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THE UNIVERSITY OF ALBERTA

NUMERICAL AND EXPERIMENTAL STUDIES OF RIME ICE ACCRETION ON
CYLINDERS AND AIRFOILS

by

C

KAREN JOAN FINSTAD

THESES

SUBMITTED TO THE FACULTY OF GRADUATE STUDIES AND RESEARCH
IN PARTIAL FULFILMENT OF THE REQUIREMENTS FOR THE DEGREE
OF DOCTOR OF PHILOSOPHY

IN

METEOROLOGY

DEPARTMENT OF GEOGRAPHY

EDMONTON, ALBERTA

Fall, 1986

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PERMANENT ADDRESS:

#901 - 10149 Saskatchewan Drive

Edmonton, Alberta

T6E 6B6

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METEOROLOGY.

Edward Webster

Supervisor

ABW.....
ERKinsel.....
SM Gates.....

J. S. Clark.....

External Examiner

Date.....July 9, 1986.....

Abstract

Four separate studies describe improvements to numerical techniques used in the modelling of ice accretion on circular cylinders and on airfoils.

1. For collision efficiencies, accretion limits and impact velocities on circular cylinders, tables of numerical data and analytical approximating functions are presented. These data replace those of Langmuir and Blodgett (1946).
2. Parameterizations of the local collision efficiency distribution on circular cylinders are derived from numerical data, and applied in an 'operational' style computer model of rime ice accretion. The model's performance is compared with experimental accretions grown in the University of Alberta FROST tunnel.
3. The distribution of local ice density in rime ice is investigated in experimental wind-tunnel accretions made on circular cylinders. A form of the distribution is derived for application to the numerical model described in the previous study.
4. A second model, for rime ice accretion on airfoils or other arbitrary shapes, combines droplet trajectory integrations and parameterizations for local collision efficiency and density distribution. By use of these parameterizations, and of simple numerical methods throughout, the model is made significantly faster than previous models have been. Verification of the model is

provided by comparisons with several wind-tunnel
experiments.

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List of Symbols

<u>Symbol</u>	<u>Definition</u>
c_d	- droplet drag coefficient
$c_{x,n}$	- constant coefficients for Equation II.11
D_c	- circular cylinder diameter
D_d	- droplet diameter
d_{42}	- diameter of droplet at the 42nd volume percentile of the droplet spectrum
E	- overall collision efficiency
h_0	- airfoil width as a fraction of chord length
K	- Langmuir non-dimensional mass parameter
K_0	- a function of K and Re , used in analytical fits to the data of Langmuir and Blodgett
lwc	- liquid water content
mvd	- median volume diameter droplet
R	- Macklin correlation parameter for rotating cylinder ice density
R'	- revised Macklin parameter for correlation of ρ_0
R_c	- circular cylinder diameter
Re	- droplet Reynolds number based on the free-stream velocity
Re_d	- droplet Reynolds number based on the relative velocity
r_d	- droplet radius
T	- temperature

Th	- thickness of ice layer
T_{L}	- tail length, added to α_{max}
T_s	- mean surface temperature of ice layer
$T_{s,0}$	- surface temperature of ice at the stagnation line
Δt	- integration time step for droplet trajectories
U	- free-stream air speed
u_x, u_y	- air speed coordinates
v	- relative speed of droplet with respect to the air
v_r	- impact speed of droplet in the direction normal to the cylinder surface
v_0	- droplet impact speed at the stagnation line
v	- absolute velocity of droplet
v_x, v_y	- droplet speed coordinates
x_0, y_0	- initial droplet coordinates
Δy	- initial separation of trajectory pair
$\Delta \ell$	- final separation in arc length of the trajectory pair
α	- angle between surface normal and the free-stream
α'	- scaled surface angle for Equation V.7
α_0	- angular position of maximum β
α_{max}	- angular position of accretion limit
α_m	- accretion limit plus tail length, T_L

- β - local collision efficiency
 β_0 - local collision efficiency at the stagnation line
 ζ - Stallabrass correlation parameter
 θ - position angle on circular cylinder
 μ - dynamic viscosity of air
 ρ - local ice density
 ρ_a - density of air
 ρ_d - density of the water droplet
 ρ_0 - local ice density at the stagnation line
 τ - ice layer time step
 Φ - angle of attack of the free-stream
 ϕ - non-dimensional Langmuir flow parameter, or Stokes number
 ψ - potential flow streamfunction

I. Introduction

'Icing' presents a problem occasionally to pastry cooks, sometimes to hockey players, and always to those concerned with the natural phenomenon of ice accretion.

