HEAT TRANSFER TO THE TRANSPIRED TURBULENT BOUNDARY LAYER

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Report HMT-14

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Abstract

This paper contains a summarization of five years work on an investigation on heat transfer to the transpired turbulent boundary layer. The experimental results are presented for friction coefficient and Stanton number over a wide range of blowing and suction for the case of constant free-stream velocity, holding constant certain blowing parameters. The problem of the accelerated turbulent boundary layer with transpiration is considered, experimental data are presented and discussed, and theoretical models for solution of the momentum equation under these conditions are presented. Data on turbulent Prandtl number are presented so that solutions to the energy equation may be obtained. Some examples of boundary layer heat transfer and friction coefficient are presented using one of the models discussed, employing a finite difference solution method.

Nomenclature

<u>English letter symbols:</u>

A^+	constant in the Van Driest Hypothesis, Eqs. (18) and (20)
A_o^+	value of A^+ for the case of no transpiration and no pressure gradient
B^+	constant in the Evans hypothesis, Eq. (22)
B_o^+	value of B^+ for the case of no transpiration and no pressure gradient
B_f	a friction blowing parameter, $(v_o/U_\infty)/(C_f/2)$
b_f	a friction blowing parameter, $(v_o/U_\infty)/(C_{f_o}/2)$
B_h	a heat transfer blowing parameter, $(v_o/U_\infty)/St$
b_h	a heat transfer blowing parameter, $(v_o/U_\infty)/St_o$
C_f	friction coefficient
C_{f_o}	friction coefficient for the case of no transpiration, at the same value of Re_{δ_2}
c_p	specific heat at constant pressure
C	lag constant in Eq. (24)
D	mixing-length damping function
F	blowing or suction fraction, v_w/U_∞ , $\dot{m}''/G_\infty$
G_∞	free-stream mass velocity, $\rho\, U_\infty$
g_c	proportionality factor in Newton's 2nd Law
H	boundary layer shape factor, δ_1/δ_2
h	convection heat transfer coefficient
K	acceleration parameter, $(\nu/U_\infty^2)(dU_\infty/dx)$
k	mixing-length constant for the inner region
ℓ	mixing-length
ℓ_t	turbulence length scale

$\dot{m}"$ mass transfer rate at fluid-surface interface

$\overline{P}$ time-averaged pressure

P^+ pressure gradient parameter, $-K/(C_f/2)^{3/2}$

Pr Prandtl number, $\mu c_p / \gamma$

Pr_t turbulent Prandtl number, ϵ_M / ϵ_H

$\dot{q}"$ heat flux at fluid-surface interface

Re_H enthalpy thickness Reynolds number, $\Delta_2 G_\infty / \mu$

Re_M momentum thickness Reynolds number, $\delta_2 G_\infty / \mu$

Re_t local turbulence Reynolds number, $\ell_t \sqrt{\overline{u'v'}} / \nu$

Re_x x-Reynolds number, $x G_\infty / \mu$

St Stanton number, $h/(G_\infty c_p)$

St_o Stanton number for $F = 0$ at same Re_H

$\overline{t}$ time averaged value of temperature

t_o temperature at fluid-surface interface

t_∞ free-stream temperature

t' fluctuating part of temperature

$\overline{t'v'}$ turbulent heat flux

u' fluctuating part of x-component of velocity

$\overline{u}$ time averaged value of x-component of velocity

U_∞ free-stream velocity

$\overline{u'v'}$ turbulent shear stress

u^+ $\overline{u}/\sqrt{g_c \tau_o / \rho}$

v' fluctuating part of y-component of velocity

$\overline{v}$ time averaged value of y-component of velocity

v_w normal direction velocity at fluid-surface interface

v_w^+ a blowing parameter, $v_w/\sqrt{g_c \tau_o / \rho}$

x distance measured along surface in direction of flow

x^+ $x\sqrt{g_c \tau_o / \rho}/\nu$

y distance measured normal to surface

y^+ $y\sqrt{g_c \tau_o / \rho}/\nu$

y^+_{crit} value of y^+ at edge of viscous sublayer

Greek letter symbols:

α thermal diffusivity, $\gamma/(\rho c_p)$

γ thermal conductivity

δ momentum boundary layer thickness

$\delta_{.99}$ 99 percent thickness of momentum boundary layer

δ_1 displacement thickness, $\int_o^\infty (1 - \bar{u}/U_\infty)\,dy$

δ_2 momentum thickness, $\int_o^\infty (1 - \bar{u}/U_\infty)(\bar{u}/U_\infty)\,dy$

Δ thermal boundary layer thickness

Δ_2 enthalpy thickness of thermal boundary layer,

$\int_o^\infty \{(\bar{u}/U_\infty)(t - t_\infty)/(t_o - t_\infty)\}\,dy$

ε_H eddy diffusivity for heat

ε_M eddy diffusivity for momentum

λ a mixing-length constant for the outer region

μ dynamic viscosity

ν kinematic viscosity, μ/ρ

ρ density

τ total shear stress

τ_o shear stress at wall surface

τ_t turbulent component of total shear stress

τ^+ τ/τ_o

Introduction - Interest in the behavior of the transpired turbulent boundary layer extends back at least to the early 1950's when transpiration was first being investigated as a means of cooling aerodynamic surfaces under high velocity flight conditions. This interest is reflected in some of the pioneer papers and reports of Rubesin [1], and the paper of Dorrance and Dore [2]. While these contributions were on the theoretical side, experimental work began to take place, notably in the Chemical Engineering Department at M.I.T., and is reflected in the papers of Mickley et al [3,4]. The early 1960's saw interest spreading with significant contributions from Great Britain, in particular those of Stevenson [5], McQuaid [6], Black and Sarnecki [7], and contributions from the Soviet Union, for example Romanenko and Karchenko [8], and Kutateladse and Leont'ev [9]. During the past five years the number of contributions has grown into dozens. A complete bibliography, especially since 1960, is outside of the scope and purposes of the present paper, and the above citations are simply noted to provide some idea of interest intensity as a function of time.

Transpiration from a solid surface over which a fluid is flowing, and on which a boundary layer is developing, is of interest in a number of quite different types of applications. In the typical transpiration cooling application, the solid surface is constructed from some kind of porous solid material. Cooling fluid, chemically the same as the free-stream, is then forced through the surface with the objective of protecting the surface from a hot free-stream. This is a boundary layer problem for which the normal component of velocity at the solid-fluid interface is non-zero, but otherwise the same momentum and energy boundary layer differential equations must be solved as for the

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non-transpired boundary layer. A variation on this problem occurs when the cooling fluid is a chemically different specie than the free-stream fluid. For example, helium might be injected as a coolant to protect a surface from a high temperature air free-stream. In this case the mass-diffusion equation of the boundary layer must be solved in addition to the momentum and energy equations. There are obviously similarities between these two types of problems, but also fundamental differences. Both are "mass transfer" problems in the sense that mass is transferred across the fluid-solid interface, but the latter is also a mass diffusion problem, while the former is not.

Another transpiration problem arises when there is evaporation or sublimation from an interface into a boundary layer, or condensation onto the interface. A further variation on the problem arises when there is chemical reaction either within the boundary layer or at the surface.

In any of the cases cited, the direction of the flow normal to the surface at the interface could be into the surface, or it could be out of the surface. The terms "blowing" and "suction" are frequently used to denote the direction of flow at the interface, while the word "transpiration" generally is taken to embrace both cases. Suction is sometimes used as a scheme for aerodynamic boundary layer control because it is possible to inhibit or prevent boundary layer separation by suction.

These various types of applications suggest why chemical, mechanical and aeronautical engineers have all made significant contributions to the theory, and the terminology to a certain extent reflects these various origins of interest.

In 1965, the author and his colleagues embarked on a comprehensive experimental investigation of heat transfer to the transpired turbulent boundary layer with the objectives of resolving some of the contradictions in the earlier data, filling in obvious holes, and ultimately of obtaining necessary information for a more fundamental theory so that adequate boundary layer predictions can be made over a wide variety of operating conditions. To date, this work has been entirely concerned with the single chemical component problem (i.e., no mass diffusion) using air as the working medium, and employing small temperature differences for the heat transfer work, so that it is effectively the constant property boundary layer that is considered. At the present time, the technical application that provides the major impetus for this program is the prospect of cooling gas turbine blades by transpiration. This application involves the transpired boundary layer with large temperature differences, as well as other complications that will be briefly discussed at the end of the paper, but a fundamental understanding of the behavior of the idealized low-velocity, constant property case is essential if any significant progress is going to be made in the more complex applications.

The objective of this paper is to present a summary of some of the more significant results and conclusions that have resulted from this program. Much of what will be presented (although not all) has appeared in various reports and papers, but significant portions of the work have now been completed, and it seems an appropriate time to attempt to summarize.

