains only the variables that are independent of the fluids viscosity, it describes the way a fluid with zero viscosity would behave. In contrast, the viscous flux \mathbf{H}^v represents the effects of viscosity. The Reynolds number $Re = \frac{\rho v d}{\eta}$ is defined to be the ratio of inertia to tenacity, which makes it very valuable for understanding turbulent flows. This is also due to the fact that two familiar objects with the same Reynolds number behave similar in turbulence. One can assume that flows with $Re \ll 1$ are laminar and

flows with $Re \gg 1$ are turbulent. To solve the Navier-Stokes-Equations, more information regarding some variables is required. For Calculating the specific inner Energy E and the heat conduction \mathbf{q} , the following equations are used:

$$E = e \frac{1}{2} |\mathbf{u}|^2 \quad (3.6)$$

$$\mathbf{q} = - \frac{\mu}{Pr(\gamma - 1)} \nabla T \quad (3.7)$$

with

$$\gamma = \frac{c_p}{c_v} \quad (3.8)$$

and the Prandtl number

$$Pr = \frac{\mu_\infty c_p}{k_t} \quad (3.9)$$

using the specific heat capacities of the fluid c_v and c_p . If one could assume that the fluid is a newtonian fluid, the linear correlation between stress and the rate of strain results in:

$$\boldsymbol{\tau} = 2\mu \mathbf{S} - \frac{2}{3} \mu (\nabla * \mathbf{u}) \mathbf{I} \quad (3.10)$$

in which $\mathbf{S} = \frac{(\nabla \mathbf{u})(\nabla \mathbf{u})^T}{2}$ denotes the rate-of-strain-tensor. Additionally, the viscosity μ can be approximated through Sutherland's law, which is based on the ideal gas-theory:

$$\mu(T) = \mu_\infty \left(\frac{T}{t_\infty} \right)^{3/2} \frac{T_\infty + S}{T + S} \quad (3.11)$$

S is in this case the Sutherland temperature. To achieve closure the caloric state equation $e = c_v T$ and the state equation for an ideal gas $p = \rho R T$ are used. The specific gas constant is determined by $R = c_p - c_v$. These equations form a set of partial differential equations, so for solving them starting values are needed. These are initialized at the first timestep of the simulation. To achieve physical solutions, 150 timesteps are computed before the particles are initialized, so the turbulence can evolve from the synthetic values to a natural flow field.

Christoph Siewert: -2.1 bis 2.6 Stephan Fritz: -Navier-Stokes-Gleichungen (Anhang B) Randbedingungen?

3.2 Particle dynamics

Siewert: -3.1a-3.14 (spherical particles) OHNE GRAVITATION Stokes Drag/Stokes Coefficient Filterung (Fritz) -; Viskosität durch numerischen Fehler, Smagorinsky nicht benutzen

To get a useful equation of motion for the particles in the flow, we use the Euler Lagrangian approach, as it is common in Direct Numerical Simulations (DNS) and Large Eddy Simulations (LES). In this work we deal with small and heavy particles, that have a spherical shape. Their radius r_p is even smaller than the Kolmogorov scale η , but also large enough to neglect the Brownian motion. The density of the particles ρ_p is much higher than that of the fluid ρ_f . In addition due to the very low particle concentration we can neglect the influence of the particles on the mean flow and on each other. This simplification is called one-way coupling. That means that the particles can cross each other without any effect. That means particle collisions are neglected. After all this simplifications we obtain a simplified version of the Maxey-Riley equations.

$$\frac{\partial \mathbf{x}_p}{\partial t} = \mathbf{v}_p \quad (3.12)$$

$$\frac{\partial \mathbf{v}_p}{\partial t} = \frac{f_D}{\tau_p} (\mathbf{u}(\mathbf{x}_p) - \mathbf{v}_p) \quad (3.13)$$

It shows that the particle motion is only depending on the hydrodynamical drag force which results out of the actual difference between the fluid velocity $\mathbf{u}(\mathbf{x}_p)$ at particle position \mathbf{x}_p and the particle velocity \mathbf{v}_p . τ_p is the particle response time and a factor to obtain the drag force in Stokes flow conditions. To take the case of an Reynolds number \Re of order 1 into account the correction factor

$$f_D = 1 + 0.15 Re_p^{0.687} \quad (3.14)$$

is used. The biggest simplification of this Lagrangian approach is that the interaction of particles coming close together is neglected.

