

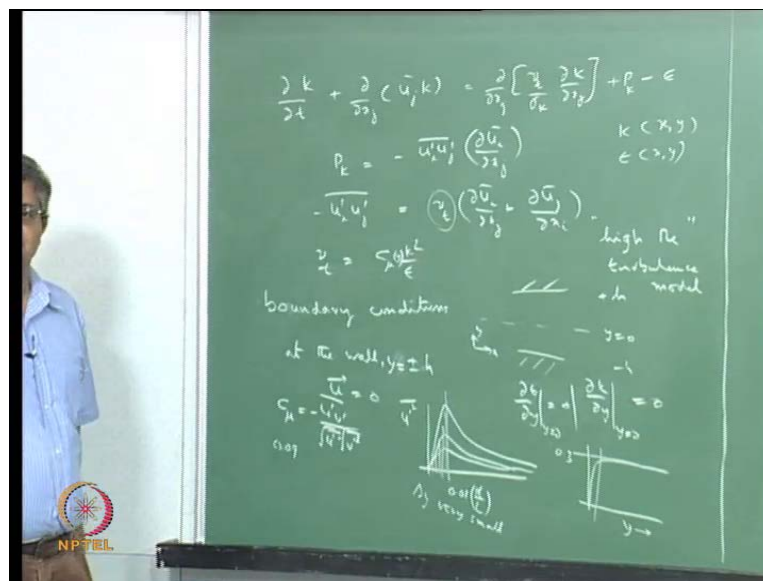
Computational Fluid Dynamics
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Module No. # 06
Dealing with complexity of physics of the flow domain
Lecture No. # 39

We have seen the two equations model for turbulence, in terms of two quantities, k the turbulent kinetic energy, and epsilon, the rate of dissipation of the turbulent kinetic energy being used to evaluate the turbulent viscosity or the eddy viscosity associated with turbulent flows. It is very important in turbulence modeling to consider the near wall turbulence, because that is a very important region, in terms of the production of turbulence and so on.

So, we have to treat the near wall turbulence properly, so as to get solution which is appropriate.

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So, what we mean by near wall turbulence is, that in our model we have written equations like; $\frac{\partial k}{\partial t} + \frac{\partial (\bar{u}_j k)}{\partial x_j} = \frac{\partial}{\partial x_j} \left[\frac{\nu_t}{\sigma_k} \frac{\partial k}{\partial x_j} \right] + P_k - \epsilon$, mean this equal to $\frac{\partial}{\partial x_j} (\nu_t \frac{\partial k}{\partial x_j}) + P_k - \epsilon$

k , $\frac{dk}{dx_j}$ representing the turbulent diffusivity of kinetic energy, plus a production term and rate of dissipation term, this is the epsilon and p_k here, it is a production term which is $-\overline{u_i' u_j'}$, in terms like this.

So where you have large mean velocity gradients, we have a production of turbulence, which goes in to this as a source term, and this itself is represented in terms of ν_t , $\overline{u_i} \frac{d\overline{u_j}}{dx_i}$ plus $\overline{u_j} \frac{d\overline{u_i}}{dx_i}$, where ν_t is given as $C_\mu \frac{k^2}{\epsilon}$. We have a similar transport equation for epsilon and therefore, we have two standard equations of the generic scalar transport equation representing k and epsilon, which go together to determine the turbulent viscosity, which goes on to determine the turbulent or Reynolds stresses, in terms of mean velocity gradients and the turbulent viscosity.

Now, what about the boundary conditions, we know that any partial differentiation equation, any differential equation cannot do without the boundary conditions, and initial conditions as appropriate. Since, we are talking about mean quantities, k and epsilon here are mean time average quantities, so wherever you have a standard boundary conditions, like asymmetry plane, plane of symmetry, and those kinds of things are still applicable.

So, when you have plane of symmetry, like flow through a duct, which is exactly identical with respect to the center plane, then we can calculate the flow field, either in this part or in this part, it is not necessary to go through this, in which case, along this plane **once can say that**, if you say that, this is x and this is y , and this is $y = 0$, and $y = \pm h$; like that, one can say that $\frac{dk}{dy}$ at $y = 0$ is 0, and similarly, one can say that $\frac{d\epsilon}{dy}$ at $y = 0$ is equal to 0.

So, these kinds of boundary conditions natural to the flow are quite valid, but what is important in many practical cases, is also the boundary conditions at the wall, and that is where we have to make a special distinctive characteristics of turbulent flow. Ideally, one would say that, at the wall, in this case, that is at $y = \pm h$. We know that the velocities has to be 0, the velocity has to be 0 by the no slip condition, in laminar flow.