The general problem considered is perhaps best illustrated by reference to Fig. 1. Air flows at a steady rate along a flat surface which is porous, and through which air can be forced into the boundary layer or withdrawn from

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the boundary layer (blowing or suction). It is presumed that the surface is aerodynamically smooth, which means that the transpiration holes are small relative to the thickness of the laminar sublayer of the boundary layer, and that their spacing is small relative to the thickness of the laminar sublayer. It is presumed that the solid surface is a heat conductor, and that heat can be conducted to or from the surface. It is further presumed that the surface construction is such that the transpired fluid is in thermal equilibrium with the solid surface at the interface. In general, we would like to consider the case where the mass transfer rate $\dot{m}''$ (i.e., the transpiration rate) is any arbitrary function of distance x along the surface, where the convection heat transfer rate $\dot{q}''$ is any arbitrary function of distance x along the surface, and/or the surface temperature t_o is any arbitrary function of x. We would like to consider the case where the free-stream velocity U_∞ may vary in any arbitrary manner with x, but we will restrict the free-stream temperature t_∞ to a constant. The boundary layer is two-dimensional, with the coordinate y being used to measure the distance normal to the surface. Under these conditions, we are interested in the development of a momentum boundary layer, characterized by a thickness δ, and a thermal boundary layer characterized by a thickness Δ.

It should be added that the problem described, but with the boundary layer laminar rather than turbulent, is of equal interest in many applications including the turbine blade cooling one. However, a wealth of closed form solutions exist for particular idealized cases, and with modern finite-difference calculation methods the laminar problem poses no particular difficulty even for very complicated boundary conditions.

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The remainder of the paper is in six sections. The case of a constant free-stream velocity is considered first. Experimental results covering the entire range of blowing and suction are presented for the case of a constant rate of transpiration and a constant surface temperature. Methods of extending these results to problems of arbitrarily varying transpiration rate and arbitrarily varying surface temperature, using the integral equations of the boundary layer, are then discussed.

Next the problem of an accelerated free-stream velocity is considered. It will be seen that simple correlations together with the integral equations of the boundary layer no longer appear adequate as tools for performance prediction, and a more fundamental theory is needed.

The two sections following contain a discussion of some models that can be used for direct solution (by finite difference methods) of the momentum and energy differential equations of the boundary layer. Constants for use in these models, derived from experimental data, are presented.

Finally, some examples of predictions using one of the models applied to some very difficult cases are presented. The paper ends with a discussion of some of the future problems which must be considered.

<u>The Case of Constant Free-Stream Velocity</u> - It is first instructive to examine the momentum integral equation of the boundary layer, particularized to the case of constant properties and constant free-stream velocity.

$$\frac{d\delta_2}{dx} = C_f/2 + v_o/U_\infty \tag{1}$$

An alternative form is obtained if the friction coefficient is factored from the right-hand side, and if the resulting term containing the transpiration velocity v_o is defined as a "blowing parameter", B_f .

$$\frac{d\delta_2}{dx} = (C_f/2)\left[1 + \frac{v_o/U_\infty}{C_f/2}\right] \tag{2}$$

or

$$\frac{dRe_M}{dRe_x} = (C_f/2)(1 + B_f) \tag{3}$$

In laminar boundary layer theory, similarity solutions to the momentum equations are obtained when B_f is maintained constant along the surface. Since for a laminar boundary layer $C_f/2$ varies as $Re_x^{-1/2}$, similar velocity profiles are then evidently obtained if the rate of blowing (or suction) decreases along the surface as the square root of distance x . For a turbulent boundary layer, an analogous situation obtains. B_f can be interpreted physically as the ratio of the transpired momentum rate to the shear force. If this ratio is maintained constant along the surface, similar profiles are obtained for a laminar boundary layer, and for a turbulent boundary layer a so-called "equilibrium" boundary layer is obtained, one that possesses outer

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region similarity of velocity profiles. However, in the case of the turbulent boundary layer, $C_f/2$ tends to vary with distance x approximately as $x^{-0.2}$, so constant B_f corresponds to a blowing rate, v_o/U_∞, which decreases in the direction of flow as $x^{-0.2}$.

The constant B_f boundary layer thus appears to be a fundamental case, and was a case studied by Simpson, et al [10]. Simpson also did extensive testing holding v_o/U_∞ constant rather than B_f. (Hereafter v_o/U_∞ will be referred to as F, the blowing fraction. Note also that since only the constant density case is being discussed, $v_o/U_\infty = \dot{m}''/G_\infty$, i.e., F is a mass flux ratio, and this is how it should be interpreted in compressible flow applications.)

One of the conclusions from Simpson's work is that $C_f/2$ can be expressed as a function of the momentum thickness Reynolds number, Re_M, and that virtually the same function or relationship is obtained for experiments at constant F as for experiments at constant B_f. In fact, this function seems to be remarkably independent of how F varies along the surface, and it is only for step changes in F that any appreciable difference can be noted.

A further point of interest that can be seen in the momentum integral equation, Eq. (1), is that when $v_o/U_\infty = F$ is negative, the possibility exists that the right-hand side of the equation will go to zero, leading to a situation where the momentum thickness of the boundary layer does not grow with x. As a matter of fact, this is a situation that will always be approached where F is maintained constant and negative. $C_f/2$ has a decreasing tendency with increasing momentum thickness, starting indefinitely high at the beginning of the plate, and so regardless of the value of F

this point of equilibrium will ultimately be reached. This is commonly called the "asymptotic suction layer". Note that for this condition B_f is equal to -1 , and thus -1 represents a lower limit on B_f .

Simpson's experiments can be correlated within the limits of the experimental uncertainty of the data itself by the following equation.

$$C_f/2 = 0.0130 \ Re_M^{-.25} \ [\ell n(1 + B_f)/B_f]^{0.77} \tag{4}$$

A plot of Eq. (4) is shown on Fig. 2. Since $C_f/2$ is contained within B_f , Eq. (4) is a little awkward to use. In Fig. 2, F is employed as a parameter as well as B_f , taking advantage of the experimental fact that the same functional relationship seems to be obtained for constant F experiments as for constant B_f experiments.

The fact that moving along a constant B_f line leads to decreasing values of F can be clearly seen. Note also that the lines for negative constant F are slightly concave upwards. If extended, these lines will approach values of $C_f/2$ equal to the negative of F .

An alternative way of presenting these results is shown in Fig. 3 where the ratio of C_f to the value of C_f , at the same momentum thickness Reynolds number for the case of no transpiration, is plotted as a function of another blowing parameter, b_f . b_f differs from B_f in that F is divided by the value of $C_f/2$ for no transpiration, at the same momentum thickness Reynolds number, rather than with respect to the local value of $C_f/2$. b_f is thus a more convenient parameter to use in presenting data on C_f because it does not contain the value of C_f sought. Interestingly, Eq. (4) plots as a single line on this diagram, and illustrates

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vividly the influence of blowing and suction on the friction coefficient.
Also superimposed on the diagram are the results of the laminar similarity
solutions mentioned earlier. The behavior is quite analogous, especially
on the suction side, but it should be noted that the laminar similarity
solution does lead to a zero value of C_f at $b_f = 3.47$. This point is
generally referred to as the "blow-off" point, and it is the value at which
the boundary layer is literally blown off the wall. However, Eq. (4) only
approaches zero and does not indicate a definite blow-off point. There is
no reason to suppose that blow-off is not just as real a phenomena for the
turbulent boundary layer as for a laminar boundary layer, but it is ex-
tremely difficult to detect experimentally because the uncertainty in mea-
suring C_f becomes indefinitely large as C_f approaches zero. The fact
that Eq. (4) does not indicate blow-off should only be interpreted as a
deficiency in the form of the empirical equation used. A popular rule of
thumb is that blow-off for a turbulent boundary layer will occur at approxi-
mately $F = 0.01$, and this rule of thumb is certainly not inconsistent with
the results on Fig. 3 when it is noted that the uncertainty in C_f/C_{fo} is
of the order of magnitude 0.1 at the right-hand extreme of the diagram.

The shape factor of the boundary layer, H , is often of interest,
and the following equation is an empirical representation of Simpson's mea-
surements for constant free-stream velocity.

$$H = 1.0/\{1.0 - 3.1\sqrt{C_f/2}[(1 + B_b)^{1/2} + (1 + 0.635 \, B_f)^{1/2}]\} \quad (5)$$

A plot of Eq. (5) is shown on Fig. 4. The significant point is that
blowing leads to high values of H , and that for F approaching 0.01 H

becomes greater than 2.0. It is perhaps worth noting that for the non-transpired boundary layer in an adverse pressure gradient, separation generally tends to occur when H is somewhere above 2.0. This is one of many instances of similarity between the behavior of the transpired flat-plate boundary layer with positive transpiration, and the adverse pressure gradient non-transpired boundary layer.

It is only fair to remark that the friction coefficient data of Simpson, upon which this entire discussion is based, while unquestionably the most extensive data available, has been the subject of some controversy. The case against it has been very well put by Squire [11], who feels that Simpson's friction coefficients are too high. There is ample room for argument because of the very considerable difficulty in measuring friction coefficients, and the different ways in which experimental data can be interpreted. The writer would only add at this time that friction coefficients close to Simpson's have been measured in subsequent work on the same apparatus, but in connection with accelerating flows, and where other methods of data interpretation have been used. Furthermore, there is an internal consistency between Simpson's data and the heat transfer data to be discussed shortly. There seems no question that the experimental uncertainty in the heat transfer data is less than for the friction coefficient data because the heat transfer rates are measured by a direct method. The internal consistency shows up particularly in the theoretical models to be discussed later; a single, reasonable model can be used to predict both the heat transfer data and the friction coefficient data equally well.

