

# Turbulent channel flow tests

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## Nomenclature

### Latin symbols

|           |   |
|-----------|---|
| $\bar{u}$ | filtered stream-wise velocity               |
| $\bar{v}$ | filtered wall-normal velocity               |
| $\bar{w}$ | filtered span-wise velocity                 |
| $U$       | Mean stream-wise velocity                   |
| $u$       | stream-wise velocity fluctuation            |
| $V$       | Mean wall-normal velocity                   |
| $v$       | wall-normal velocity fluctuation            |
| $W$       | Mean span-wise velocity                     |
| $w$       | span-wise velocity fluctuation              |
| $L_x$     | Length of the domain, stream-wise direction |
| $L_y$     | Height of the domain, wall-normal direction |
| $L_z$     | Width of the domain, span-wise direction    |

### Greek symbols

|              |  |
|--------------|--|
| $\mathbb{H}$ | Quaternions                                |
| $\mathbb{C}$ | Complex Numbers                            |
| $\mathbb{R}$ | Real Numbers                               |
| $\delta$     | Half-channel height, wall-normal direction |

### Brackets, Superscripts, Sub-scripts

|        |                 |
|--------|-----------------|
| $\rho$ | Friction Index  |
| $V$    | Constant Volume |

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## **1 Introduction**

### **1.1 Literature survey**

### **1.2 Objectives and Motivation**

## **2 Lattice Boltzmann theories**

### **2.1 BGK**

### **2.2 MRT**

### **2.3 Cumulant LBM**

### 3 Turbulence theories

#### i DNS.

Direct numerical simulation (DNS) method uses a modelling-free approach for approximation of the exact solution, as the navier-stokes equation comprises of all turbulence mechanisms. Since the turbulence is three dimensional and inherently instationary it has to take in to account all the components present in the navier stokes equation. Turbulence involves the interaction in between the scales of various different sizes and thus to represent the exact solution or the characteristics of the turbulent flow considered correctly, the DNS method emphasises on the fine spatial and temporal resolution of the meshes. Apart from that this method requires the discretization methods to have low level of numerical dissipation and dispersion. In order to correctly determine the effects of the turbulence intensity, the fluctuating flow variables which satisfy the navier stokes equation are provided as inputs.

#### ii DES.

Detached eddy simulation, DES, combines the approach of RANS and the LES method; it therefore belongs to the class of the hybrid RANS-LES simulations. The basic idea of this method is that the near wall boundary layer would be resolved using the RANS approach, while the usage of LES to (correctly) resolve such regions requires a very fine spatial and temporal resolution, as the turbulent scales to be resolved in the boundary layers are very small. On the other hand, the detached regions, wakes and the free shear layers are the regions, where the RANS model frequently fails, are handled by the LES.

DES was first introduced by *Spalart et al.*. The goal was to compute the aerodynamic flows involving massive separation with a moderate increase in the grid resolution compared to the RANS method. The idea behind the DES is to segregate the flow, using suitable detectors, in to the domains where the entire energy spectrum will be modelled using the RANS (near wall boundary layer) or the large energy containing structures are resolved using the LES method (detached flow region). Therefore in this approach, a suitable method is chosen, for each domain, in terms of efficiency and accuracy.

In comparison with LES, for the computation of the flows involving the boundary layer, DES reduces the computational effort significantly. Since the flows involved are unsteady and three dimensional turbulent flows, the computational effort will be relatively higher when compared with RANS.

A very good compilation of the hybrid RANS-LES approach is found in *Fröhlich and von Terzi (2008)*

#### 3.1 RANS

#### 3.2 LES



## 4 Implementation

### 4.1 Taylor-Green vortex simulations

## 5 Model set-up and theories

Apart from the free shear flows, most turbulent flows are bounded by one or more solid surfaces depending on whether they belong to class of the internal or external flow of fluids i.e. flow through pipes and flow over a car respectively. In this investigation one of the simplest internal flow has been chosen for the validation & testing of the implementation: fully developed channel flow. This flow is considered to be of prime importance as it has played a prominent role in the development of the study of the wall bounded turbulent flows [3]. [3] has described the fundamental theories of the full developed channel flow in great detail. [4] also has a good compilation on the fully developed channel flow.

In this section a brief description of the turbulent channel flow with some basic theories is presented. Later on it is followed by description of the computational domain, meshing, initial conditions, boundary conditions, pressure gradient modelling, solver settings and in the end by post-processing.

### 5.1 Turbulent channel flow

Fully developed turbulent channel flow is a classic benchmark case and has been studied extensively to investigate and understand the mechanics of wall-bounded turbulent flows. It is the geometrical simplicity of this case that has lured to a large number of, computational (DNS) and experimental, investigations being carried out for the turbulent channel flow over the years and as a result, an adequate database (DNS data) for certain Re-numbers is already available [5]. Historical progression in the turbulent channel flow simulations have been very well documented in the literature by *Kim et al.* [6]. Also, Kim has pointed out several difficulties related to the experimental measurements and how numerical simulations serve as an essential tool, alongside the experimental results, for the study of the wall-bounded turbulence. The recent development along with the comparison of the DNS results of the turbulent channel flow simulations for higher Re-numbers is documented in [7].