Even in turbulent flow this no slip condition is valid, but we cannot use this boundary condition with this kind of turbulence model, because this kind of turbulence model assumes that, the turbulence is fully developed, and it is strong enough that most of the

diffusion and the production terms, these are uninfluenced by the molecular properties. For example, we have said that, τ_{ij} these Reynolds stresses, are functions only of ν_t which is of a function only of k and ϵ , so this is purely a turbulent quantity, and therefore, this model is appropriate only for high Reynolds numbers

So, it is a high Reynolds number turbulence model, and so I would expect the Reynolds number of the flow to be of the order of 20,000 or 50,000 or greater, for this flow to be valid, and also very close to the wall, we have the special feature that the turbulence is damped.

We have seen the typical plots of, for example u_{rms} , that is $\overline{u'^2}$, it would actually go something like this, and then it would come out like this, and this maximum is very close to the wall. It may be point 0.1 of the non-dimensional length. So, that is let us say that distance divided by correct length, it can be something like that

So within that, you have a very steep variation of typical turbulence quantities, and even the other things will go like this, and towards, and they all may become, as you move away from the wall, you may have isotropy, but very close to the wall, we have severe non isotropy. And, in this model we are not considering the non isotropy of the turbulence, because we have an isotropic eddy viscosity, which is being and brought in here, and it is a function of only the scalar quantities k and ϵ , and there is no possibility of introducing anisotropy, that is very strong, very close to the wall, as you move other from the wall, then you have more of isotropy. So, that is one reason, why we cannot make use of this model very close to the wall.

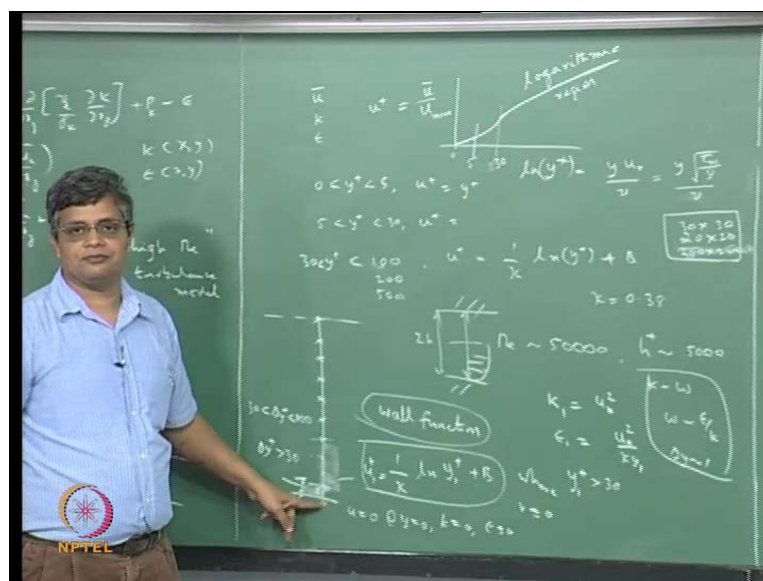
You also have significant damping of turbulence very close to the wall, and finally, we have said that C_{μ} is a constant, and we have said that this is something like, $\overline{u'v'}$ by $\overline{u'^2} \overline{v'^2}$, square roots. So, it is a non-dimensional variation of $\overline{u'v'}$ like this, with a minus sign, and we have said that this quantity is mostly constant, like this and this has a value of a point 3 with respect to distance from the wall, distance from the wall increasing, this reaches a constant value, based on this we have put C_{μ} value of 0.09. We have made this a constant, and it is not a function of a distance from the wall, but very close to the wall this changes drastically, again like here.

So, that means that when you come very close to the wall, the turbulent viscosity is no longer directly proportional to k and ϵ , but this constant of functionality itself is a function of y , the distance to the wall. And, also the k and ϵ although they are 0 at the wall, they show a very steep gradient very close to the wall, in the region very near the wall.

So, that means that if you want to resolve a very steep gradient going like this, you need to have fine Δy here, so Δy needs to be very small. In order to resolve the very rapid variation of the turbulence quantities both k and ϵ in the region close to the wall, and that means that the number of grid points, which are needed to discretise the given cross section will become very large.