Turning now to the case of heat transfer with transpiration, but with free-stream velocity constant and surface temperature constant, consider first the energy integral equation of the boundary layer, particularized to the case of constant properties, constant free-stream velocity, and constant surface temperature.

$$\frac{d\Delta_2}{dx} = St + v_o/U_\infty \qquad (5)$$

A heat transfer blowing parameter, B_h , is defined in a manner completely analogous to the definition of the friction blowing parameter.

$$\frac{d\Delta_2}{dx} = St \left[1 + \frac{v_o/U_\infty}{St} \right] \qquad (6)$$

$$\frac{dRe_H}{dRe_x} = St(1 + B_h) \qquad (7)$$

The blowing parameter B_h can be interpreted physically as the ratio of the transpired thermal energy to the conducted thermal energy. Again, it is not surprising that holding this parameter constant along a surface leads to similarity of temperature profiles in the laminar boundary layer case, and apparently leads to outer region temperature profile similarity in the turbulent boundary layer case. There is also another physical significance to constant B_h . In the case of blowing, if the transpired fluid originates from a constant temperature source or plenum below the surface, and if there is no independent energy source within the surface wall

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itself, constant B_h yields a constant surface temperature t_o. In other words, to obtain a constant surface temperature by transpiring a fluid which starts out at a single uniform temperature, it is necessary for the blowing fraction F to decrease along the surface in such a manner as to hold B_h constant. Experimentally it is found that F must decrease as approximately the 0.2 power of x, as in the case of constant B_f.

Note further that a constant negative value of F must lead to an asymtotic suction layer for the same reasons as were discussed for the momentum boundary layer. In other words, under conditions of constant suction, Stanton number approaches a minimum, equal to the negative of F.

The experiments of Moffat and Kays [12], and of Whitten, et al [13], indicate that although B_h probably represents the more fundamental case, constant F yields to all intents and purposes the same value of Stanton number for a given value of enthalpy thickness Reynolds number, Re_H, and F. In fact, Whitten's results indicate that F can vary rather considerably along the surface with negligible effect upon this relationship. The very extensive data of Moffat, and of Whitten, for uniform surface temperature, can be represented quite adequately by the following empirical equation.

$$St = 0.0128 \ Pr^{-0.5}[\ell n(1 + B_h)/B_h]^{1.25}[1 + B_h]^{0.25} \tag{8}$$

Equation (8) is plotted on Fig. 5 where both B_h and the more convenient F are used as parameters. It should be emphasized that strictly speaking these data apply only for cases where the virtual origin of both the thermal and momentum boundary layers is the same, which means that the

thermal boundary layer thickness is of the same order magnitude as the momentum boundary layer thickness (actually about 10% greater for $Pr = 0.7$ and $F = 0$). If the thermal boundary layer thickness is substantially different from the momentum boundary layer thickness, and this can occur when the surface temperature varies or when the virtual origin of the thermal boundary layer differs from that of the momentum boundary layer, somewhat different results will be obtained, although the effects are not large. This is a problem that can best be discussed later in connection with a theoretical model of the thermal boundary layer.

An alternative method of presentation of these results is shown on Figure 6. Here everything is quite analogous to the very similar results for friction coefficient on Figure 3. A more convenient blowing parameter, $b_h = F/St_o$, is introduced, and the laminar similarity solution results are superimposed. In addition, the following equation proposed by Kutateladze and Leont'ev [9], based on analytic considerations, is included.

$$St/St_o = (1 - b_h/4)^2 \tag{9}$$

Figure 6 illustrates very graphically how blowing decreases the heat transfer coefficient, and this, of course, is the primary reason for the interest in transpiration as a method of protecting surfaces against high temperature fluids.

The same comments regarding blow-off as were made in connection with the friction coefficients apply here. It is, however, of interest to note that the laminar similarity solution for heat transfer does not really indicate

a blow-off point, because for $Pr = .70$ the Stanton number is still greater than zero at the point where the friction coefficient has reached zero. The Kutateladze and Leont'ev solution predicts blow-off at $b_h = 4.0$, but over most of the range it is very close to the laminar solution. The experiments of Moffat, and of Whitten, include extensive amounts of data for b_h equal to 4 and greater for which St is measurably above zero. It thus appears that the Kutateladze and Leont'ev solution predicts blow-off a little early.

It has been noted previously that both the friction coefficient and the Stanton number tend to be rather strong functions of local conditions along the surface rather than the history of the boundary layer up to the point in question. In other words, it has been stated that Stanton number tends to be primarily a function of local enthalpy thickness Reynolds number, Re_H, and F, and is relatively independent of how F has varied up to the point in question. Figures 7 and 8, which show some of the experimental data of Whitten et al [13] illustrate this point very well. Both a step down and a step up in F are shown. Although there is in both cases a detectable recovery region, most of the adjustment to the new boundary condition occurs very rapidly. These figures represent extreme cases, so it is not surprising that when there are continuous variations of F the recovery is virtually undetectable.

The friction coefficient behavior is substantially the same, although it is much more difficult to accurately measure friction coefficients just following a step in blowing.

The experimental facts that $C_f/2$ is primarily a function only of Re_M and F, and St is a function primarily only of Re_H and F, lead to

a relatively simple approximate procedure for calculating C_f and St along

a surface where F varies in some arbitrary manner. One need only sub-

stitute Eq. (4) into Eq. (3) and Eq. (8) into Eq. (7), and one has simple

differential equations for Re_M and Re_H which may be integrated, at least

numerically, to yield momentum and enthalpy thickness Reynolds numbers as

functions of x . Then Eqs. (4) and (8) are used to determine $C_f/2$ and

St at each local point. This procedure works remarkably well, although it

is, of course, restricted to the constant property boundary layer with

constant free-stream velocity, and in the case of the energy equation, we

have so far restricted consideration to the constant surface temperature

problem.

Whitten [13] also presents data for step changes in surface temperature

with transpiration, and includes step changes in transpiration rates. With

these results it is possible to use the method of superposition to build up

solutions to the energy equation for cases where both F and surface tempera-

ture vary in any arbitrary manner along the surface, and indeed Whitten has

done just this [14]. However, at some point one must question whether it is

not more fruitful to attempt to devise a more fundamental model of the

turbulent transport processes so that all manner of variations in boundary

conditions, and including varying fluid properties and viscous dissipation,

can be handled with a single consistent theory based on the differential

equations of the boundary layer rather than the integral equations. The

constant property, constant free-stream velocity problem with arbitrarily

varying F and arbitrarily varying t_o , probably represents about the limit

of integral methods, at least where reasonable precision is desired, although

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integral methods can be pushed considerably farther where low precision is adequate (see for example, Kays [15]).

<u>The Case of Varying Free-Stream Velocity</u> – The momentum and energy integral equations, for constant properties and constant surface temperature, but now including a variable free-stream velocity, may be written in the following form:

$$\frac{dRe_M}{dRe_x} = C_f/2 + F - K(1 + H)Re_M \tag{10}$$

$$\frac{dRe_H}{dRe_x} = St + F \tag{11}$$

where $\quad dRe_x = U_\infty dx/\nu \quad ; \quad K = (\nu/U_\infty^2)(dU_\infty/dx)$

$$dRe_M = d(U_\infty \delta_2)/\nu \quad ; \quad dRe_H = d(U_\infty \Delta_2)/\nu$$

One thing that is immediately apparent from these equations is that the pressure gradient, which is incorporated implicitly in the acceleration parameter K, has a direct influence on the development of the momentum equation, but any influence on the energy equation must be an indirect one coming in through the turbulent transport properties and the velocity profile. An important simplifying feature of the previously discussed constant free-stream velocity, constant surface temperature problem is that the momentum and thermal boundary layers tend to grow together, and this is one reason why relations such as Eq. (8) hold up remarkably well in the face of arbitrarily varying F along the surface. It should be apparent from Eqs. (10) and (11) that such will not be the case when free-stream velocity varies.

Before examining the kind of heat transfer behavior that is obtained

when free-stream velocity varies, it is worthwhile to look at some more
of the implications of Eqs. (10) and (11). A particularly interesting
case arises when the acceleration parameter K is a positive number and
constant. This is a case of an accelerated flow, and in fact, it is the
accelerated flow that is obtained when the fluid is confined between two
straight but converging planes. The possibility exists, according to
Eq. (10), that an equilibrium will be reached such that Re_M remains
constant, and indeed such flows are observed experimentally. There are
obvious experimental advantages in being able to hold such an important
parameter as Re_M constant, and so the constant-K boundary layer has
been extensively studied. Most of the experimental data obtained by the
author and his colleagues for varying free-stream velocity have been for
constant-K accelerating boundary layers. According to Eq. (11), however,
the thermal boundary layer will continue to grow even though the momentum
boundary layer has reached the equilibrium, or asymptotic condition.