In cold climates, ice accretion may result from atmospheric cloud droplets, rain, snow, or sea spray. It may cause serious hazards for aircraft, power lines and transmission towers, ships, and off-shore oil rigs. All of these represent human enterprises that involve many lives and dollars, which may be jeopardized during icing events. So, the primary aim in studying icing is the prediction of its occurrence, whether of a particular event or of the probability of extreme events, and ultimately the prevention of costly and often disastrous accidents.

The 'icing' phenomenon may appear straightforward at first glance: water is collected on a cold surface and freezes there. However, a deeper look reveals many layers of complexity in the physical processes involved, and serious practical difficulties in the observational, experimental, and theoretical aspects of the problem. The development of a predictive model is therefore not an easy task, but it is nonetheless one of some importance.

As with other physical problems, there are two approaches that may be taken in model building. The first is a statistical approach; in which a very large number of events (predictands) are observed, as well as their antecedent or concurrent conditions (predictors). A

statistical analysis will then look for correlations between the predictors and predictands. Considerable computational effort may be required to establish a 'best guess' for the most significant predictors, and for the form of the predictive relation.

The reliability of the outcome is highly dependent on the number and quality of the observations. The advantage of this method is that a predictive relation may be derived and applied in the absence of any further understanding of the physical processes involved, which may be complex. But if the observations are adequate, and a good correlation can be found, it is the quickest way of achieving practical forecasting results.

However, the predictive skill of this method is limited by the strength of the discovered correlation, and its applicability is limited to the range of conditions defined by the original observations. For icing studies, this of course includes the geographical range, since the character of icing problems has a large regional variability.

The other approach to modelling is the physical one. This usually means a mathematical model, solved numerically using a digital computer, in which an attempt is made to simulate the intermediate steps, and to understand the causal relationships between the input conditions and the output predictions. Producing such a model does take considerably more time and effort than producing a statistical correlation, and the requirements for the

observational data are just as stringent. Furthermore, no mathematical model can perfectly reproduce a given physical process or situation - there must always be simplifying assumptions.

But the knowledge gained from a physical model is less restricted in its application, within those assumptions, than is any purely statistical relationship, and it can provide a useful testing ground for theory. Even the simplest physical model might at least suggest what the most effective predictors will be for a statistical study; thus if it is found that liquid water content and cloud droplet size are of prime importance in determining model ice accretion rates, then collectors of field data should take note, even if these things are very difficult to observe.

A. The Icing Problem

A model of the complete ice accretion process involves the solution of a series of problems. First, an approximation to the airflow around the (accreting object must be found. If potential flow is assumed, and there is no analytical solution (as there is for circular cylinders or Joukowski airfoils) then the partial differential equation for the velocity potential must be solved numerically, as in Hess and Smith (1967), or Kennedy and Marsden (1976), for example.

Once it is possible to calculate the air velocity at any point, the trajectories of water droplets carried by the

airstream toward the object may be determined. Again, simplifying assumptions are applied to the droplet's equation of motion in order to reduce the computational time needed to integrate it numerically.

An ensemble of droplet trajectories is calculated, and their points of impact on the accretion surface determined. This information is necessary to derive the impingement characteristics of the accretion shape. These include: the distribution of the local collision efficiency, the overall collision efficiency, the maximum impingement angle, and impact speed of the droplets. The collection efficiency is the product of the collision efficiency and the sticking efficiency. Since the latter is unknown, but usually assumed to be unity, only the collision efficiency is dealt with. These quantities and their symbols are defined for a circular cylinder in Figure II.1.

Langmuir and Blodgett (1946) authored the classical solution to this problem, for collision efficiencies of cylinders, spheres, and ribbons in a potential flow field. More recent, but less extensive calculations have been made by McComber and Touzot (1981), and by Egelhofer, et al. (1984). The experimental work of Ranz and Wong (1952) qualitatively validates the theory of Langmuir and Blodgett, if not their numerical calculations.