As shown in figure 1, a 3D rectangular duct has been chosen as the compu-

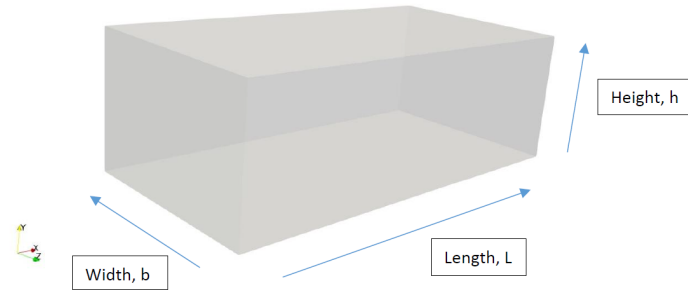


Figure 1: 3D rectangular duct

tational domain. The positive x-axis is the stream-wise direction or the flow direction with the flow varying in the y-direction, the height  $h = 2\delta$ , the wall-

normal direction. The  $z$ -axis is the span-wise direction, the width  $b$ . The domain is long and has a large aspect ratio ( $b/\delta \gg 1$ ). The bottom ( $y = 0$ ) and top ( $y = 2\delta$ ) surfaces of the domain are walls. The mid-plane ( $y = \delta$ ), which is frequently termed as the half-channel width in the literature, is ambiguous with the assigned Cartesian coordinate directions in this case and hence term half-channel height will be used ( $h = \delta$ ).

The mean velocities in the three directions are  $U$ ,  $V$ ,  $W$ , the stream-wise, wall-normal and span-wise velocities, and the corresponding fluctuating velocities are  $u, v$  and  $w$ . The mean span-wise or cross-stream velocity is zero,  $W = 0$ . In the turbulent channel flow test case, fully developed flow region in the channel is of our interest i.e. it is the region where the velocity statistics of the flow remain invariant to spatial changes,  $x$ -direction in our case. Also, the domain is large enough in the span-wise direction,  $z$ , compared to the half-channel height,  $\delta$ , so that the flow remains statistically independent of the  $z$ -direction. In short, we have two homogeneous flow directions along  $x$  &  $z$ -axis. Thus, the flow is said to be statistically one- dimensional as the flow statistics vary only in one-direction i.e.  $y$ -direction.

### 5.1.1 Governing equations

Turbulent channel flow is fully developed, where the velocity statistics along the stream-wise direction remain constant i.e.  $\partial U / \partial x = 0$ , and statistically one dimensional, where the flow statistics vary only in the wall-normal( $y$ ) direction. There is no mean flow in the span-wise direction i.e  $W = 0$  and thus the gradient of the span-wise velocity in the span-wise direction is zero i.e.  $\partial W / \partial z = 0$ . The mean continuity equation for the incompressible flow results in the mean wall normal velocity gradient to be zero:

$$\partial V / \partial y = 0 \quad (1)$$

The eq. 1 when integrated, within the wall limits  $[0, 2\delta]$ , gives  $V = \text{const}(y)$  and using the no-slip boundary condition,  $U_i = 0$ , applied at the walls results in  $V = 0$ . Now to satisfy the resulting continuity equation eq. 1,  $V$  has to be zero everywhere in the domain. Using the above made considerations and applying it to the mean stream-wise momentum equation we will have the following force balance:

$$\frac{\partial \tau}{\partial y} = \frac{\partial P}{\partial x} \quad (2)$$

where  $\tau = \left( \mu \frac{\partial U}{\partial y} - \rho \overline{u_i u_j} \right)$  is the total shear stress.

The eq. 2 shows that the total shear stress gradient and the pressure gradient balance each other in the full developed channel flow. It shows an equality of the form  $f(x) = g(y)$  and for this equality to hold it is important that both  $f(x)$  and  $g(y)$  are individually constants. In eq. 2 both the gradients are dependent on different variables and thus they must be individually constants for this equality to hold.

$$\frac{\partial P}{\partial x} = \text{const} \ \& \ \frac{\partial \tau}{\partial y} = \text{const}$$

Total shear stress is the summation of the viscous stress,  $\mu \frac{\partial U}{\partial y}$ , and the Reynolds

stress,  $-\rho \overline{u_i u_j}$ . At the wall with the no-slip boundary condition applied, the Reynolds stresses are reduced to zero and thus the shear stress near the wall or to be specific at the wall would entirely be dominated by the viscous stresses. At the walls the Reynolds stresses drop to zero and the stress at the walls can be defined as the *wall shear stress*, the region where the viscous stresses are dominant:

**wall shear stress**

$$\begin{aligned}\tau_w &= \mu \frac{\partial U}{\partial y} \\ \tau(0) &= \tau_w = -\tau(2\delta)\end{aligned}\tag{3}$$