So, there are number of reasons, one is the non constancy of the constants, the non isotropy of the turbulent viscosity, and the need for very small Δy ; for all these reasons we cannot take this model very close to the wall, so we start calculating variation of k and ϵ . For example, k as a function of x and y and ϵ as a function of x and y , are calculated only from after a small distance from the wall, within the small distance the turbulence quantities change very rapidly, but fortunately that change can be expressed in terms of local variables or what are known as inner variables, that means that the variation here

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The variation of the mean velocity u and ϵ are very close to the wall, can be expressed in terms of dimensionless quantities y^+ , which is defined as $y^+ = u^* y / \nu$, where u^* is the friction velocity, and that is given as $u^* = \sqrt{\tau_w / \rho}$, this is square root of wall shear stress divided by the kinematic viscosity of the fluid. So, we are bringing in, the kinematic viscosity of the fluid here, and we are non-dimensionalizing with respect to this friction velocity, which is a function of this, and now if you have to plot, this mean velocity, which is again non-dimensionalized with respect to for example, u_{max} or u_{mean} as in the average velocity in a pipe like that.

Then, it shows a characteristic variation, it goes through like this, and then becomes a fairly constant thing on long plot, where y^+ is plotted like this, and within a distance of say over a distance of y^+ from 0 to 5, you have a linear variation between u^+ and y^+ , so this we call as u^+ . So, we know that $0 < y^+ < 5$, this u^+ varies linearly with this, and between 5 and y^+ of thirty or so, u^+ is a fairly strong function of y^+ , it is a non-linear expression and for y^+ greater than 30, and so we can put it like this. $u^+ = \kappa \ln(y^+ / b)$, where κ is a constant. We are familiar it is around point 4.38. 4.142 like that, so this has a slope of about 2.5 when plotted as u^+ versus $\ln y^+$.

So, this is that this is known as the logarithmic region, and this region here, where your u^+ is y^+ is known as the viscous sub layer, and this is a buffer layer, so through this there is a more complicated variation here, but once you reach a distance of y^+ greater than of around 30, you have almost a linear variation with $\ln y^+$, and this is known to be valid, this logarithmic thing is valid for, it is a very universal velocity profile, this particular variation. It is universal, because it is valid for flows with pressure gradient and with adverse pressure gradient, and with favorable pressure gradient for boundary layer flows, for flows inside pipes, and for all these cases, this sort of thing is valid up to a certain distance.

So, in this sense very and this y^+ of 30 is typically very small, because the maximum y^+ , for example if you are looking at flow, between 2 plates and at a Reynolds number

of the order, say 50,000, then if you say that this is $2h$, then h plus, which corresponds to the maximum distance, is of the order of 5000 or something like that.

So, this 30 which is the beginning of the logarithmic here is a small fraction of this whole thing, and we take advantage of this to say that, we know the variation of the velocity, parallel to the wall, so in this direction you have a velocity profile like this, we know this variation in the region very close to the wall, and it is related to the local shear stress in this particular way. So, we put, if this is our overall domain, this is the center and here, and this is wall here; so we put a first grid point here, second point, third point like this, and instead of saying that u is equal to 0, at y equal to 0, which is the no slip boundary condition. We say that at this point, say u_1 is given by $1/\kappa \ln(y_1/b)$. So, the boundary condition that we apply here, and this is u_1 plus.

So, this u_1 plus is the mean velocity at that point divided by the average velocity or the maximum velocity like that. So, in that sense, we specify the values of the velocity at the first grid point, taking advantage of the fact that within this point here, where in such a way y_1 plus is greater than 30, so we choose the Δy here.

We choose the Δy here, such that Δy plus is greater than 30, or let us say it is between 30 and 100 or 200 like this. So, we choose a value which is not very very close to the wall, but it is at a sufficiently long distance from the wall, such that it falls in the logarithmic region, because in the logarithmic region we have an expression simple expression for u plus as a function of y plus.

So, if you know the distance when using this we can get the non dimensional velocity here, so this is the boundary condition that we use, so this is known as a wall function, because close to the wall, we are making use of an algebraic function, u in terms of y here to represent the velocity, it is not given by the momentum equation, nor is it given by the real boundary condition that at y equal to 0, u equal to 0

So, we are saying that very close to the wall, we have a velocity which is non zero and it is given by this universal velocity profile and the velocity normal to the wall, v is equal to 0 and w is equal to 0, is the standard boundary condition that we would get. What about k and ϵ , even k and ϵ are given by a similar functional values, one can show that k_1 , so that is the turbulent kinetic energy at the first point, provided the first point is between 30 and 100 in terms of y_1 plus is equal to u^*^2 , where u^*

square is given in terms of this and one can show something like this, and ϵ at this point is given in terms of u^2 by κy^2 , something where κ is point 38 and y is the distance of the grid point from the wall. So, in that sense, we specify the values of u , v , w and k and ϵ at the first grid point, locating the first grid point at a sufficient distance from the wall, such that the y lies between 30 and 100.