The impingement characteristics depend on the cylinder size, droplet size, the free-stream air speed, and the air temperature and pressure. If more than one droplet size is

present, the impingement characteristics of each size, or size interval, must be determined. A common simplification to this requirement for droplet spectra is made through the substitution of the impingement characteristics of a single droplet size, that of the median volume droplet of the size spectrum. The usefulness of this simplification is examined in Appendix VI, although it is used throughout the thesis.

Only when all of these calculations are completed can ice begin to accrete in the model. For obvious reasons then, most modelers have turned to standard accretion shapes for which these calculations are simpler, for which much data already exist, and for ease of comparison between models. Thus we have the ubiquitous horizontal circular cylinder in a uniform cross-flow, which appears in the models of Macklin and Payne (1967), McComber and Touzot (1981), Lozowski, Stallabrass and Hearty (1983), Horjen (1983), Egelhofer, et al. (1984) and Makkonen (1984). The more difficult problem of accretion on airfoils has been treated by Oleskiw (1982), Gent and Cansdale (1985), and MacArthur, et al. (1982).

The cloud water is usually assumed to be at the temperature of the air stream, and to arrive at the accreting surface in a uniform flux determined by the liquid water content, or lwc, (defined as the mass of water contained in a unit volume of air), the air speed and the object's collision efficiency. The density of the accreted ice also determines the thickness and shape of the deposit.

On rotating cylinders, accreted rime ice densities have been examined by Macklin (1962), and by Makkonen and Stallabrass (1985). On non-rotating objects, the shape of the deposit is also partly dependent on the variation of the local ice density along its surface. Bain and Gayet (1982) have suggested one form for this variation.

If the flux of super-cooled water is large enough so that the latent heat of freezing cannot be radiated, conducted or convected away from the surface fast enough, then the surface temperature may rise to the melting point. When this happens the ice may become 'wet', rather than 'dry', meaning that all of the impinging water is no longer able to freeze on impact. Now the problem becomes even more complex since a detailed heat balance must be calculated for the accretion surface, in order to determine how much of the impinging water is able to freeze, and how much is shed, or remains in the accretion as liquid water. Calculation of the heat balance is made especially difficult since the heat transfer properties of the ice surfaces and even of rough circular cylinders are very poorly known (Makkonen, 1985).

Models in which a computation of the heat balance on cylinders is attempted using empirical heat transfer data include those of Lozowski, Stallabrass and Hearty (1984), Macklin and Payne (1967), Egelhofer, et al. (1984), and Makkonen (1984). However, such data are scarce, especially for non-cylindrical shapes, and for rough surfaces. Some modelers have thus used numerically calculated heat transfer

for cylinders (Makkonen, 1985) or for airfoils (Cansdale and McNaughton, 1977, Cansdale and Gent 1983).

Models developed specifically for marine applications must contend with the further complications of salinity, droplet distributions and liquid water contents that vary with height, and intermittent sprays. Models including these effects are currently under development by various researchers.

The models developed for this thesis are intended as 'operational' icing models, which would be efficient and inexpensive enough to be used regularly in the laboratory for predictive purposes, or for numerical experiments into specific aspects of icing. Previous models have often been much too costly to be used in such a way. To reduce computational time several 'shortcuts', or parameterizations, have been introduced which help to simplify some of the calculations needed in this problem. Specifically, the local collision efficiency and local ice density are approximated by analytical functions of the surface slope. Such an approach naturally involves some loss of accuracy, but this loss must be balanced against unavoidable uncertainties in the input data and the required precision of output predictions.

Two models are described, for circular cylinders in Chapter III, and in Chapter V for airfoils, that could also be applied to other arbitrary shapes. Both models are limited to rime ice accretion, i.e. there is no heat balance

calculation, and a uniform air speed and flux of pure water droplets are assumed. This is so that the performance of the parameterizations can be easily compared with real ice accretions whose shape and mass are not dependent on the poorly known heat transfer distribution.

In the course of model development, significant improvements were made in two specific aspects, the theoretical calculation of collision efficiencies on smooth circular cylinders, (Chapter III) and the distribution of local density in non-rotating circular cylinder rime accretions (Chapter IV).

Experiments for model verification were carried out in the University of Alberta's original icing wind-tunnel, known as the Facility for Research On Solidification and Thawing, or FROST tunnel. A full description of the facility may be found in Gates (1981), or Lozowski and Gates (1984). All wind-tunnel experiments were performed under the direction of Professor Gates.