Now we can integrate the eq. 2, from 0 to  $2\delta$ , for the pressure gradient and express the solution in terms of the  $\tau_w$ :

$$-\frac{\partial P}{\partial x} = \frac{\tau_w}{\delta}, \tag{4}$$

From the fully developed channel flow it is known that the mean velocity profile (time-averaged) remains constant. It is clear from eq. 3 that wall shear stress is linearly proportional to the velocity gradient near the wall. A constant velocity profile implies that the wall shear stress will also be constant in a fully developed channel flow. Thus from eq. 4 we can say that the pressure gradients is constant. Integrating the eq. 2, from 0 to some  $y$ , with the no-slip boundary condition at the wall results in the following solution:

$$\tau(y) = \tau_w \left(1 - \frac{y}{\delta}\right) \tag{5}$$

It is this pressure gradient that drives the flow through the infinite parallel plates. The pressure gradient in eq. 4 is negative as the viscous effects decreases the pressure along the stream-wise direction. For a given pressure gradient and the half-channel height we get a linear shear stress profile from eq. 4 and 5 independent of the fluid properties and the state of the fluid motion (laminar or turbulent) [3]. Experimental and numerical results [6] show that the total shear stress profile is anti-symmetric about the channel mid height  $y = \delta$  and thus  $\tau(\delta) = 0$ .

### 5.1.2 Scaling units

The Reynolds stresses decreases towards the wall, remain zero at the wall and so does the eddy viscosity (*Boussinesq hypothesis*). The value of the eddy viscosity ranges from zero, at the wall, to several orders of magnitude higher than the molecular viscosity, in the core region of the turbulent boundary layer [1]. From the figure. 2 it can be seen that the velocity profile changes more slowly (more flatter compared to the parabolic flow profile) in the core of the turbulent boundary layer, where the eddy viscosity is higher than the molecular viscosity, and there is an abrupt change in the velocity profile near the wall, where the molecular viscosity is dominant. The turbulent flow profile has the largest gradient near the wall.

Turbulent flow along the wall can be considered to consist of four different layers, characterised by their distance from the wall. All four layers have been

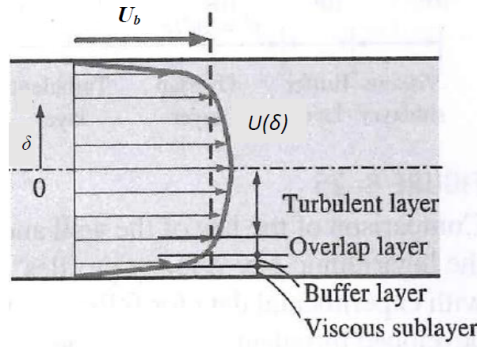


Figure 2: Velocity profile in the fully developed channel flow [1]

approximately shown in figure 2. The layer closest to the wall is the viscous sub-layer, where the viscous effects are dominant. The velocity profile changes linearly with the distance from the wall and the flow is streamlined. Next to the viscous sub-layer is the buffer layer, where the turbulent stresses start to develop, but still the viscous stresses are dominant. Above the buffer layer is the overlap layer, where the turbulent stresses are significant, but still not dominant enough as the viscous stresses. The layer above the overlap layer for the rest of the flow is the outer or turbulent layer, where the turbulence stresses dominate over the viscous stresses.

A single analytical relation for the entire turbulent velocity profile is not suitable, as different regions or layers execute different flow characteristics. The best possible option is to find the key parameters or the functional forms using the dimensional analysis and later on experimental data can be used to obtain the numerical values of the any constants.

**i. Viscous sub-layer :** With the no-slip boundary condition applied at the wall,  $u_i(0, t) = 0$ , the entire contribution to the shear stress is purely via viscous stresses i.e.  $\tau_w$ . From eq. 3 it is clear that shear stress is proportional to the molecular viscosity. The thickness of this layer is very small, but it plays an important role on the flow characteristic because of the presence of large velocity gradients in it. The velocity profile remains linear in this layer and also it has been shown in certain experiments [1]. Since the velocity profile is linear the gradient in the equation 3 remains constant i.e.  $\partial U / \partial y = U / y$ . Thus the wall shear stress can be written as:

$$\frac{\tau_w}{\rho} = \frac{\nu U}{y} \quad (6)$$

Near the wall the important parameters are :  $\tau_w$ ,  $\rho$ ,  $\nu$ . The square root of the term on the left hand side has the dimensions of velocity and thus we define a new parameter as the **friction velocity**,  $u_\tau$ :

$$u_\tau = \sqrt{\frac{\tau_w}{\rho}} \quad (7)$$

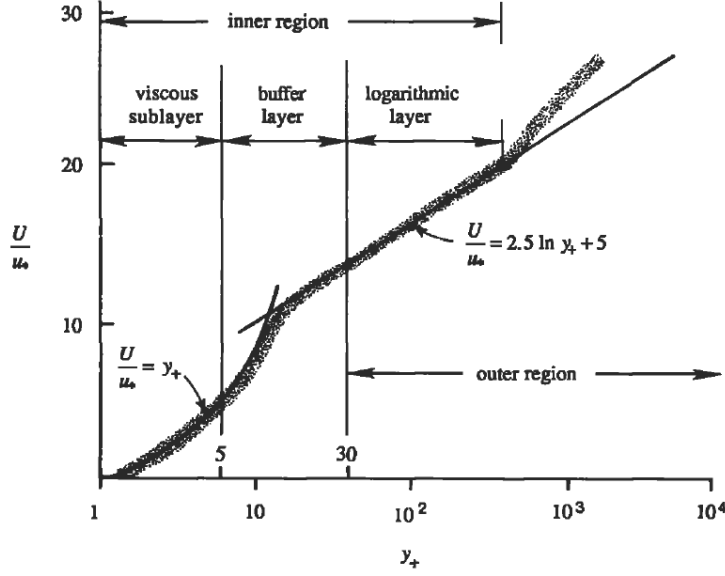


Figure 3: Nondimensional velocity  $u^+$  vs. the nondimensional distance  $y^+$  (plotted in logarithmic scale) [2]