4 Numerical methods

To simulate flows like those described above we have two options. The direct numerical simulation (DNS) is the easier one to understand, although it is numerically very expensive. The Large-eddy simulation (LES) is numerically more capable, still we must accept certain inaccuracies. These two numerical methods are now discussed in the following chapter.

4.1 Direct numerical simulation

The basis of the direct numerical simulation (DNS) are the Navier-Stokes equations as described above. The idea is that the computer is very good at calculating and solves these equations completely. This provides us a very accurate result, as all scales of motion are being resolved. Still it requires an immense level of computational resources which increases rapidly with the Reynolds number. These computational resources were not available until the 1970s [7].

Simulating only the larger-scale motions is also called filtering, which means that the smaller-scale motions are filtered out. To model the filtered smaller-scale motions usually a subgrid-scale (SGS) model is used and this method is also called an implicit LES (ILES) [7].

4.2 Particle clustering

The high number of point particles require even more computational resources for the particle-laden simulation. The main idea to reduce this requirement is to create clusters of point particles, meaning that a new variable m_c Pojektion

(noComputationalParticles) implizite LES (Motivation fuer LES - Pope Chapter 9, Bild 9.4), DNS

5 Results

5.1 Boundary conditions and simulation properties

The simulations were carried out using ZFS, the simulation tool developed and implemented at the Institute of Aerodynamics at RWTH Aachen University [2] [3]. The tool is capable of simulating finite-volume flows of compressible fluids. In this case the turbulence was simulated on a cubic grid using 64^3 , 96^3 , 128^3 and 256^3 . The first three cases were simulated using LES, the case in which 256^3 cells were used is carried out as DNS. Further information can be gained by looking at [7, p.344-357 for DNS and p. 558-639 for LES]. For simplification, the special case of isotropic turbulence was used. For this idealised flow form the statistical velocities are invariant in all directions of the grid. It follows that the flows velocity are invariant for rotations and reflections. The turbulence was initialised using a seed-based random generator. To achieve physical results, the simulation was carried out to timestep 150, at which a restart file was written out. This procedure insures a fully developed turbulent flow, which has emancipated from the initialisation. In this flow field, a specific number of spherical particles were injected. The velocities of the fluid were interpolated for each particle to match the velocities of particles and fluid as accurately as possible.

boundary conditions, for example Reynolds number?

Graphen (particleFree rot, Laden gruen)

6 Conclusion

7 References

- [1] Edited by A. Prosperetti and G. Tryggvason. *Computational Methods for Multiphase Flow*. Cambridge University Press, 2009.
- [2] W. Schröder D. Hartmann, M. Meinke. An adaptive multilevel multigrid formulation for cartesian hierarchical grid methods. *Comput. Fluids*, 2008.
- [3] W. Schröder D. Hartmann, M. Meinke. A strictly conservative cartesian cut-cell method for compressible viscous flows on adaptive grids. *Comput. Meth. Appl. Mech. Eng.*, 2010.
- [4] S. Elghobashi. Particle-laden turbulent flows: direct simulation and closure models. *Applied Scientific Research*, 1991.
- [5] S. Fritz. Simulation isotroper turbulenz. study work, 2003.
- [6] M. Meinke K. Fröhlich, L. Schneiders and W. Schröder. Validation of particle-laden large-eddy simulation using hpc systems. In *Sustained Simulation Performance 2017*, 2017. unpublished conference document.
- [7] S. B. Pope. *Turbulent Flows*. Cambridge University Press, 2010.
- [8] W. Schröder. Fluidmechanik. lecture notes 'Fluid mechanics', 2010.