It has been observed experimentally that strong acceleration, i.e.,
large K , will lead to a "retransition" of the turbulent boundary layer
to a laminar boundary layer. One might predict that this would occur simply
from examination of Eq. (10) alone. A large positive value of K will
cause Re_M to decrease, and if K is sufficiently large Re_M will be
pushed down into the region of stable laminar boundary layers. This is
not to say the Re_M is necessarily the sole criterion for "retransition",
but it certainly must be an important one. Assuming for the moment that
some critical value of Re_M (say, 200-400) is the dominant criterion,
Eq. (10) immediately suggests that positive values of F will tend to

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inhibit "retransition", while negative values will aid it. This is actually

what is experimentally observed. For $F = 0$ "retransition" or "laminarization"

will occur whenever K exceeds about 3×10^{-6} for an extended distance along

a surface. For F positive (blowing) a turbulent boundary layer can be main-

tained at considerably higher values of K without "laminarization". For F

negative (suction), both "laminarization" and the asymptotic suction layer are

approached at values of K lower than 3×10^{-6}, and it is very difficult to

tell which comes first when examining experimental data.

It is experimentally observed for the case of no transpiration,

$F = 0$, that laminarization causes a very substantial reduction in Stanton

number, but that this phenomena is a continuous function of K rather than

an abrupt change at some critical value of K. In other words, at any posi-

tive value of K there will be some reduction in Stanton number (relative to

what the Stanton number would have been at the same value of Re_H for $K = 0$).

On the basis of these facts, one might be led to conclude that blowing, under

accelerating free-stream conditions, would lessen the decreasing Stanton num-

ber tendency, while suction would enhance it.

On Fig. 9, the results of two test-runs reported by Thielbahr, et al [16]

are plotted. In the upper part of the diagram, there is a case of mild blow-

ing, $F = 0.001$, starting with constant free-stream velocity, followed by a

relatively strong acceleration, $K = 1.47 \times 10^{-6}$, and finally reverting to

constant free-stream velocity. The acceleration does, indeed, appear to cause

a small reduction in Stanton number. At the end of the acceleration, there

appears to be a start of a recovery back toward the line for $K = 0$, but it

is only a partial recovery. A reason for the failure of recovery would be

found if one examined the values of Re_M and Re_H in the recovery region.

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Re_H is very much greater than Re_M , and Re_M catches up only very slowly, a fact that an examination of Eqs. (10) and (11) would make apparent.

The second set of data on Fig. 9 shows a case of a rather strongly blown boundary layer, $F = 0.006$, that develops at constant free-stream velocity, and then is subjected to a relatively mild acceleration, $K = 0.8 \times 10^{-6}$. Surprisingly, the effect of acceleration is precisely the reverse of the previous case. Stanton number tends to increase above the value for no acceleration.

Attention has been drawn to these examples merely to illustrate the fact that the combination of transpiration and varying free-stream velocity results in complex interactions that are not going to be understood unless the basic turbulent transport mechanisms are studied in more detail. It does not appear practicable or fruitful to attempt to describe the kind of behavior shown on Fig. 9 by use of the integral equations and integral correlations. Let us examine a more fundamental approach.

Some Models for Solution of the Momentum Differential Equation of the
Boundary Layer - The digital computer has, in a period of no more than
five or six years, revolutionized turbulent boundary layer theory and
turbulent boundary layer prediction methods. Stable, accurate finite-differ-
ence solutions to the momentum, energy, and mass diffusion equations of the
boundary layer have become practicable. Fast, economic procedures are
available which include the possibility of an infinite variety of boundary
conditions as well as the inclusion of variable fluid properties, viscous
dissipation, chemical reaction, etc. It is not the present intention to
extoll the virtues of any particular procedure; actually there are several
very good ones in common use. However, the accuracy of any of the calcu-
lation procedures depends upon the basic physics introduced into it, and
this is our present concern.

The time-average momentum equation of the boundary layer, particular-
ized for the moment to constant fluid properties, and neglecting normal
turbulent stresses, may be written as follows:

$$\bar{u} \frac{\partial \bar{u}}{\partial x} + \bar{v} \frac{\partial \bar{u}}{\partial y} - \frac{\partial}{\partial y} \left[\nu \frac{\partial \bar{u}}{\partial y} - \overline{u'v'} \right] + \frac{g_c}{\rho} \frac{d\bar{P}}{dx} = 0 \tag{12}$$

If the turbulent shear stress $\overline{u'v'}$ were known at all points in the
boundary layer, the momentum problem simply becomes one of solution of Eq.
(12) for any desired boundary conditions, including transpiration. As dis-
cussed above, numerous adequate finite-difference methods are available to
do this.

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Although progress continues to be made in turbulent transport theory in general, and turbulent boundary layer theory in particular, it is still fair to say that there is as yet no truly fundamental turbulence theory that may be used as a universal starting point for solution of turbulence problems. Turbulent boundary layer theory has gone through, and continues to go through, a series of stages involving successively higher orders of sophistication. Each step in this process involves the correlation of experimental data at a more fundamental level, opening up the possibility of solving successively broader ranges of problems with a single consistent set of empirical constants. The information and calculating procedures to be presented here do not represent any very bold steps toward a more general theory. Higher order models are presently being investigated by numerous researchers, and hopefully will lead to theories that embrace still broader classes of applications, although probably at the price of complexity and computation cost. In the meantime, it will be demonstrated here that a relatively simple model can be used to adequately predict the behavior of the transpired boundary layer over a sufficiently wide range of conditions to make it a very practicable tool for engineering design.

We will first introduce the concept of eddy diffusivity for momentum, ε_M , as a convenient way of expressing the turbulent shear stress.

$$\overline{u'v'} = -\varepsilon_M \frac{\partial \overline{u}}{\partial y} \tag{13}$$

Already, we are in the realm of theoretical controversy, but the eddy diffusivity concept has the computational virtue of allowing one to use the

same computation program for both laminar and turbulent boundary layer. Since most turbulent boundary layers grow out of laminar boundary layers, the advantage is obvious.

It is convenient to visualize the turbulent boundary layer as consisting of an inner wall-dominated region, and an outer region which physically occupies most of the thickness of the boundary layer. However, for most applications the inner region turns out to be by far the more important one, and it is to this region that we will now focus primary attention.

The inner region is characterized by a region immediately adjacent to the wall in which viscous forces predominate ($\overline{u'v'}$ approaches zero), and a region farther out in which momentum transfer is almost entirely by turbulent transport processes, but in which the scale and intensity of the turbulence is still strongly dependent upon the proximity of the wall. Between these is a transition region in which both mechanisms are important in varying degrees. Some analysts prefer to represent the entire inner region by an empirically established "law-of-the-wall", but probably the most significant conclusion that has come out of the experimental work upon which this paper is based is that the thickness of the viscous-dominated part of the inner region is strongly dependent upon both transpiration and free-stream pressure gradient, and thus no single law-of-the-wall can be adequate.

The Prandtl mixing-length theory, despite much criticism for many years, still provides a remarkably adequate basis for describing the turbulent momentum transport process in the inner region. The mixing-length, ℓ , is defined such that it is related to the eddy diffusivity for momentum and the mean velocity gradient by the following equation:

$$\varepsilon_M = \ell^2 \; \frac{\partial \overline{u}}{\partial y} \tag{14}$$

Outside of the viscous-dominated region immediately adjacent to the wall, the mixing-length in the inner region of the boundary layer seems to be proportional to distance y from the wall, with a proportionality factor, k , which can be assumed to be independent of either transpiration rate or pressure gradient. Whether or not k is a truly universal constant is not highly important, because another constant to be described shortly can be used to absorb any dependence on transpiration or pressure gradient; the experimental data are not sufficiently accurate to justify finer discrimination. Thus we will model the region outside of the viscous near-wall region (which we will now term the viscous sublayer), but inside of the outer, or "wake", region, by:

$$\ell = ky \tag{15}$$

The viscous sublayer can be modelled in a simple way by introducing a damping function that forces the mixing-length ℓ to zero at the wall. Designating the damping function as D , the mixing-length may then be expressed as:

$$\ell = kyD \tag{16}$$

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The damping function D can be any function that is equal to zero at the wall and equal to unity at large values of y , but obviously some functions will model the actual experimental data better than others. The simplest possible damping function idealizes the sublayer as a finite, purely viscous region as follows:

$$D = 0.0; \qquad y^+ \stackrel{\leq}{=} < y^+_{crit}$$

$$D = 1.0; \qquad y^+ > y^+_{crit}$$

(17)

y^+ is the non-dimensional distance from the wall expressed in so-called wall coordinates. y^+_{crit} is then the effective viscous sublayer thickness. For the flat-plate case with no pressure gradient and no transpiration, y^+_{crit} is measured to be approximately 11.0 . For accelerating flows, y^+_{crit} increases, for transpiration y^+_{crit} decreases, and for suction y^+_{crit} increases. A reasonably adequate scheme can be developed using this very simple damping function, its main deficiency being that it does not model the velocity profile very well in the y^+ range 5 to 30.

Another scheme which is very popular today was first suggested by Van Driest [17]. The Van Driest damping function is exponential, so that the influence of the viscous region decays smoothly from the wall and is still felt as far out $y^+ = 50$. The simplest version of the Van Driest damping function is that given in the following equation:

$$D = 1.0 - \exp(-y^+/A^+)$$

(18)

In this case A^+ becomes an effective viscous sublayer thickness that must be established experimentally. For the flat-plate case, $A^+ = 26.0$ is approximately what is obtained. However, the precise value of A^+ depends to a certain extent upon the value of the mixing-length constant k. Fig. 10 shows the values of A^+ that are obtained for differing values of k to yield identical values of u^+ at an arbitrary $y^+ = 80.0$. There is some evidence that at low Re_M, i.e., below 6000, k tends to increase (see Simpson [18]), but the data upon which this paper is based are not conclusive in this regard, and $k = 0.44$ is used for all of the illustrative calculations to follow. Those preferring other values of k can use Fig. 10 as a basis for changing all of the functions, as will be seen later.