Replacing  $u_\tau$  in the eq. 6 we have *the law of wall*:

$$\frac{U}{u_\tau} = \frac{u_\tau y}{\nu} \quad (8)$$

The quantity  $\nu/u_\tau$  has the dimensions of length and it is called **viscous length**,  $\delta_\nu$ , and it is used to nondimensionalize the distance from the wall,  $y$ . In the analysis of the boundary layer flow, it is convenient to work with the nondimensionalised distance and the velocity. The nondimensionalised quantities are represented with a +.

**Nondimensionalised variables :**

$$y^+ = \frac{yu_\tau}{\nu} \text{ and } u^+ = \frac{U}{u_\tau} \quad (9)$$

**Normalised law of the wall :**

$$u^+ = y^+ \quad (10)$$

This law shows a very good correlation with the experimental data for smooth surfaces for  $0 \leq u_\tau y/\rho \leq 5$ . Thus for the  $y^+ = 5$  we will have the linear velocity profile in the viscous sub-layer.

**ii. Overlap layer:** From the dimensional analysis and the experimental results it proved that the velocity in the overlap region is proportional to the logarithm of distance and it can be expressed as:

$$u^+ = \frac{1}{\kappa} \ln y^+ + B \text{ or } u^+ = 2.5 \ln y^+ + 5.0 \quad (11)$$

where  $\kappa$  and  $B$  are constants and equal to 0.40 and 5.0 respectively. The eq. 11 is called the *logarithmic law* and the overlap layer is also called as the *logarithmic layer* or *log-law layer*. It is clear from the figure 3 that the eq. 11 represents the experimental data quite satisfactorily from  $y^+ \geq 30$  till outer region. Neither eq. 11 nor the eq. 8 represents the experimental data satisfactorily in the buffer layer. From Figure. 3 it is seen that :

**viscous sub-layer** :  $y^+ \leq 5$

**buffer layer** :  $y^+ \geq 5$  till  $y^+ \leq 30$

**logarithmic layer** :  $y^+ \geq 30$  till  $y^+ \leq 10^3$

iii. **Outer layer:** The velocity law for this region is:

$$\frac{(U_{max} - U)}{u_\tau} = \frac{1}{\kappa} \ln \left( \frac{y}{\delta} \right) \quad (12)$$

The difference between the mean center line velocity,  $U_{max}$ , and the mean velocity,  $U$  is called the **velocity defect** and thus this law is called the **velocity defect-law**. In this region the turbulent stresses are dominant compared to the viscous stresses and this is also seen from the eq. 12: the normalised velocity profile in the core region of the turbulent channel flow is dependent on the half-channel height and is not dependent on the viscosity.

## 5.2 Reynolds number

Since the Reynolds number is used to characterize the flow, great care was taken to simulate the same flows i.e. same Reynolds number as that of the DNS data was used. The dimensionless Reynolds number requires three parameters for its definition: characteristics length ( $\delta$ ), characteristic velocity ( $V_{ch}$ ) and the kinematic viscosity ( $\nu$ ) of the fluid.

$$Re = \frac{V_{ch} \delta}{\nu}$$

Several choices of  $V_{ch}$  are possible, but in this investigation two different  $V_{ch}$  have been chosen and the respective Reynolds number are as follows :

i. **Friction Reynolds number ( $Re_\tau$ ):**

$$Re_\tau = \frac{u_\tau \delta}{\nu} = \frac{\delta}{\delta_\nu} \quad (13)$$

ii. **Bulk Reynolds number ( $Re_b$ ):**

$$Re_b = \frac{U_b 2\delta}{\nu} \quad (14)$$

The bulk mean velocity is defined as:

$$U_b = \frac{1}{h} \int_0^h U dy$$

### 5.3 Computational Domain and Meshing

#### 5.3.1 Computational Domain

We are interested to investigate the fully developed turbulent channel and that implies a sufficient length and width of the domain for the flow in the channel to be fully developed. Since the stream-wise and the span-wise directions are homogeneous, periodic boundary conditions are applied on either side of the two homogeneous directions. All flow properties along with the state of the flow are set to be equal for the equivalent points on the periodic boundary pairs. Use of periodic boundaries is equivalent to having an infinitely long domain in that direction. This implies that we have a flow between two infinite parallel planes, Figure 4. Since the artificial periodic boundaries are enforced great care