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II. Cloud Droplet Trajectories Updated

A. Introduction

For many years the standard reference work for cloud droplet trajectories in an airstream flowing around an obstacle has been Langmuir and Blodgett (1946), hereafter LB. This was one of the first published treatments of the problem which used an analog computer to integrate the equations of motion for water droplets in a steady potential flow about an infinitely long circular cylinder.

These calculations are of interest in studies of atmospheric and marine ice accretion, in which the characteristics of droplet impingement on cylinders and other objects are important factors in modelling the icing process. However, since the calculations were done so long ago, on what would now be considered a primitive machine, and because the original reference can be difficult to obtain, a need was perceived to recalculate the droplet trajectories and to publish the results in the open literature. The effect of replacing the drag coefficient formulation used by LB with a more recent one, that of Beard and Pruppacher (1969), has also been investigated.

Functional fits have been devised by several authors to the original tabulated data of LB. A number of these are compared to show the need for improving and standardizing them. Toward this end, a non-linear regression fit to the new data has been found for the range of conditions

pertaining to airborne, land structure and marine icing.

B. Impingement Parameters and Droplet Trajectories

From trajectory calculations may be derived the overall collision efficiency, E , the local collision efficiency at a point on the surface represented by its local surface slope, $\beta(\alpha)$, the local collision efficiency at the stagnation line, β_0 , the maximum impingement angle on the cylinder surface, α_{\max} , and the non-dimensional impact velocity at the stagnation line, V_0 . V_0 is normalized by the free stream speed, U . The local surface angle, α , is defined as the angle between the local surface normal and the free stream direction. The other quantities are defined in Figure II.1.

Values of E , β_0 , α_{\max} and V_0 are given by LB for a large number of cases which vary in droplet size, cylinder size, and air speed. These results are tabulated in terms of the non-dimensional parameters, K :

$$K = (\rho_d D_d^2 U) / (9 \mu D_c), \quad \text{II.1}$$

and ϕ :

$$\phi = Re^2 / K, \quad \text{II.2}$$

with the droplet Reynolds number based on the free-stream velocity, Re , given by:

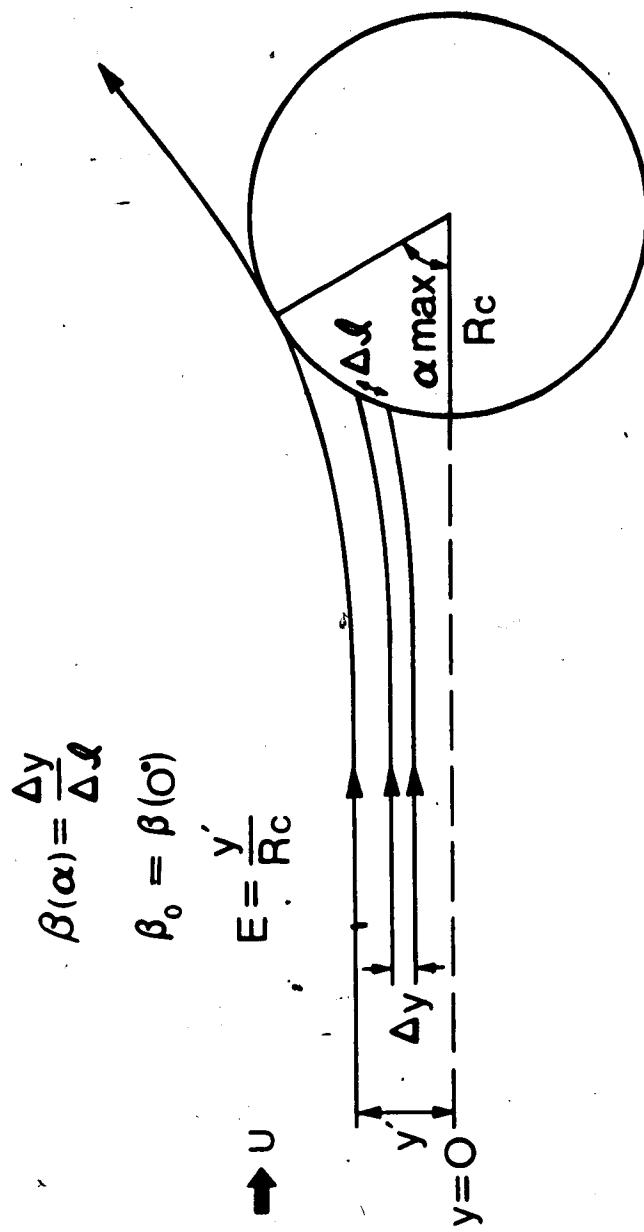


Figure II.1 - Droplet trajectories impinging on a circular cylinder, giving the definitions of $\beta(\alpha)$, β_0 , E and α_{max} . On the accretion surface α is the angle between the local surface normal and the free stream direction.