The reason for this is that, we do not want to get too close to the y , because then we have steep variation, slightly more complicated expression for this, and if you put too much, and if you take a y plus and may be 200 500 or 1000 in to that, then this first grid point will be somewhere here, and in all this region, you are not actually solving the momentum equation, you are making use of a mole function. So, in that sense, we want to make sure that the region, where we make use of where we impose the functional variation is very thin compared to the overall floor domain, we would like to make use of as much as the floor domain to be computed by the momentum equations, and it is only very close to the wall where we have deficiencies with respect to the way that modeling turbulence here, we make use of the wall functions.

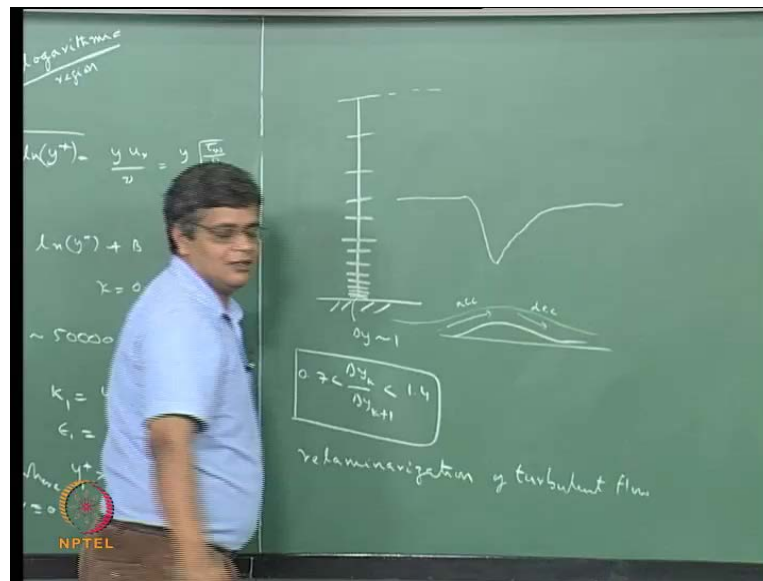
So, this approach enables us to use, for example if you have a duct of a cross section like this, we can make use of the grid which may be a 20 by 20, grid to represent the flow through this, or may be 30 by 30, but if we do not have this wall function, and if we want to resolve this velocity profile all the way through this, right up to this, then we have to make changes to this turbulence model, there are definitely those kind of turbulence models which enable us to calculate right down to y equal to 0, for which the boundary condition will be u equal to 0 at y equal to 0 and k equal to 0 and ϵ equal to 0.

So, you have the natural boundary conditions, for such models, those turbulence models which make modifications to the standard high Reynolds turbulence model in such a way that you take account of the (ω) trophy, that you take account of the damping of turbulence close to the wall and all that. So, there are such kinds of low Reynolds number turbulence models, and there are so models which like k , ω type of model, where ω is related to ϵ by k , so instead of solving for k and ϵ like what we are solving here, we solve for k and ω and we get ϵ from this. So, these kind of approaches are more amenable to go right down to the wall, so using those models we can resolve the full layer, but in such a case we cannot use only this

kind of grid, we may have to have Δy for these kind of models, Δy close to the wall should be of the order of one.

So, that means that if you have a flow, such that the overall the maximum h plus is the order of 5000, then your grid Δy here should start with one here, and then if you have uniform spacing, you have 5000 grid points, so it should be 5000 by 5000, you probably do not need that, but overall you'd have in such a case.

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A grid which is geometrically expanding, this is the center line, and here you'll have a grid spacing which can increase in geometric progression, but with the restriction that Δy close to the wall is of the order of 1, and also from, in order not to bring in numerical accuracies associated with a non uniform grid, you would like Δy_k divided by Δy_{k+1} plus 1

So, that is, this Δy_k divided by this Δy_{k+1} , so that sort of things, successive values to lie between point 7 and roughly 1.4 is a general thumb rule, so as not to lose too much accuracy of discretization or approximation, finite difference approximations when we have this sort of expanding grid.

So, because of this you may not have to use 5000 or 5000 but, definitely the number of grids to represent something like a square duct at a Reynolds number of 50,000 like this, may increase up to 200 by 200, and that means it is a hundred fold increase in the

number of grid points, and the computational time will also increase much more, may be a factor of 100 squarelike that.

So, that is why in the general case, we would try to use the wall functions in order to represent the turbulence variation very close to the wall, but there are special cases where we may want to these kinds of models which enable us to compute the flow variables all the way to y equal to 0.

For example, if you have a case of a strongly accelerating flows, in such a case it is possible that the flow near the wall may become laminar, so that is called relaminarization of a turbulent flow, and one example is that, let us say that you have a wall here, and then you have a protuberance like this, so the flow is supposed to go through over this and come out, there's a small hill associated with this, and in this you have an accelerating region and then you have a decelerating region. So, this acceleration of the flow, if it is strong enough, it is supposed to damp out the turbulence, and if the acceleration is sufficiently strong and sustained, the flow the flow may become laminar as it goes over the hill, and once it becomes laminar, the turbulent viscosity is 0.