Like y^+_{crit}, A^+ is found to be a function of both a pressure gradient parameter and a transpiration parameter (and perhaps other things yet to be investigated).

The non-dimensional distance from the wall y^+ contains within it the wall shear stress τ_o. A^+ is simply the value of y^+ at the effective outer edge of the sublayer, and thus is also normalized with respect to τ_o. It can be argued that in the region under consideration, a local Reynolds number of turbulence can be defined as

$$Re_t = \ell_t \sqrt{\overline{u'v'}}/\nu = \ell_t \sqrt{g_c \tau_t / \rho}/\nu \qquad (19)$$

ℓ_t, the turbulence length scale, is effectively equal to the mixing-length $\ell = ky$. Thus,

$$Re_t = ky\sqrt{g_c \tau_t / \rho}/\nu \qquad (19a)$$

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In the flat-plate case, the shear stress τ is essentially equal to τ_o out to a distance well beyond the viscous sublayer, so that Re_t is equal to ky^+. Thus A^+ can be interpreted as a critical value of the local Reynolds number of turbulence, which determines the viscous sublayer thickness.

If this idea has any merit, however, account should be taken of the fact that for flows with pressure gradients in the flow direction, or for transpired boundary layers, the local shear stress can vary markedly with y in the region near the wall, in which case the _local_ shear stress ought to be used in the damping function rather than the wall shear stress. This notion has led to the following alternative form for the Van Driest damping function

$$D = 1.0 - \exp\left(-y^+ \sqrt{\tau^+} / A^+\right) \qquad (20)$$

$$\text{where} \quad \tau^+ = \tau / \tau_o$$

Eq. (20) is appealing because it immediately suggests that for strongly accelerating flows, where τ^+ decreases very sharply with increasing y^+, it would be possible to have a sufficiently strong pressure gradient so that D would remain a very small number throughout the boundary layer. This, of course, would correspond to the observed phenomena of laminarization. In other words, this would be a situation where Re_t never exceeds a critical value of the Reynolds number of turbulence, A^+, and the entire boundary layer remains laminar. Similarly, the thinning of the laminar sublayer that is observed for blowing is correctly modelled, as well as the opposite trend for suction. These features have led some analysts to prefer Eq. (20) over

Eq. (18) as a basis for defining A^+. It is important to recognize this difference (and there are still other schemes) because otherwise it is difficult to compare the results of different workers.

Despite the fact that Eq. (20) with $A^+ = 26.0$ yields the _trends_ that are observed experimentally, the observed effects are considerably greater. Thus for transpiration and/or pressure gradient it still becomes necessary to express A^+ as a function of a transpiration parameter and a pressure gradient parameter. Using wall coordinates, these are respectively v_w^+ and P^+ (defined in the Nomenclature). All that is accomplished by using Eq. (20) in preference to Eq. (18) is to somewhat lessen the dependence of A^+ upon these parameters.

Values of A^+ as defined by Eq. (20) have been determined from a wide range of experimental velocity profiles for blowing, suction, acceleration, and combinations of acceleration and transpiration, from the work of Simpson et al [10], Julien et al [19], and Loyd et al [20]. The following equation is an empirical correlation of these results and fits them all with a tolerance of about ±10 percent.

$$A^+ = \frac{A_o^+}{5.15 \left[v_w^+ + \frac{5.86\ P^+}{(1 + 5.0\ v_w^+)} \right] + 1} \qquad (21)$$

where A_o^+ is the value of A^+ for no transpiration and no pressure gradient, as taken from Fig. 10. Eq. (21) is plotted on Figure 11 for the case of $k = 0.44$. Here the effects of a favorable pressure gradient, and the effects of transpiration are clearly seen. Although the equation itself extends to positive values of P^+, and may well be valid there, the experimental data upon which Eq. (21) is based are all for the case of accelerating flows (negative P^+).

Another damping function, proposed by Evans et al [21], utilizes a linear function as follows:

$$D = y^+/B^+ \quad ; \quad y^+ \leq B^+$$

$$= 1.00 \quad ; \quad y^+ > B^+ \tag{22}$$

In this case B^+ becomes an effective viscous sublayer thickness. This function involves a discontinuity, but only in D, not in either mixing-length or the resulting velocity profile. For very high Prandtl number fluids, the heat transfer behavior turns out to be very sensitive to the behavior of the damping function at small values of y^+. It is important to note that the Van Driest function is also linear at small values of y^+, so either scheme seems equally attractive from this point of view. (Both schemes result in ε_M varying as $y^{+^{\frac{1}{4}}}$ as the wall is approached). The Evans damping function seems to give velocity profiles that are a little closer to the experimental data at values of $y^+ < 100$, and therefore is especially attractive for low Reynolds number work.

Like A^+, the value of B^+ depends to a certain extent upon the value of mixing-length constant k used. Fig. 10 shows this dependence.

Values of B^+ have been determined from the experimental data cited above, and these results are correlated within about ± 10 percent by the following equation.

$$B^+ = \frac{B_o^+}{\left[9.0 \ v_w^+ + \dfrac{3.35 \ P^+}{1.0 + 4.0 \ v_w^+} \right] + 1} \tag{23}$$

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Again, B_o^+ is taken from Figure 10. Eq. (23) is plotted on Fig. 12 for the case of $k = 0.44$. The behavior is very similar to that of A^+; there is a somewhat stronger effect of v_w^+ and P^+, but this is due to the fact that Eq. (20) contains the shear stress ratio τ^+. Had the comparison been made with A^+ defined as in Eq. (18), it would be found that B^+ differs from A^+ by almost a constant factor.

Figs. 13 and 14 show examples of the inner region velocity profiles (i.e., the law-of-the-wall) as predicted by integration of Eq. (12) using the B^+ scheme, and a comparison with experimental data, in the y^+ range 10 to 100 (the temperature profile on Fig. 13 will be discussed presently.) The data on Fig. 13 are from several test runs for $v_w^+ = 0.0$ and $P^+ = 0.0$; they correspond closely to the data that are generally found in the literature for this basic case. Fig. 14 shows two cases of blowing alone, and one case of a combination of blowing and acceleration. As can be seen, the prediction is excellent; the A^+ scheme is also quite adequate, but slightly underpredicts u^+ in the y^+ range 10-50.

A word should now be said about the reasons for the observed dependence of A^+ or B^+ upon P^+ and v_w^+. It was suggested earlier that the outer edge of the viscous sublayer might be characterized by some critical value of the local Reynolds number of turbulence, Eq. (19), and indeed this hypothesis does yield the correct <u>trends</u> with v_w^+ and P^+. This scheme applied to A^+ yields somewhat better results than if applied to y_{crit}^+. If applied to B^+, it works better still. Since each of these three schemes employs an effective sublayer thickness that is greater than the previous one, it appears possible that the value of y^+ at which fully developed self-sustaining turbulence obtains, and therefore at which the Reynolds number of

turbulence has exceeded some critical value, is at a value of y^+ greater

than the effective sublayer thickness used in these models. An investigation

by Andersen [22] based on the experimental data cited above, has revealed

the interesting fact that if y^+ equal to about 2 times B^+ is chosen as

the critical point, the local Reynolds number of turbulence is very close

to the same number for all of the data examined. These data include some

37 test runs covering a wide range of transpiration, acceleration, and

combinations of transpiration and acceleration. To be more precise, at

$y^+ = 2B^+$, the Reynolds number of turbulence, Re_t , is equal to about 33.0

over the entire range of tests. In other words, with this fact alone it is

possible to calculate B^+ as a function of P^+ and v_w^+ and to obtain

results that correspond remarkably closely to those represented by Eq. (23),

or Fig. 12. Thus the concept that the thickness of the viscous sublayer is

determined by a critical Reynolds number of turbulence seems at the present

time to be an attractive one.

A^+ and B^+ as represented by Eqs. (21) and (23), were obtained under

what might best be described as equilibrium conditions, i.e., conditions under

which v_w^+ and/or P^+ are invariant or at worst are varying only slowly along

the surface. Under non-equilibrium conditions where these parameters are

changing rapidly or abruptly, it has been observed that the sublayer does

not change abruptly to its new equilibrium condition, i.e., B^+ does not

immediately assume its new equilibrium value. Given the stability nature of

the problem, this is probably not surprising. It can be hoped that some of

the higher order models of turbulence will predict this effect, but in the

meantime, a reasonably satisfactory expedient is to use a rate equation of a

type suggested by Launder [23]:

$$\frac{dB^+_{eff}}{dx^+} = (B^+_{eff} - B^+_{eq})/C \tag{24}$$

$$(\text{or} \quad A^+_{eff} \quad \text{and} \quad A^+_{eq})$$

B^+_{eff} is the effective value of B^+, while B^+_{eq} is the equilibrium value obtained from Eq. (23). A value of C of about 4.0 has been found to be reasonable.