$$Re = U D_d \rho_a / \mu.$$

II.3

Here ρ_d is the density of water, ρ_a the density of air, μ the dynamic viscosity of air, D_d the droplet diameter, D_c the cylinder diameter, and U the free stream air speed.

The K parameter measures the inertia of the droplet, and is also known as the Stokes number. To aid in solving problems in which the droplet size is the unknown quantity, LB presented their data in terms of a second non-dimensional parameter, ϕ , that is independent of D_d .

The case $\phi = 0$ corresponds to Stokes, or creeping flow:

$$C_d Re_d / 24 = 1,$$

II.4

where C_d is the droplet drag coefficient, and Re_d is the droplet Reynolds number based on the relative velocity with respect to the air, V , so that:

$$Re_d = |V| Re / U.$$

II.5

In component form, the droplet equations of motion are:

$$dv_x / dt = -(v_x - u_x) (U C_d Re_d) / (12 D_c K)$$

$$dv_y / dt = -(v_y - u_y) (U C_d Re_d) / (12 D_c K),$$

where t is time, v_x , v_y and u_x , u_y are the droplet and air

speed components, respectively. This set of equations is in dimensional form, with the origin at the cylinder axis and U in the positive x direction. Gravity and the effects of droplet acceleration on the drag have been neglected, as in the original theory of LB.

The air speed components for potential flow around a cylinder (Green, 1937) are:

$$u_x(x,y) = U \left(1 + (R_c^2 (y^2 - x^2) / (x^2 + y^2)^2) \right)$$

II.7

$$u_y(x,y) = -(2 U R_c^2) (x y) / (x^2 + y^2)^2.$$

In order to permit a direct comparison with the original solutions by LB some trajectories have been integrated using their empirical formulation for the steady-state drag coefficient:

$$C_d R_{ed} / 24 = 1 + 0.197 R_{ed}^{0.63} + 2.6 \cdot 10^{-4} R_{ed}^{1.38} \quad II.8$$

More recent studies have improved the accuracy of C_d approximations at low Reynolds numbers, however, so in further calculations over the complete range of conditions the formulation of Beard and Pruppacher (1969) has been substituted when R_{ed} is less than 200. This formulation is:

$$C_d R_e / 24 = 1 + 0.102 R_e^{0.955}, \text{ for } 0.2 \leq R_e \leq 2.0$$

$$C_d R_e / 24 = 1 + 0.115 R_e^{0.802}, \text{ for } 2.0 \leq R_e \leq 21.0 \quad \text{II.9}$$

$$C_d R_e / 24 = 1 + 0.189 R_e^{0.632}, \text{ for } 21.0 \leq R_e \leq 200.0$$

A comparison of a variety of theoretical and experimental results for the drag coefficients of spheres is given by Beard (1976), in which the above formulation is shown to be an accurate analytical approximation.

The integration is carried out by the Heun method as described in Mesinger and Arakawa (1976). The droplet trajectories are begun at ten cylinder radii upstream of the origin with speeds at the initial position (x_0, y_0) assumed to be:

$$v_x(x_0, y_0) = u_x(x_0, y_0) \quad \text{II.10}$$

$$v_y(x_0, y_0) = 0.5 u_y(x_0, y_0)$$

At this initial upstream distance, the final results are not greatly influenced by the choice of initial droplet speed as long as it lies in between the local air speed and the undisturbed upstream air speed. Tests indicate differences in collision efficiency of less than 0.5 percent, when the initial droplet speeds are varied within this range.