So, that means, that your friction factor will be very different, and the heat transfer coefficient will be different and all that. So, if you were to plot, for example the Nusselt number, it may be something like this value here but, as you go over the hill, it may come down and then it may go back like this.

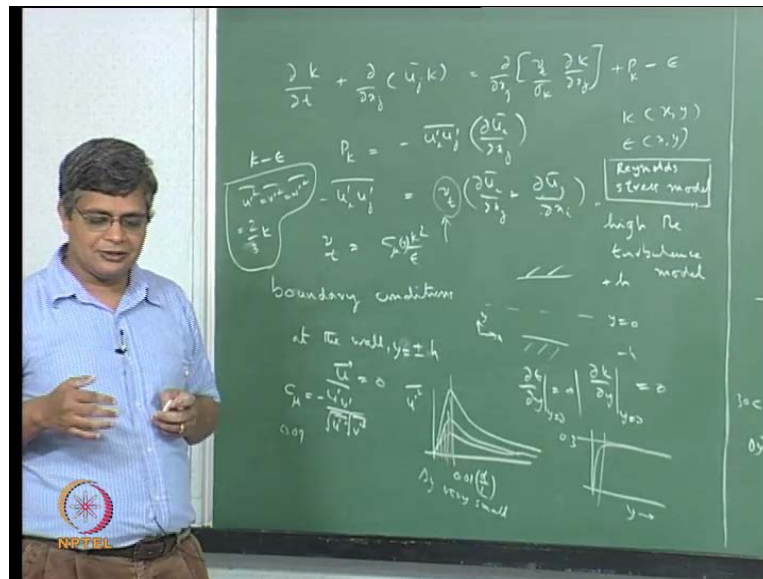
So, this drop in the heat transfer coefficient the Nusselt number is associated with the relaminarization of the flow, and if you want to capture those kind of effects, then you have to resolve the turbulence right up to the wall, so in such a case you may have to use low Reynolds version of this, these type models or the k ω type of models like this, to get these special effects.

Otherwise we have this kind of fairly universal model, the k ϵ model, which can be used as a first to investigate to calculate the turbulent flow without having to fix any constants, but even though this is considered as a fairly general model, there are modifications to this, for example strongly accelerating force for buoyant flows, and so in such cases people have come up with different set of constants for this, and so if we know such information, if you have such information then we can make use of those

ad hoc fixes to the constants that going to the determination of k and epsilon as a function of x y z and then the corresponding determination of the turbulent viscosity and so on.

Now, in that sense k epsilon model can be considered as a good enough, completely specified, self consistent and a model which we can use straight away without having to give ad hoc fixes, like we have to do for the case of mixing length model, but this has some deficiencies, one deficiency is the near wall treatment that we have seen, and that is that we have to make use of wall functions, and that we cannot make use of the natural boundary conditions that at the wall we have k equal to 0 and epsilon equal to 0, but we have to assume a certain variation or we have to impose certain variation close to the wall, some non dimensional variation close to the wall in order to get a solution.

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Another problem with this, is the assumption of isotropic viscosity here, the turbulence stresses here are proportional to the mean velocity gradients through an isotropic viscosity, isotropic viscosity, because this viscosity is independent of which of the 6 stress that we are considering here, and this is isotropic, because this is determined only by k and epsilon, which are again scalar quantities, which are therefore, in that sense isotropic.

So, this isotropic viscosity is known to be not so good when you are considering, for example a swirling flow, where you have flow going like this or within a flow with a strong curvature, you have flow going and then turning, so you have flow in a bend or

flow in a coiled tube or flows with strong buoyancy. All, this are cases in which, the additional force field arising out of this special effect, the swirling nature and the buoyancy, these are directional, these are these are aligned only in certain components of the velocity.

So, you can say that these bring in source terms, additional terms only in certain momentum equations, and not in others other equations, so in such a case one can expect the different turbulence stresses to respond in different ways to the imposed external force field like buoyancy.

So, this assumption of isotropic viscosity, it will give us problem in such cases, and then this model is also unable to represent the different levels of u' prime square v' prime square and w' prime square at the wall, because once we know ν there, and once we know these velocities, one can calculate the different stresses here, and for this kind of plane flow in which only du by dy is present, so that this is not present, we find that the turbulence stresses are considered to be the same, and one would say that u' prime square is equal to v' prime square is equal to w' prime square is equal to two thirds of k , is the only thing that we can say about the level of these turbulence quantities fluctuating components in each of the directions.