All of the discussion up to now has been concerned with the inner region of the boundary layer. The outer region, comprising the greater part of the boundary layer thickness, is of considerably less importance in predicting performance, and thus can be handled successfully using more gross approximations. This statement may not be valid for very non-equilibrium boundary layers under adverse pressure gradient conditions, but how valid it actually is for accelerating flows and for transpired boundary layers with and without acceleration, will be demonstrated later. In any case, for equilibrium or near equilibrium boundary layers, either the assumption of a constant value of eddy diffusivity over the entire outer region, or the assumption of a constant value of mixing-length over the entire outer region yields approximately the same result. If constant eddy diffusivity is used, an empirical correlation of eddy diffusivity as a function of momentum thickness Reynolds number can be obtained. However, if mixing-length is used in the inner regions, it is probably computationally simpler to use the mixing-length concept for the entire boundary layer. A satisfactory scheme is to express the outer region mixing-length as a fraction, λ of the total boundary layer thickness.

$\lambda = 0.085$ or 0.080, based on the 99% boundary layer thickness, works remarkably well for the entire range of test data on which this paper is based. One simply evaluates ℓ from Eq. (15) until the value obtained exceeds $\ell = \lambda \delta_{.99}$, and then uses the latter value for the remainder of the boundary layer.

At values of Re_M less than about $6,000$, there is some evidence that λ is greater. The following is an empirical fit that works reasonably satisfactorily.

$$\lambda = 0.25 \ Re_M^{-1/8} \quad ; \quad Re_M \leq 6000 \tag{25}$$

Boundary layer predictions for strong blowing are somewhat improved if this latter correction is not made. This may actually be indirect evidence that the mixing-length constant, k , is lowered by strong blowing, but this kind of discrimination is very difficult to make from the available data.

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A Model for Solution of the Energy Differential Equation of the Boundary

Layer - The time-averaged energy equation of the boundary layer, particularized to constant fluid properties and negligible viscous dissipation, and neglecting turbulent conduction in the stream direction, may be written as follows:

$$\bar{u} \frac{\partial \bar{t}}{\partial x} + \bar{v} \frac{\partial \bar{t}}{\partial y} - \frac{\partial}{\partial y} \left[\alpha \frac{\partial \bar{t}}{\partial y} - \overline{t'v'} \right] = 0 \tag{26}$$

This equation can be solved for any desired boundary conditions providing that the velocity field has been established first by solution of the momentum equation, and providing that we have information on $\overline{t'v'}$. For convenience we introduce the concept of eddy diffusivity for heat, ϵ_H.

$$\overline{t'v'} = -\epsilon_H \frac{\partial \bar{t}}{\partial y} \tag{27}$$

Although it might be fruitful to attempt to evaluate $\overline{t'v'}$ or ϵ_H on the basis of assumptions that are independent of the turbulent shear stress, it seems apparent that there is some kind of relationship between $\overline{t'v'}$ and $\overline{u'v'}$, or ϵ_H and ϵ_M. Therefore most analysts have found it convenient to introduce the concept of turbulent Prandtl number, Pr_t, defined as follows:

$$Pr_t = \frac{\epsilon_M}{\epsilon_H} \tag{28}$$

Introducing Eqs. (27) and (28) into Eq. (26) we obtain:

$$\bar{u}\,\frac{\partial\bar{t}}{\partial x} + \bar{v}\,\frac{\partial\bar{t}}{\partial y} - \frac{\partial}{\partial y}\left[(\alpha + \varepsilon_M/Pr_t)\,\frac{\partial\bar{t}}{\partial y}\right] = 0 \qquad (29)$$

If Pr_t were known, Eq. (29) could be solved for any desired boundary conditions so long as the momentum equation must be solved anyway. Evaluation of the turbulent Prandtl number is then one of the central problems of turbulent heat transfer.

A very simple physical model of the turbulent momentum and energy transfer process leads to the conclusion that $\varepsilon_H = \varepsilon_M$, i.e., $Pr_t = 1.00$ (the "Reynolds Analogy"). Slightly more sophisticated models suggest that $Pr_t > 1.00$ for $Pr < 1.00$, and still other models suggest that Pr_t equals 0.7 or 0.5 in turbulent wakes. The experimental data are not abundant, but Figs. 15 and 16 show the measurements, respectively, of Simpson, Whitten, and Moffat [24], and of Kearney, Moffat, and Kays [25], with air as a working substance. These were all evaluated from measurements of the slopes of mean velocity and temperature profiles, together with estimates of shear stress and heat flux distributions, and the experimental uncertainty is high. The data on Fig. 15 are all for constant free-stream velocity, but cover a wide range of blowing and suction conditions. The data on Fig. 16 are for accelerated boundary layers with a considerable range of blowing and suction.

The choice of y^+ as a basis for comparison is made only because a slightly better correlation is obtained than with other parameters, but even this fact is debatable. However, despite the very considerable scatter of data, two conclusions seem definitely warranted. First, the turbulent Prandtl number, at least for air, has an order of magnitude of unity. This fact alone is significant, for it provides considerable justification for

use of the concept of eddy diffusivity in the first place. The second conclusion is that Pr_t is evidently less than 1.00 in the wake or outer region, while it is greater than 1.00 near the wall. Very near the wall the experimental uncertainty becomes excessive, so that it is not possible to establish any limiting value in this rather critical region.

Another way of evaluating turbulent Prandtl number from experimental temperature profiles is to determine by computer experiments the turbulent Prandtl number function that must be used in order to predict given experimental temperature profiles using Eq. (29) and any one of the models for ε_M previously discussed. In effect the experimental velocity profile is partially removed from consideration, but different results for Pr_t will be obtained depending upon how well the model for ε_M reproduces true velocity profiles.

Using this scheme, the results on Fig. 17 for the inner region were obtained using the B^+ damping function, Eq. (22), and experimental data for several cases of blowing and acceleration, as well as some data based on the average of several non-transpired, non-accelerated runs. B^+ for each run was evaluated from the corresponding velocity profile by a computer experiment that forced the computed and measured velocity profiles to match at an arbitrary $y^+ = 80$. The results are thus independent of velocity measurements at very small values of y^+ . For $y^+ > 30$ there seems no question that the most probable value of Pr_t is simply 1.00, i.e., the Reynolds Analogy, regardless of transpiration or acceleration. On the other hand there is a definite indication of much higher Pr_t for lower y^+ . Note of course that since only $y^+ < 100$ is being considered, the lower Pr_t

that seems characteristic of the outer region is not seen.

The same computer experiment performed with the A^+ damping function tends to yield values of Pr_t that average about 0.9 for $y^+ > 30$, but otherwise the same conclusions are reached. However, since the B^+ scheme generally predicts the velocity profiles more accurately in the region $y^+ < 100$, it would seem that the results on Fig. 17 more nearly represent the truth. (The term "truth" here must be qualified, because if A^+ is being used in a boundary layer prediction scheme, it is necessary to use values of Pr_t based on A^+, even though these values differ from what might be measured directly.)

The absolute necessity for a high Pr_t at very low values of y^+ is well illustrated by the data and calculations for temperature profile shown on Fig. 13 for the case of no transpiration and no pressure gradient. For y^+ greater than about 30 two parallel lines describe the velocity and temperature profiles very well. This can be interpreted as meaning that $Pr_t = 1.00$. However, if an attempt is made to predict the temperature profile with $Pr_t = 1.00$ throughout, the result is as shown. A higher Pr_t is needed for $y^+ < 30$ to avoid badly underpredicting t^+ in the y^+ range 30 to 100. Closer examination reveals that it makes little difference how high Pr_t is for $y^+ < 10$, and at $y^+ = 15$ it is already too late to introduce the correction. Apparently for $y^+ < 15$ turbulent velocity fluctuations result in a considerably greater rate of transport of momentum than of heat for reasons that are not yet fully understood. The heat conduction model of Jenkins [25] would predict this trend, but not the concentration of the effect in this narrow region.

At the present time an empirical correlation of this effect seems the

most practicable expedient if heat transfer rates are to be calculated. The assumption of a constant value of $Pr_t = 0.9$ throughout the boundary layer will yield overall heat transfer rates that are quite satisfactory; temperature profiles can be more accurately predicted if a variation of Pr_t with y^+ is introduced. The following correlations have been used successfully by the writer for air:

If A^+ is used as the damping function (Eq. 20):

$$Pr_t = (1.43 - 0.17 \, y^{+1/4})(1 + v_w^+)$$

(30)

$$\text{If } Pr_t < 0.85 \; ; \; Pr_t = 0.85$$

The dependence upon v_w^+ in Eq. (30) is not based on the raw Pr_t data in Figs. 15 and 16, but rather upon computer experiments using the A^+ scheme for evaluation of ε_M.

If B^+ is used as the damping function:

$$Pr_t = 1 + .35[1 + \cos(\pi y^+/37)]; \; y^+ < 37$$

$$Pr_t = 1.00 \; ; \; y^+ > 37$$

(31)

$$Pr_t = 0.60 \; ; \; y > (\lambda \delta_{.99}/k)$$

Figure 13 shows an example of an inner region temperature profile prediction using the B^+ damping function, and Eq. (31). Actually there are any number of Pr_t functions that work equally well, and the cosine function in Eq. (31) has no physical significance.