Other factors which influence the accuracy of the integration are the time step, Δt , and the initial upstream distance, x_0 . A value of $x_0 = 10$ cylinder radii has been chosen, which, when compared to the results with $x_0 = 20$ cylinder radii, agrees to within 0.7 percent. At this initial distance, the size of the time step made less than a 0.5 percent difference in the results, as long as $\Delta t < 0.0033 (D_c/U)$. The value used is 0.0025 (D_c/U). Double precision is used throughout the program.

Values of the air viscosity and air density used are based on an assumed air temperature of -10° C. Within the temperature range of interest for most dry icing, i.e. -20 to 0° C, the collision efficiencies vary by a maximum of ± 1 percent.

An air pressure of 100 kPa was assumed. The calculations of LB were carried out for an atmospheric pressure of 78.5 kPa, for application to Mt. Washington, or to airborne icing in general. At this pressure value, the present results differed from the 100 kPa results by less than 0.5 percent. The new results should therefore be applicable to both ground-level and airborne icing.

Two complete trajectories must be integrated in order to calculate one value of the local collision efficiency $\beta(\alpha)$, where α is the mean angle of the two impact points. As illustrated in Figure II.1, $\beta(\alpha)$ is defined as the ratio of the initial vertical separation of the two trajectories, to the final separation in arc length of the two impact

points. In the present model, the initial y separation of the two trajectories is always 10^{-8} metres, an arbitrarily small distance for use in the finite difference approximation to $dy/d\zeta$, the definition of local collision efficiency. The stagnation line value, β_0 , is calculated from the case where the first trajectory's initial position $y_0 = 0$.

The maximum impingement angle, α_{\max} , is found from the droplet trajectory which is tangent to the cylinder surface at its point of impact. This 'grazing' trajectory is found by increasing the initial position y_0 by successively smaller amounts until the tangent trajectory is reached.

The total collision efficiency E may then be calculated from the ratio of the initial vertical distance $y_0(\text{graze})$ to the cylinder radius.

The Fortran77 code written for these calculations is included in Appendix I.

C. Results

In Figure II.2 three curves of β_0 vs. K are shown for the value $\phi = 10^3$, to illustrate the differences between LB's original calculations, the re-calculation of their results, and the re-calculation using Beard and Pruppacher's formulation for drag coefficient. The data for this Figure are listed in Table II.1.

A comparison between the first two curves shows that the LB results are too large for small K ($K \leq 0.4$), by up to

K	β_0 (LB), old C_d	β_0 (new), old C_d	β_0 (new), new C_d
.144	0.048	0.007	0.011
.196	0.095	0.058	0.073
.256	0.146	0.116	0.137
.4	0.233	0.224	0.247
.625	0.323	0.335	0.359
.9	0.398	0.425	0.447
1.6	0.513	0.558	0.575
3.6	0.680	0.714	0.724
6.4	0.778	0.798	0.806
10.	0.837	0.848	0.854
19.6	0.901	0.902	0.906
32.4	0.932	0.929	0.932
62.5	0.959	0.953	0.955
90.	0.969	0.962	0.963
160.	0.980	0.972	0.973
360.	0.989	0.981	0.981
640.	0.993	0.984	0.984
1000.	0.995	0.986	0.986

Table II.1 - $\beta_0(K)$ for $\phi = 10^3$. A comparison between Langmuir and Blodgett's results and the new results.