So, this equivalence of these things, is implicit in this model, whereas experiments show that these can vary in different ways, and there are also experiments, which show that, this linear relation between the turbulence stress and the strain rate, in which we have made use of the, for example the similar linear relation between the viscous stress, and the strain rate in a fluid, if this linearity assumption is also not correct.

So, there are certain experimental data, which can be obtained, which can be explained only by a non-linear variation between the turbulence stresses, and the strain rates, and so we have also models non-linear two equation models, and to represent the general anisotropy and responsive different turbulences from arbitrary force field, we have the so called Reynolds stress model.

I am just writing in this box, in which separate transport equation is solved for each of the 6 Reynolds stresses, and in such a case you have the 4 equations representing, the time averaged continuity equation, and 3 time averaged momentum equations, and then there are 6 Reynolds stress components, so you have 6 additional transport equations for the 6

Reynolds stresses and you still have an equation for ϵ , that is dissipation rate which appears in the Reynolds stress equations.

So, you have total of eleven equations to be solved for a turbulent flow calculation, not just four, as we do in laminar flow. In the case of k-epsilon model, we have to solve 4 time averaged momentum equations in the continuity equation, and then 2 equations for k and epsilon, we have 6 equations.

So, by increasing the number of equations, one can bring in more and more sophistication into calculation, and thereby we can become more accurate in our calculations, especially for certain cases where we have sufficient data to determine fully the constants that appear in the Reynolds stress equations, but there is also the side effect of trying to solve more and more turbulent equations, because the side effect is, that the number of equations that we want to solve increases from 4 in laminar flow to 6 in the two equation model to 11 in the case of Reynolds stress model, this increases the computational time and it also increases the coupling among all these equations, say therefore, the convergence of these coupled equations it becomes compromised so that overall solution becomes less robust.

So, that means that, if you try to solve all the equations, you may not be able to get an acceptable solution, sometimes you may get into non convergence, divergence of the ϵ scheme like that. So, there is a loss of robustness as you go to, as you move away from the standard k-epsilon model, towards Reynolds stress model or towards non-linear models and all that. So, we have to make a balance between the accuracy that we want, and the kind of numerical effects that we are willing to spend in getting a solution.

If there is strong reason to believe that k-epsilon model is not correct for a particular flow application, only then we should venture out into the special effects, either for better treatment of the turbulence close to the wall, in terms of the k-omega type of models, which require more number of grid points or the Reynolds stress type of models which do not depend on the isotropic nature of the turbulence viscosity as modeled in this or the non-linear models which also do not make this relation linear. So, all those special effects should be considered in, if we know, if we anticipate a problem with the k-epsilon model.

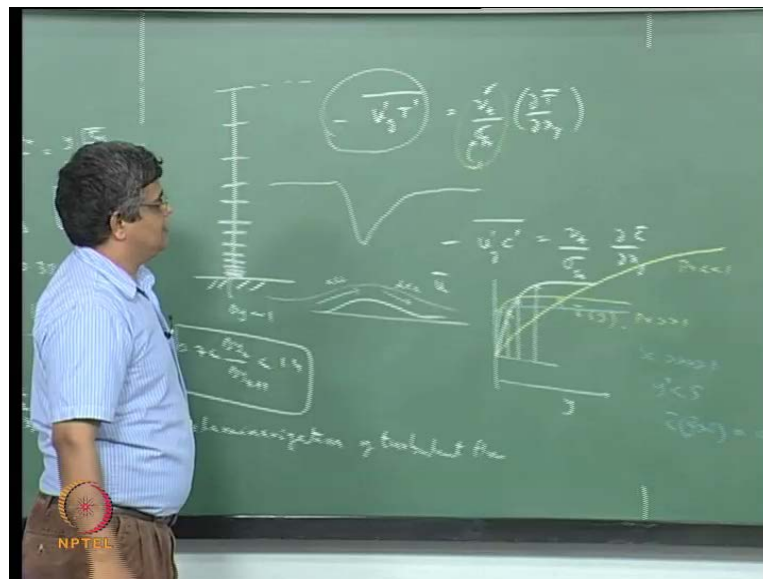
Now, we have said enough about the flow equations to especially the solution of the Navier-Stokes equations in case of turbulent flow, but typically we not only have the flow,

but we also have the heat transfer, the mass transfer and the chemical reactions associated with the flow. What about those things, how are those equations modified, when we consider turbulent flow.

The simple answer is that, the approach that we have used for the Navier-Stokes equations, for the momentum equations in terms of Reynolds averaging, and then trying to come up with a simplified form of an equation, representing the turbulence effect on the overall time averaged variation of the quantity.