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<u>Some Examples of Predictions of Difficult Cases</u> - To illustrate the quality

of turbulent boundary layer predictions which can be made with a finite

difference program using some of the material presented in the preceding

sections, two rather difficult cases have been chosen. A modification of the

Spalding/Patankar [27] program was used, although any good finite difference

should yield similar results.

The first of these is illustrated on Fig. 18. Here is a case of rather

strong blowing, $F = 0.0058$, subjected to a moderately strongly accelerated

free-stream along about half the length of the test section, with constant

free-stream velocity thereafter. The variation of U_∞ is shown at the top

of the diagram. This figure then shows a comparison between measured and

predicted values of Re_M , $C_f/2$, and St , plotted as functions of distance

x along the test surface. (The fluid used was room temperature air with small

temperature differences,as was the case for all of the data considered in this

paper). Note that Re_M is almost constant throughout the accelerated region,

a consequence of a constant-K accelerated boundary layer, as discussed earlier.

Following acceleration Re_M then increases rapidly. The comparison between

experiment and prediction must   be considered excellent, especially when

experimental uncertainty is taken into consideration. This prediction was made

using the A^+ damping function, but the B^+ scheme will do equally as well.

Figures 19 and 20 are in some ways more impressive, for they show two

velocity profiles and two temperature profiles from the same test run, in real

dimensional coordinates, and a comparison between experiment and predictions.

In each case the profile at $x = 45.6$ inches is one taken in the

accelerated region, while the other profile at $x = 69.7$ inches is taken
in the recovery region farther downstream.

Another difficult case is shown on Fig. 21. In this run, the flow
starts at constant free-stream velocity but with moderately strong blowing,
$F = .004$. This flow is then subjected to a very strong acceleration starting at
$x = 2$ feet. In approximately the middle of the accelerated region the
blowing is removed entirely. Then at about $x = 3.4$ feet the acceleration
is removed, and for the remainder of the test section there is no blowing
and no acceleration. The Stanton number prediction shown here was obtained
using the A^+ scheme, although the outer region of the boundary layer was
calculated using a higher order turbulence model than the simple mixing-length
model described here. The turbulent kinetic energy equation was solved, and
ε_M was determined from an assumed relationship between ε_M and the kinetic
energy of turbulence. However, it has been found that in flows of this type,
use of a turbulent kinetic energy model in the outer part of the boundary
layer contributes very little, and to all intents and purposes identical
results will be obtained with the simple mixing-length model.

The important thing to note here is that the model responds remarkably
to the abrupt changes in boundary conditions, and predicts the resulting
non-equilibrium boundary layer very well indeed. Of particular significance
is the abrupt rise in Stanton number following the removal of blowing. The
ability of the prediction to follow the data at this point is heavily dependent
upon the use of the rate equation and lag constant, Eq. (24).

<u>Future Work</u> - The needed future research on heat transfer to the transpired turbulent boundary layer can be subdivided into two general categories. The first of these is research that is concerned with the turbulent boundary layer in general, while the second is research particularized to the transpired case. All advances in the general category will contribute to a better understanding of the transpired boundary layer. Mention has already been made of investigations into higher order models of turbulence, and as time goes on the results of these investigations can be expected to find their way into transpired turbulent boundary layer theory and prediction methods.

The problem of turbulent Prandtl number, or more generally, the problem of the turbulent transport of thermal energy, should provide a particularly fruitful area for research in the near future. This is obviously not a new problem; it was of great interest two decades ago when liquid metal heat transfer was first being actively investigated. However, it is clear that a better theory is needed than is presently available.

For the particular case of the the transpired turbulent boundary layer, the first obvious problem for which there is presently inadequate experimental data is the case of an adverse pressure gradient. Actually, the case of blowing and adverse pressure gradient is going to lead to early stall, so it is probably the suction problem with an adverse pressure gradient that provides the most scope for experimental work. It should be noted that Eqs. (21) and (23) will yield some numbers for use with adverse pressure gradients, and since these numbers can be predicted by a theory that also predicts the other cases relatively satisfactorily, one might be temped to use these equations for predicting the adverse pressure gradient cases.

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However it should be emphasized that no experimental data for the adverse pressure gradient cases has been used to generate these empirical equations.

In attempting to apply the transpiration results and theories presented, two obvious practical problems arise. Real transpiration surfaces frequently are aerodynamically rough, and real transpiration surfaces frequently are constructed by drilling a large number of small discrete holes in a surface rather than constructing the surface from a sintered powder as was done for all of the experimental data considered here. The roughness problem is presumably not totally unlike the roughness problem for the non-transpired case, although the effects of transpiration are at the present time completely unknown. The large hole problem opens up an entire field of investigation because the number of possible geometrical variables increases enormously. There is an entire spectrum of problems lying between what might be called pure transpiration, with surface holes and spacing small relative to the laminar sublayer thickness at one end, and film cooling at the other end.

In a sense, the present paper has been concerned with only an idealized case lying at one end of a very broad spectrum of turbulent boundary layer problems.

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Acknowledgements

Figure Captions

Fig. 1   Diagram of the Boundary Layer System Considered

Fig. 2   Friction Coefficients for the Case of Constant Free-Stream Velocity
With Various Values of Constant Blowing Fraction F, and Constant
Blowing Parameter B_f

Fig. 3   Ratio of Friction Coefficient to Friction Coefficient Without
Transpiration but at the same Momentum Thickness Reynolds Number,
Plotted as a Function of a Modified Blowing Parameter b_f. Comparison
of Turbulent Experiments With the Exact Laminar Similarity Solution
Results

Fig. 4   Shape Factor Plotted as a Function of Blowing Fraction for Various
Values of Momentum Thickness Reynolds Number, Free-Stream Velocity
Constant

Fig. 5   Stanton Number Plotted as a Function of Enthalpy Thickness Reynolds
Number for Various Constant Values of Blowing Fraction, and Various
Constant Values of Blowing Parameter for a Fluid With $Pr = 0.7$
and Constant Free-Stream Velocity

Fig. 6   The Ratio of Stanton Number to Stanton Number for No Blowing at the
Same Value of Enthalpy Thickness Reynolds Number Plotted as a Function
of Modified Blowing Parameter. Comparison of the Experimental Results
With the Exact Laminar Similarity Solutions, and the Equation Proposed
by Kutateladze and Leont'ev.

Fig. 7   The Effect of a Step-Down in Blowing on the Stanton Number, Constant
Free-Stream Velocity

Fig. 8   The Effect of a Step-Up in Blowing on the Stanton Number, Constant
Free-Stream Velocity

Fig. 9   The Effects of a Combination of Blowing and Acceleration on Stanton
Number

Fig. 10   Values of Effective Sublayer Thicknesses A^+ and B^+ for the Case of
No Transpiration and No Pressure Gradient Plotted as a Function of
Mixing-Length Constant, k . The Indicated Combinations of A^+ or B^+
and k will Predict Identical   Values Of u^+ at $y^+ = 80$.

Fig. 11   The Effect of Pressure Gradient and Transpiration on A^+, Curves
Generated from an Empirical Fit to Available Data

Fig. 12   The Effect of Pressure Gradient and Transpiration on B^+, Using the
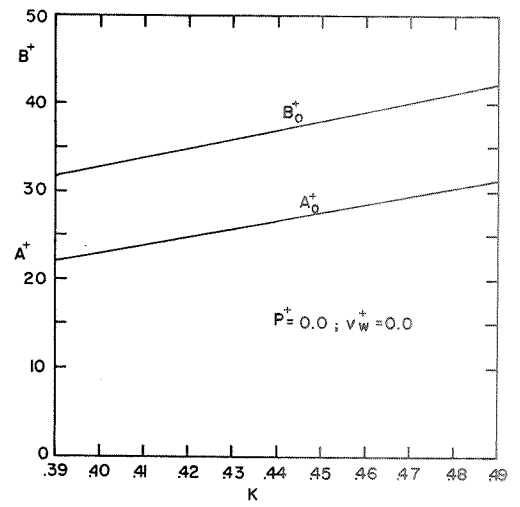Same Basic Data as for Fig. 11
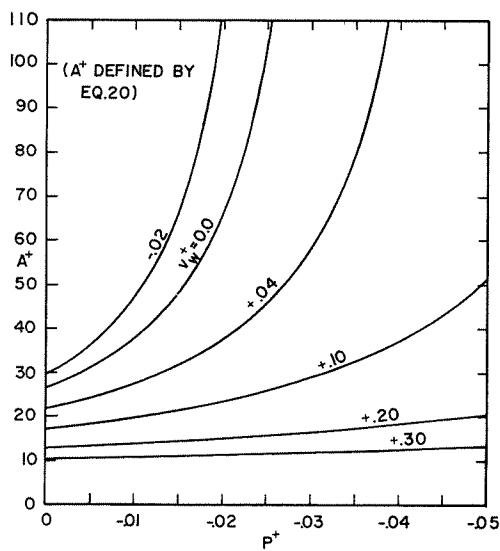
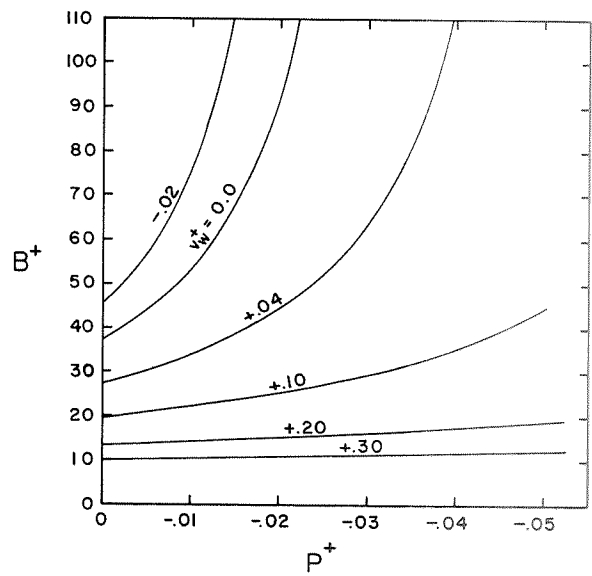FIG. 1