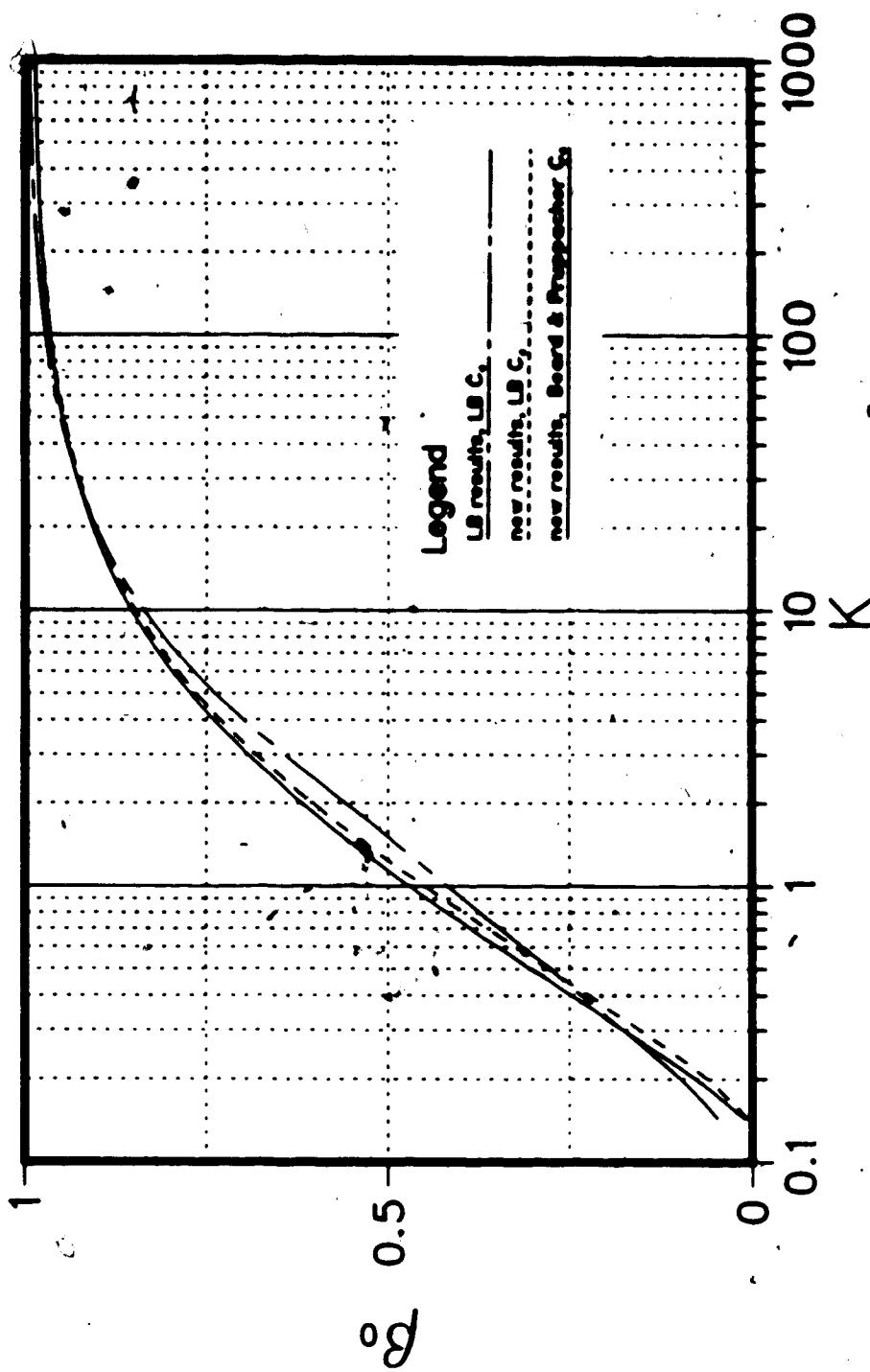


Figure II.2 - Comparison of β_0 vs. K for one value of $\phi = 10^3$.

several hundred percent. In the range $0.4 < K < 20.0$ their results are too small, by amounts near 10 percent in the middle of the range. For larger K , their values are again larger than the new results, but only by a few percent. The use of the revised drag coefficients (dotted curve, Figure II.2) gives another small increase over the LB results for all $K < 200.0$. A noticeable difference is that the new results for β_0 reach an effective value of zero for $K \approx 0.14$, rather than at LB's theoretical value of 0.125. This difference is as yet unexplained. Results for other ϕ values, and for α_{\max} , E and V_0 , are quantitatively similar in differences between the old and new results.

It is concluded from the comparisons in Figure II.2 and Table II.1, that the original LB calculations are accurate only for very large K , but that when K is less than about 200 they may be significantly in error. The new results are suggested for future use. These are shown for β_0 , α_{\max} , E and V_0 for a complete range of ϕ and K in Tables II.2 to II.3, and in Figures II.3 through II.6, where the tabulated data have been fit by spline curves.