So, that approach is extendable readily to the energy equation for non-isothermal flows, and to the species conservation equation for the mass transfer and also chemical reactions. So, we can come up with time averaged energy equation, which look very similar and in such case instead of having a term like $u_i' u_j'$, $u_j' T'$.

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We will have a term like $u_j' T'$, minus $u_j' T'$ as the equivalent fluctuating energy flux, temperature flux, time c_v will give us the energy flux, associated with turbulence and this is modeled in very similar way, like this is represented as ν_t by something like the turbulent Prandtl number, we will just put it as Prandtl number here, subscript Prandtl number, times $\frac{d\bar{u}_j}{dx_j}$, $\frac{d\bar{u}_j}{dx_j}$.

So, this is how we can do, this is already calculated as part of the solution of the momentum equations, and this value here is roughly the order of 1, \bar{t} is calculated from the time averaged energy equation. So, this is one way of closing this particular equation.

So, this is the additional term, that appears in the time average energy equation, and that can be solved like this. Similarly, when we are looking at the species conservation equation, when they come up with u_j' , say c' , where c represents concentration species, and this again can be represented as u_j by σ Schmidt number or something like that, representing the mass transfer, $\sigma = \frac{\rho \bar{c}}{\rho' c'}$.

So, in that sense, when we time average the corresponding conservation equation, we do get into terms which are similar to the turbulence stresses that we have obtained by time averaging the Navier-Stokes equations. We can extend a similar kind of modeling of this turbulence fluxes in terms of the turbulent viscosity, and turbulent Prandtl number, turbulent Schmidt number and the mean gradients of the temperature, and of the concentration gradient like this, and thereby come up with time averaged scalar transport equations for the energy. And, when we come to the wall variation, variation of the temperature and concentration flows to the wall, we have to come up with equivalent wall functions. As long as, the Prandtl number of the fluid, and the Schmidt number of the fluid for the energy flux, and the mass flux are around 1, we can make use of very similar kind of wall functions for this, but typically when we consider heat transfer, the Prandtl number of the fluid need not be equal to 1, for many gases it is of the order of point 7, so it is fairly close to 1, but if you consider some organic liquids, it can be very high, of the order of 100, and if you consider liquid metals, Prandtl number can be very low of the order of point 0.1.

So, you can have wide variation of the Prandtl number, and in such cases when Prandtl number deviates significantly from 1, you have to be very careful about the correctness of the wall functions that we use to represent the turbulent fluxes close to the wall.

Similarly, when we consider mass transfer, we have the Schmidt number which represents mass diffusivity to the momentum diffusivity, so that can be very large and for some organic fluids it can be of the order of 10000 even 50000 like that.

So, again in such a case you have a variation of the concentration profile which is very different from the variation of the velocity profile, in turbulent flow, and especially when the Prandtl number is much much greater than 1, or when the Schmidt number is much much greater than 1, you have the strange case that, if you are looking at a velocity profile which is going like this.

So, this is the velocity, and this is the distance from the wall, so this is y here, and this is the average velocity, so this is the typical let us say that this is the turbulent typical turbulent velocity profile, the Prandtl number variation can be like this, it is most of the variation of the temperature, so this is \bar{u} as a function of y . It can be like this, if the Prandtl number is much much greater than 1, and it can be like this, if Prandtl number is much much less than 1, and what does it mean here, your first grid point in the wall functions may be such that, your first grid point is here, this is your this is the region in which you have assumed a certain variation of velocity, and you see that, for when Prandtl number is much much greater than 1, most of the variation in the temperature, mean temperature profile is already over within that region.

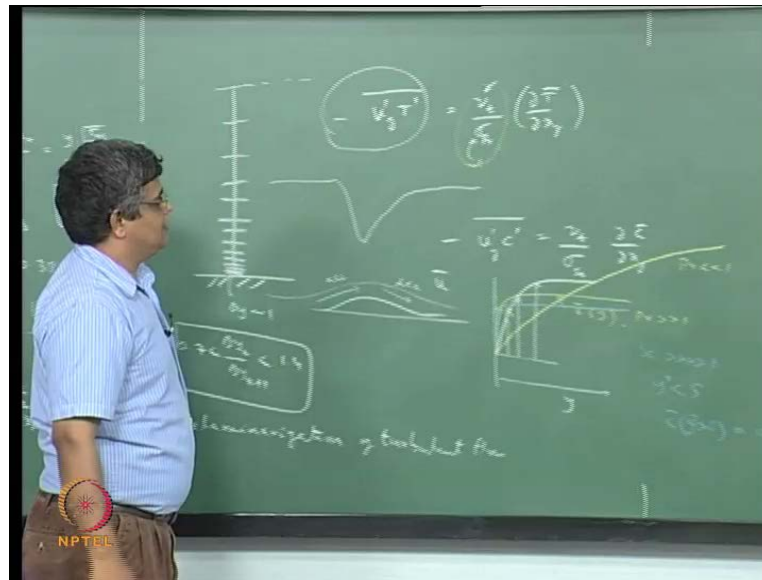
So, you are imposing an uncalculated turbulence profile on the heat transfer here, and there may be some discrepancy coming from that, and when you are considering this, when you use a wall function which is standard wall function, that again may be quite wrong here.