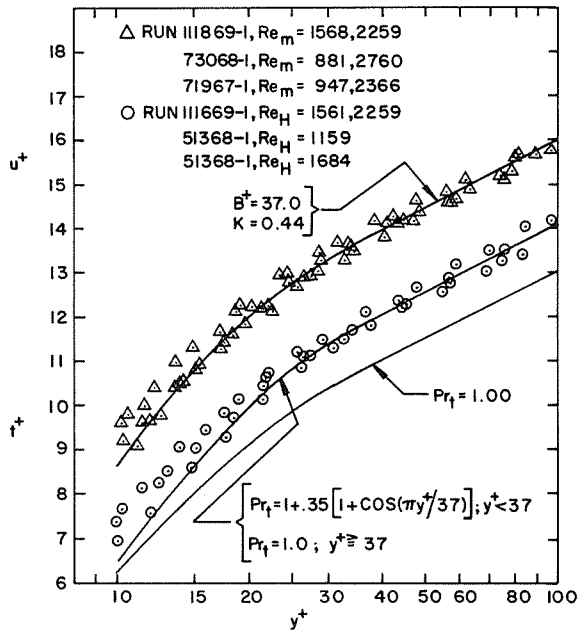FIG. 2

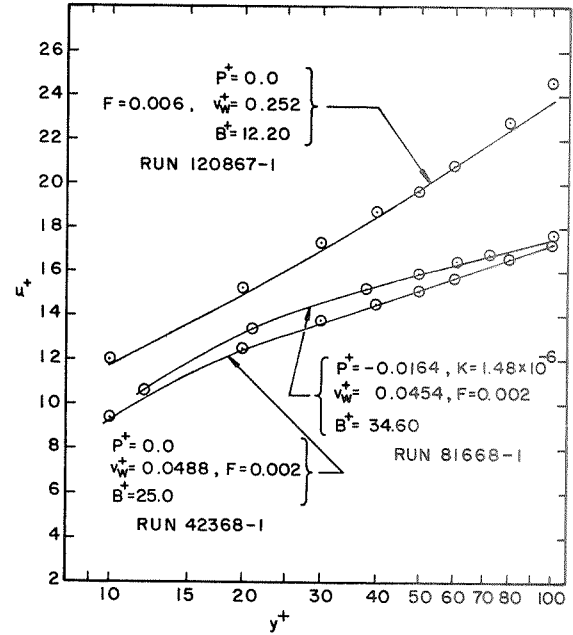FIG. 3

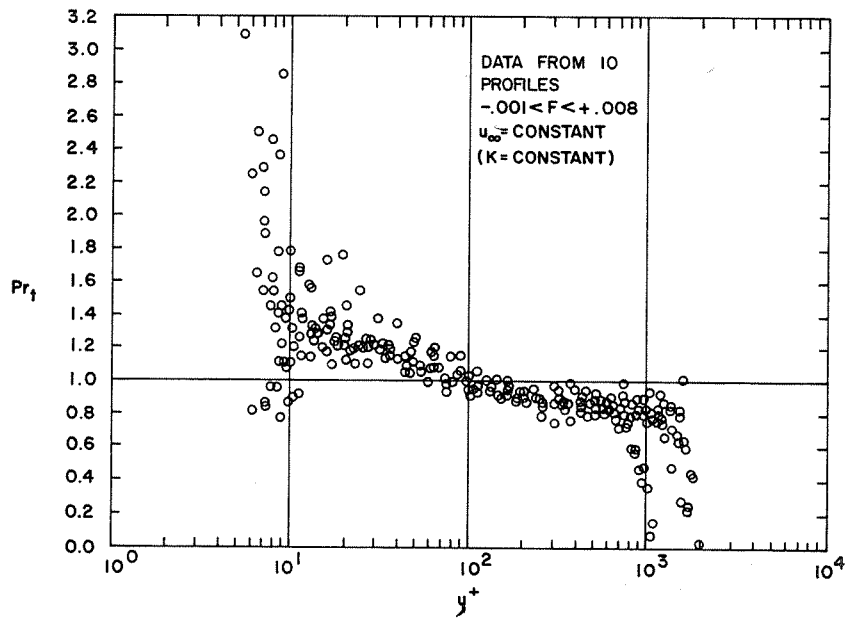FIG. 4

FIG. 5

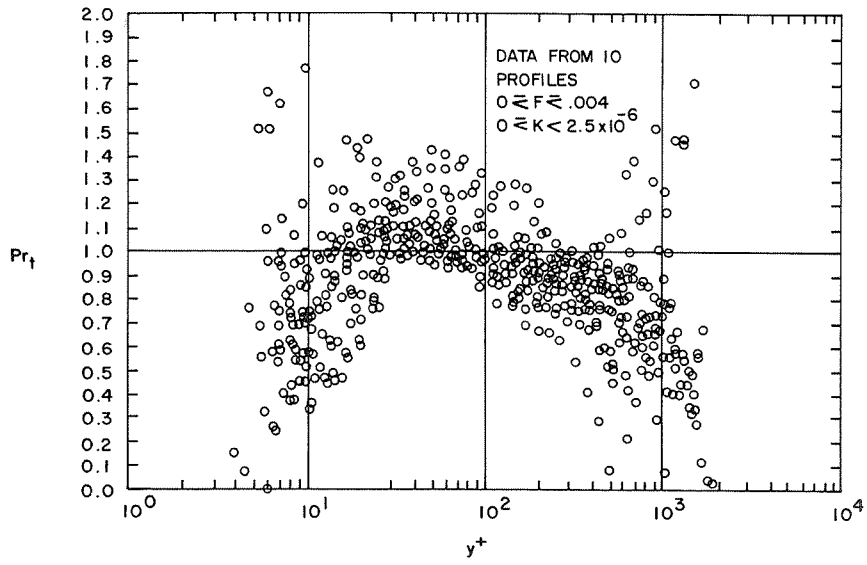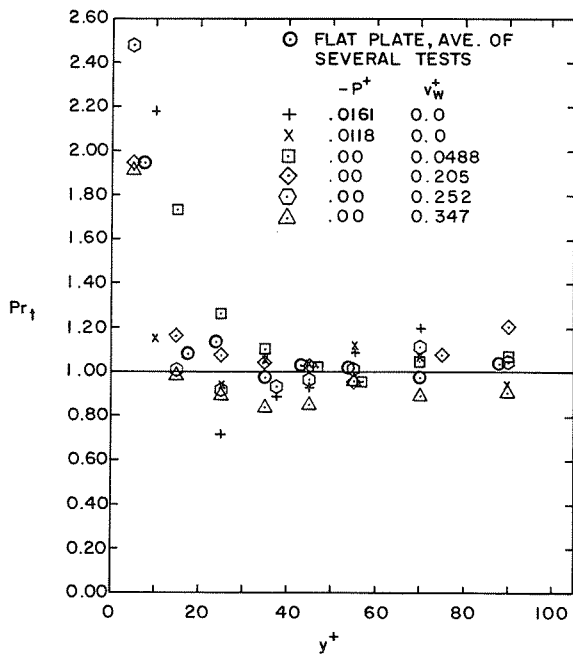FIG. 6

FIG. 7

FIG. 8

FIG. 9
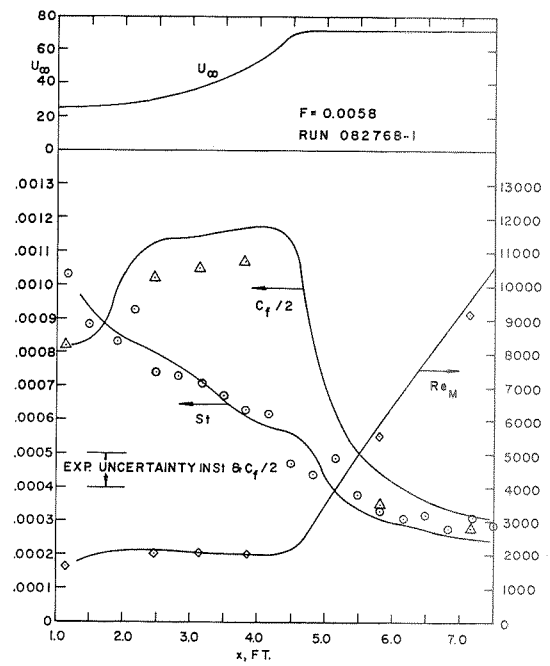

FIG.10

FIG.11

FIG.12

FIG.13

FIG. 14

FIG.15

FIG.16

FIG.17

FIG.18

FIG. 19

FIG. 20

FIG. 21