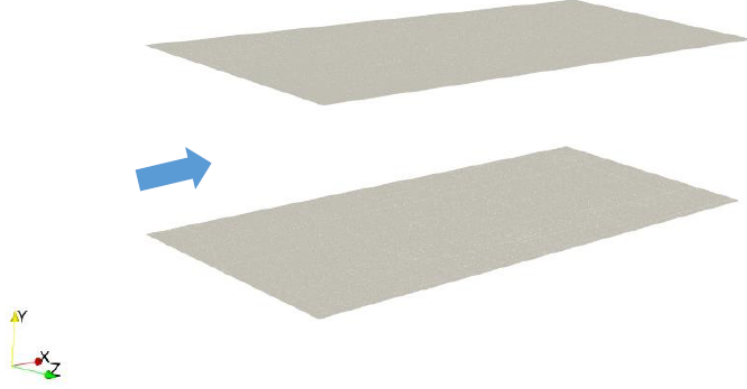


Figure 4: Flow between the infinite parallel plates

has to be taken that the domain is long enough to avoid this artificial boundary conditions from influencing the results. *Moin et.al* [9] has specified certain aspects to be considered in the process of choosing the dimensions of the flow domain for the numerical simulation of the turbulent channel flow. According to [9] the experimental two-point correlation functions have to be considered for choosing the stream-wise ( $L_x$ ) and the span-wise ( $L_z$ ) lengths. The usage of periodic boundary conditions is justified if the distance between the pair of periodic boundaries is at least twice the distance, at which the correlation function becomes zero. This means that the domain size is enough to accommodate the largest eddy without the pair of periodic boundaries influencing each other.

To validate the LES results of the fully developed turbulent channel flow the DNS data from the database of [8] for the  $Re_\tau = 395$  is used. The authors have shown the adequacy of the domain by performing the two-point spatial correlations. Same domain size used in [8] has been chosen for performing the LES of channel flow, Table 1. Thus, no separate study for two-point correlation has been performed in this investigation.

Several authors [10, 11] have performed the turbulent channel flow simulations with smaller domain dimensions compared to the dimensions used in the refer-



| $Nominal Re_\tau$ | $L_x \times L_y \times L_z$                  |
|-------------------|--|
| 395               | $2\pi\delta \times 2\delta \times \pi\delta$ |

Table 1: The size of the domain in x, y, z directions

ence data, [8], in order to either save the computational time or to statiate the upper limit of the total gird points on a Graphic card [11]. But *Bespalko* [12] suggests that for performing the validation studies the dimensions of the domain should be similar to that used in the reference data. Else, the usage of smaller domain dimensions might add some kind of error and this might be than attributed to the error originating from the implementation under test.

### 5.3.2 Physical parameters

The physical parameters that are necessary for the model set-up and later on for the simulations are pressure gradient(to drive the flow), discussed in the later section,  $u_\tau$ ,  $\nu$  and  $U_b$ . In [8] the authors have not specified  $u_\tau$ ,  $\nu$  and  $U_b$  used in the simulations. [13] has computed and specified these parameters obtained from the DNS results of [8] for  $Re_\tau = 395$ , Table 2 :

| <i>Physical quantity</i>      | <i>Value in S.I Units</i> |
|-------------------------------|---------------------------|
| Half channel height, $\delta$ | 1.0 m                     |
| Friction velocity, $u_\tau$   | 0.0079 m/s                |
| Bulk velocity, $U_b$          | 0.1335 m/s                |
| Kinematic Viscosity, $\nu$    | 0.00002 m <sup>2</sup> /s |

Table 2: Physical quantities

In order to compute the bulk mean velocity from the DNS data (Mean stream-wise velocity), trapezoidal rule can be used to solve the equation of  $U_b$  numerically. i.e. *trapz* command in Matlab.

### 5.3.3 Meshing

In order to see the effects of mesh resolution a hierarchy of meshes was created. Three uniformly spaced meshes, in all directions, were generated namely: Mesh 1 (coarse), Mesh 1.5 (medium) and Mesh 2 (fine). The resolution of Mesh 1 was refined by a factor of 1.5 in all directions for Mesh 1.5 and similarly by a factor of 2 for Mesh 2. To compute the mesh spacing the dimensionless distances in wall normal and wall parallel directions (stream-wise and span-wise) were used. The  $u_\tau$  used for spacing computation is taken from the DNS data. Initially the plan was to go for Mesh 3, refinement by factor 4 in all directions, but the total number of mesh points were  $\sim 88$  million ,so the idea was dropped and an intermediate mesh resolution was chosen and hence the name Mesh 1.5.

$$\begin{aligned}
 x^+ &= \frac{\Delta x u_\tau}{\nu} \\
 y^+ &= \frac{\Delta y u_\tau}{\nu} \\
 z^+ &= \frac{\Delta z u_\tau}{\nu}
 \end{aligned} \tag{15}$$

Because of the uniform meshes the spacing is same in all three directions and from now onward it will be referred as  $\Delta^+$ . The bottom and top faces of the domain were defined as the wall, no-slip boundary condition. The no-slip boundary condition was realized by using the half-way bounce back boundary condition. Because of the half-way bounce back boundary condition the distance of the nearest cell to the wall,  $y^+$ , is halved [23]. The  $y^+$  will always be a smaller than  $\Delta^+$  by a factor of 2. The following table 3 shows the mesh statistics used in the investigation, where  $N_x, N_y, N_z$  denote the number of nodes in the stream-wise, wall-normal and span-wise directions. In order to facilitate the grid generator the size of domain was reduced to  $6\delta \times 2\delta \times 3\delta$ .

|          | $N_x \times N_y \times N_z$ | Total nodes | $\sim \Delta^+$ | $\sim y^+$ |
|----------|-----------------------------|-------------|-----------------|------------|
| Mesh 1   | 192 x 64 x 96               | 1179648     | 12              | 6          |
| Mesh 1.5 | 288 x 96 x 144              | 3981312     | 8               | 4          |
| Mesh 2   | 384 x 128 x 192             | 9437184     | 6               | 3          |

Table 3: Mesh statistics

The domain size has been slightly altered in this investigation and this leads to the necessity for performing the two-point correlations to prove the adequacy of the domain. Premnath *et.al* [14] suggests doing it otherwise.