If we consider a Schmidt number, which can be much much higher than 1, so Schmidt number much much greater than 1, let us say a 10,000, then the variation of the concentration profile, it may be like this, that means that typically for these values, the even for y^+ , within y^+ of greater than 1 itself, that is within the viscous sub layer itself. So, mean concentration for y^+ greater than y^+ greater than 5, it may become constant.

So, the variation of the concentration with y will be within the viscous sub layer, and what is importance of that, generally when we talk about viscous sub layer here, we say that k is equal 0, that is, it is purely viscous dominated thing, there is no turbulence within that viscous sub layer. So, there's a general thing that we normally ascribe doing this.

So, that means that in the case of turbulent diffusivity of the species for a Schmidt number, which is much much greater than 1. In reality there may be some effect of turbulence, but you are imposing no effect of turbulence here. So, this is something that we have to consider here.

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So, as per your wall function which is somewhere here, there is no further variation of concentration profile within this region. So, this leads to some discrepancy, one has to be careful about evaluation of the temperature profiles and the concentration profiles for fluids which have a Prandtl number or a Schmidt number much greater than 1.

One has to pay special attention to the near wall treatment of these quantities in turbulent flows. So, if necessary one may have to go to models like this, which can take us right up to y equal to 0; in which case, they can bring in faithfully, the effect of molecular viscosity and molecular diffusivity in the flux of mass flux, species mass flux and energy flux close to the wall. So, that is one thing or we have to alter the standard wall functions that we use for the energy and species concentration.

So, we have to be careful, otherwise we cannot expect to get good agreement between the measured heat transfer coefficients or measured mass transfer coefficients and the computed completed one, all not because of the discretization errors, and all computation errors, because you have made wrong assumptions about the influence of turbulence close to the wall for these cases. So, there you have wall dominated heat transfer, wall

dominated mass transfer, it is important to look carefully at the near wall treatment of the turbulent diffusion of these quantities, in case of turbulent flow.

In case of laminar flow, we do not have any such problem, but turbulent flow we have problem, because we are solving only the approximate form of the equations with assumed model for the turbulent transport of these quantities. So, if this is accurate, if this is appropriate, we get good profiles for temperature and concentration. If, this is not accurate, then we can get wrong values of the profiles, and therefore wrong values of the heat transfer and mass transfer coefficients.

So, this is about turbulence modeling, one can see that in the general case we have to use some sort of Reynolds time averaging of the governing equations, so that we do not have to worry about very fine turbulence fluctuations, temporal fluctuations, and spatial fluctuations, but in the process we are coming up with additional unknown quantities like the Reynolds stresses, and these are only approximately modeled in using a vast array of models like this, and therefore, the equations that we are solving for turbulent flow are no longer, fundamental equations which have to be ϵ , these are only approximate form of equations.

So therefore, we have to be varying of the errors introduced in modeling, when we are looking for tumbling flows, and there is an array of models from which, we can choose that model which is appropriate for our flow. Although, the standard k epsilon model has proved to a robust model, it may not always give the correct values, and if so we have to anticipate what kind of problems we have, what kind of phenomenon we are looking at, flow phenomenon we are looking at, and choose to either go for a Reynolds stress model, in case we are looking at strong curvature effects, strong buoyancy effects playing a part in this or if you are looking at strong wall influence, we may have to go for these type of models or we may have to take careful look at the wall functions that we are using.

So, one has to judiciously select turbulence model, so that we can solve for the variables of interest, that is velocities pressure, temperature and concentration and so on.

Finally, when we are looking at chemical reactions, we have to specify the mean rate of reaction and that is very contentious issue, and there is a vast array of models, when we have a large set of chemical reactions which are happening together, because usually

this chemical reaction rates are dependent exponentially on the temperature, and temperature fluctuations may cause a wide range in terms of the chemical reaction rates.

So, there is an entirely different set of models developed for that, one has to consider those things, before one attempts a serious calculation of turbulent reacting flows. So, we have to be informed to make good choices and I hope this particular discussion has helped in highlighting some of the aspects.