## 5.4 Initial conditions

It is known from the theory of the internal fluid flow that there exists a hydrodynamic entry length ( $L_h$ ), starting from the inlet ( $x = 0$ ), that the fluid has to travel, in the channel, for the flow to become fully developed. This means when we integrate the flow governing equations, in time, the fluid has to travel the distance  $L_h$  ( $\sim x = 10H$ ) in every time-step, to generate a fully developed flow, which is unnecessary considering the computational efforts and also the storage requirements. The idea is to consider only that portion of the channel which will be fully developed, turbulent and neglecting the initial part ( $L_h$ ) of the channel. This implies that specific initial conditions for the flow to become fully developed and turbulent has to be provided.

Different approaches have been used by different authors to generate the initial condition for the incompressible, three-dimensional, turbulent channel flow simulations. Moin *et.al* [9] has specified few consistency criteria for initial velocity field. Sagaut [15] & Fröhlich [4] have a concise description of most commonly used approaches with their respective advantage, disadvantages and their applications to suitable flows. Two of the several approaches mentioned in [15] & [4] will be discussed here.

**i. First approach:** In this approach a simulation is performed for some Reynolds number smaller than the one to be simulated. The resulting fields of this simulations are then used as the starting point for the required Reynolds number and is simulated further in time until a steady state, where the statistics of the flow do not change with time, is achieved. This approach is termed as *Precursor simulation* in several literature.

**ii. Second approach:** In this approach a laminar flow is chosen as the base flow and certain set of disturbances are applied on the base flow. The base flow super-imposed with disturbances is then integrated in time until the flow

transitions to a turbulent flow. This approach is computationally intensive, as the flow takes time to become turbulent. There are several options available to impose the disturbances on to the base flow. Several authors have used this approach for the generating appropriate initial conditions. In this investigation this approach has been used.

For the initialisation of our simulations the approach of *Eugene de Villiers* [13] was followed. *Eugene de Villiers* introduced some sort of near wall streak cycle, where he uses the fundamentals of the turbulent structures near the walls. A parabolic flow profile is chosen as the base flow. To generate the parabolic flow profile the bulk mean velocity,  $U_b = 0.1335 \text{ m/s}$ , which was computed by [13], from the mean velocity DNS data of [8] for  $Re_\tau = 395$ , is used.

Parabolic flow profile has been applied in the stream-wise direction,  $u(y, 0)$ , on all the fluid nodes in the domain. Figure 5 shows the parabolic profile used as the base flow. A combination of sine and cosine waves have been used to break the flow symmetry and to develop the turbulence in the flow. Initially the the

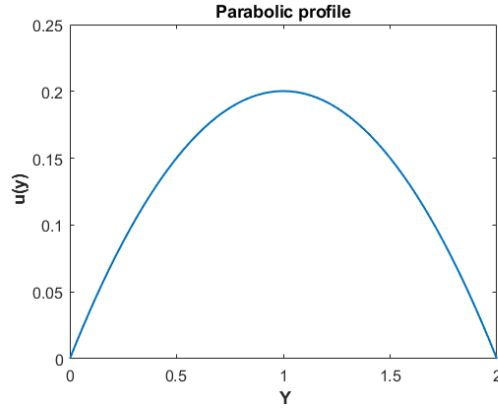


Figure 5: Parabolic flow profile with  $U_b = 0.1335 \text{ m/s}$

disturbances were applied in the stream-wise direction, but that approach was quite time consuming. It took approximately 4.5 million time-steps for the flow to become turbulent. In the next approach the disturbance was applied in the wall-normal direction,  $v$ . This approach reduced the time it took for the flow to become turbulent, 100,000 time-steps. The exact equations for the initialization have been specified in the eq. 18 in Appendix.

## 5.5 Modelling of pressure gradient

In the theoretical section it was shown that the pressure gradient drives the flow through a channel. With the periodic boundary conditions applied in the both streamwise and spanwise direction no information on the pressure gradient is available and thus some kind of external forcing is required to drive the flow through the channel. This requires an additional step, to model the pressure gradient, in the numerical simulation. An additional force term will be added on the right hand side of the Navier-stokes equation. Forcing is only applied in

the streamwise direction,  $F_x$ . Two most commonly used modelling strategies will be discussed here.

**i. Constant Pressure gradient :**

In this approach we set  $F_x$  to some constant value, thus explicitly setting the driving force, which becomes constant and uniform. Since the forcing is constant the mass flow rate will be varying in time, which is attributed to the wall shear stress, until the flow is fully developed (see eq. 4).

$$F_x = -\frac{\partial P}{\partial x} = \frac{\tau_w}{\delta} = \frac{\rho u_\tau^2}{\delta} \quad (16)$$

Thus the simulation a posteriori yields the mean flow rate,  $U_b$ , required to generate the prescribed frictional losses [16]. [17, 10] have used this forcing strategy.

**ii. Constant Flow rate :** In this approach the forcing is not constant, but it is time dependent. The forcing is adjusted in every time step to keep the mass flow rate in the channel equal to the specified flow rate. Flow rate per channel cross-sectional area is  $U_b$ . Thus  $U_b$  is provided as the input and we get the mean friction velocity,  $u_\tau$  as the output which we compare with that of the DNS result. This is the most widely used approach in the turbulent channel simulations [13, 11, 18] and it has been utilised in this investigation. Detailed description of this approach is specified rarely and the authors [16, 19] describe the approach in detail. [16] has shown a comparison between different forcing strategies applied on the turbulent channel flow.

## 5.6 Solver settings

Once the channel flow becomes turbulent the flow is allowed to further develop in time until the statistically steady state is reached i.e. the statistics of the flow variables are invariant to temporal changes. Once the statistically steady state is reached the time averaging of the flow variables is carried out and these time averaged quantities are further averaged over in space along the two homogeneous directions i.e spatial averaging over the x-z planes.

To determine the statistically steady state in the channel one can observe the evolution of the several quantities viz. turbulent kinetic energy (volume averaged),  $\tau_w$ ,  $u_\tau$  etc. over some non-dimensional time. Refer to [20] for the detailed description for steady state determination and spatial averaging. Here two non-dimensional time will be specified *large-eddy turn-over-time (LETOT)*,  $tu_\tau/\delta$ , and the flow through the domain time,  $tU_b/\delta$ . The  $u_\tau$  obtained from DNS data is used for computing the LETOT. LETOT is the time required by the large eddy to breakdown and transfer the energy to the small eddies [12]. In this investigation no specific quantity had been monitored to determine the statistically steady state. An estimation that  $\sim 10$  LETOT is sufficient for the flow to be fully developed and statistically steady [21] had been followed.

The lb simulations are generally performed in *lattice units*, where all the physical parameters are represented by dimensionless numbers. This implies that converting the dimensional quantities into dimensionless lattice quantities and vice versa to interpret the simulation results. Exploiting the law of similarity we can ensure that we simulate the same kind of flow in both the unit systems by maintaining certain non-dimensional parameters equal in both unit systems i.e. Reynolds number.

LB is mostly used for simulating the incompressible fluids where the Mach number is smaller ( $Ma < 0.3$ ). The flow in this investigation is also incompressible. Smaller Mach number  $Ma \ll 1$  implies that the flow in the channel evolves with smaller time-step,  $\Delta t$ , i.e. flow evolves slowly and thus renders the simulation as computationally expensive. In order to avoid that,  $Ma$  has been artificially increased to speed-up the simulation i.e. coarser  $\Delta t$  by keeping the  $Re$  number constant. The computation of the parameters required to increase the  $Ma$  is specified in the Appendix. The table 4 shows the actual,  $Ma_{act}$ , and the artificial,  $Ma_{art}$ , Mach number.

| Mach number | Value       |
|-------------|-------------|
| $Ma_{act}$  | 0.000389213 |
| $Ma_{art}$  | 0.180647487 |

Table 4: Mach number

There are two stages in the simulation of turbulent channel flow: Flow development stage and the averaging stage. In the flow development stage the flow is allowed to develop till it reaches the statistically steady state and in the averaging stage the statistics of the flow variables are sampled. Table 5 shows the time, dimensional and non-dimensional values, required by different mesh resolutions for each stage. As suggested by [21]  $\sim 10$  *LETOT* is sufficient for

|                       | $\sim time$      | Mesh1 | Mesh1.5 | Mesh2 |
|-----------------------|------------------|-------|---------|-------|
| Flow developing stage | $t [s]$          | 7324  | 9806    | 12871 |
|                       | $tu_\tau/\delta$ | 57    | 77      | 101   |
|                       | $tU_b/\delta$    | 977   | 1309    | 1718  |
| Averaging stage       | $t[s]$           | 65917 | 76002   | 69873 |
|                       | $tu_\tau/\delta$ | 520   | 600     | 552   |
|                       | $tU_b/\delta$    | 8800  | 10146   | 9328  |

Table 5: Solver settings

the flow development stage, but the table 5 shows some higher values as it also involves the duration over which the flow becomes turbulent. In the averaging stage the averaging was carried out for longer duration to make sure that we have symmetric Reynolds stress component profiles. This decision was made after observing the profiles generated from every temporal and spatial averaged data set. Mean velocity profiles relatively require less time to have a steady symmetric profile.

## 5.7 Computational details & Post-processing

Simulations have been carried out on the Phoenix cluster [22] of TU Braunschweig. Computations have been performed on the one of the 8 available GPU nodes. Visualisation of the large data sets has been performed using the one of the 6 available visualisation nodes on the cluster. Time averaging of the flow

variables along with the other turbulence statistics is carried out during the on-going simulations within the solver itself i.e. running time average as referred in the literature [9]. The fluctuating quantities are computed in the solver itself. Paraview together with python scripting is used to manipulate and extract the data from the resulting time-averaged data sets. It is also used for the visualisation of the instantaneous data i.e. velocity contour plots, iso-volumes etc. The data extracted by the python scripts is then used by the Matlab scripts to further manipulate and visualise the data in the form of profile plots i.e mean velocity profiles, Reynolds stresses. Spatial averaging was carried out in Matlab and just to point out that it can be time-consuming for the finer mesh resolutions.

## 6 Results and discussion

In this section the results of the simulations carried out will be presented and discussed. Firstly, the results of the LB simulations carried out without any turbulence modelling will be presented and discussed. The result of the simulations with no-model will be termed as the *under resolved DNS (UDNS)*. Under resolved because the resolution is coarser compared to the reference mesh resolution and DNS because no additional modelling is done for turbulence. The DNS results of [6], for  $Re_\tau = 395$ , will be used as a reference for comparison. It will be followed by the discussion with the results from the LB-LES results of WALE model and will be compared with the UDNS and DNS results.

## A Appendices

### A.1 Initial conditions

The following are the initial conditions, for density and the velocity components resp, used for fully developed channel flow simulations:

$$\rho = 27 U_b^2 \left( \cos(x) \frac{4\pi}{x} + \cos(y) \frac{4\pi}{y} \right) \frac{y}{x} + 27 U_b^2 \left( \cos(x) \frac{4\pi}{x} + \cos(z) \frac{4\pi}{z} \right) \frac{z}{x} \quad (17)$$

$$\begin{aligned} u &= 3 * U_b \left( \frac{y}{h} - 0.5 \left( \frac{y^2}{h^2} \right) \right) \\ v &= U_b * \cos(x) \frac{2\pi}{x} * \sin(y) \frac{2\pi}{y} \\ w &= 0 \end{aligned} \quad (18)$$

### A.2 Lattice units

As mentioned earlier the lattice units are the dimensionless lattice values of the corresponding physical parameters. Chapter 7 in [23] provides the detailed explanation on how to non-dimensionalise and several considerations while performing that to the physical units. For the further discussion the lattice units will be marked with an asterisk, \*, sign and the unmarked symbols will represent the physical parameters. As mentioned earlier to simulate the same flow, the Reynolds number in both the unit systems has to be equal. Here the  $Re_b$  is taken as the reference.

$$\begin{aligned} Re^* &= Re_b \\ \frac{U^* l^*}{\nu^*} &= \frac{U_b l}{\nu} \end{aligned} \quad (19)$$

The total number of grid nodes are same in both unit system, thus :

$$\begin{aligned} N^* &= N \\ \frac{l^*}{\Delta x^*} &= \frac{l}{\Delta x} \\ l^* &= \frac{l}{\Delta x} \Delta x^* \end{aligned} \quad (20)$$

The usage of the uniform grid here implies that  $\Delta x^* = 1$ . With the same assumption as of Reynolds number the lattice and the physical Mach number must also match and this implies:

$$\begin{aligned} Ma^* &= Ma \\ \frac{U^*}{c_s^*} &= \frac{U_b}{c_s} \\ U^* &= Ma c_s^* \end{aligned} \quad (21)$$



where  $c_s^*$  is the lattice speed of sound and equal to  $\frac{1}{\sqrt{3}} \approx 0.577$ . When simulating the incompressible flow simulations the Mach number is small and hence it is not necessary to equate the Mach number until and unless it is small. Mach number in lattice units is considered small when  $Ma^* < 0.3$  [23]. Now  $l^*$  and  $U^*$  are known the only free parameter  $\nu^*$  can be computed by using the eq. 19.

As mentioned before the Mach number of the simulation is very small i.e.  $Ma \ll 1$ , as the flow is incompressible. This means that the  $U^*$  will be very small and thus  $\Delta t$  which scales as  $\Delta t \propto Ma^*$  will also be very small. The total simulation time,  $T$ , is expressed as [23]

$$T \propto \frac{1}{\Delta x^n \Delta t}$$

where  $n$  is the spatial dimension. This shows that smaller  $\Delta t$  requires a large number of time-steps for the flow to develop. Hence it is suggested to somehow increase  $Ma^*$  and eventually  $U^*$  as  $U^* \propto Ma^*$ . Let us consider an example of Mesh 1 resolution to show how the Mach number was increased.

$$\begin{aligned} U^* &= \frac{Re_b \nu^*}{l^*} \\ Ma^* &= U^* \sqrt{3} \end{aligned} \tag{22}$$

$l^* = 64$  and  $\nu^*$  is chosen equal to 0.0005 and the resulting  $U^* = 0.104296875$ . With the increase in the mesh resolution the  $\nu^*$  also increases to keep the Reynolds number same. Thus the eq. 22 shows how the  $Ma^*$  has been increased. The intrinsic restrictions of LB algorithm however puts a limit on the simulation parameters i.e.  $Ma^*$  [23]. Thus there is an upper limit till which the  $Ma^*$  can be increased. The table 6 shows the parameters in lattice units to increase the Mach number.

|         | Mesh1       | Mesh1.5     | Mesh2       |
|---------|-------------|-------------|-------------|
| $U^*$   | 0.104296875 | 0.104296875 | 0.104296875 |
| $\nu^*$ | 0.0005      | 0.00075     | 0.001       |
| $l^*$   | 64          | 96          | 128         |

Table 6: Parameters in lattice units

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