D. Curve Fitting

LB also drew smooth curves through their data, and derived analytical expressions in K and ϕ to fit them. Subsequently other authors have made attempts to improve on or to simplify these relations, and have then applied them

Table II.2 - Numerically calculated values of β_0 , α_{\max} and E as a function of K and φ

K	$\varphi = 0$			$\varphi = 10$			$\varphi = 10^3$			$\varphi = 10^5$		
	β_0	α_{\max}	E	β_0	α_{\max}	E	β_0	α_{\max}	E	β_0	α_{\max}	E
144	0.014	0.018	0.0002	0.013	0.009	0.0001	0.011	0.015	0.001	0.004	0.007	0.0002
196	0.094	0.170	0.012	0.091	0.159	0.011	0.073	0.131	0.007	0.030	0.054	0.001
256	0.175	0.328	0.043	0.170	0.310	0.040	0.137	0.247	0.025	0.058	0.105	0.004
4	0.313	0.580	0.131	0.304	0.553	0.122	0.247	0.436	0.079	0.113	0.195	0.016
625	0.446	0.799	0.249	0.432	0.760	0.233	0.359	0.612	0.157	0.179	0.302	0.039
9	0.545	0.953	0.5	0.530	0.905	0.333	0.447	0.742	0.234	0.244	0.400	0.070
1.6	0.680	1.150	1.1	0.663	1.095	0.492	0.575	0.921	0.365	0.353	0.559	0.142
3.6	0.821	1.345	0.718	0.805	1.298	0.686	0.725	1.136	0.553	0.508	0.764	0.274
6.4	0.886	1.433	0.818	0.872	1.383	0.789	0.806	1.255	0.670	0.601	0.898	0.376
10.	0.921	1.476	0.873	0.909	1.439	0.848	0.854	1.329	0.745	0.670	0.997	0.459
19.6	0.954	1.521	0.927	0.945	1.489	0.908	0.906	1.413	0.832	0.759	1.130	0.580
32.4	0.968	1.533	0.951	0.961	1.513	0.936	0.932	1.457	0.878	0.812	1.214	0.661
62.5	0.979	1.552	0.970	0.974	1.532	0.959	0.955	1.496	0.920	0.866	1.303	0.749
90.	0.983	1.556	0.976	0.979	1.539	0.967	0.963	1.512	0.935	0.889	1.347	0.789
160.	0.986	1.563	0.982	0.984	1.545	0.976	0.973	1.530	0.953	0.917	1.400	0.840
360.	0.989	1.567	0.987	0.987	1.553	0.983	0.981	1.546	0.969	0.945	1.454	0.892
640.	0.990	1.568	0.988	0.989	1.559	0.986	0.984	1.553	0.976	0.957	1.481	0.917
1000.	0.990	1.568	0.989	0.989	1.557	0.987	0.986	1.557	0.980	0.965	1.499	0.932

K	$\varphi = 10^2$			$\varphi = 10^4$			$\varphi = 10^6$		
	β_0	a_{\max}	E	β_0	a_{\max}	E	β_0	a_{\max}	E
.16	0.034	0.054	0.001	0.021	0.035	0.0006	0.005	0.011	0.00004
.25	0.153	0.282	0.032	0.094	0.168	0.012	0.026	0.051	0.001
.36	0.257	0.466	0.088	0.164	0.287	0.034	0.050	0.089	0.003
.64	0.415	0.720	0.212	0.281	0.468	0.095	0.103	0.176	0.013
1.	0.528	0.889	0.325	0.374	0.609	0.163	0.157	0.260	0.029
1.96	0.676	1.097	0.500	0.518	0.817	0.296	0.251	0.380	0.067
3.24	0.763	1.219	0.617	0.617	0.958	0.406	0.319	0.488	0.112
6.25	0.848	1.336	0.743	0.726	1.115	0.547	0.424	0.633	0.191
9.	0.882	1.387	0.797	0.777	1.187	0.618	0.482	0.712	0.243
16.	0.921	1.446	0.863	0.838	1.278	0.711	0.570	0.832	0.333
36.	0.955	1.501	0.922	0.899	1.379	0.812	0.680	0.982	0.463
64.	0.968	1.523	0.946	0.927	1.434	0.864	0.745	1.080	0.550
100.	0.975	1.536	0.959	0.944	1.465	0.894	0.787	1.143	0.612
196.	0.982	1.548	0.972	0.960	1.497	0.927	0.839	1.227	0.693
324.	0.985	1.554	0.978	0.969	1.517	0.944	0.869	1.282	0.744
625.	0.987	1.558	0.983	0.977	1.533	0.959	0.900	1.337	0.799
900.	0.988	1.559	0.985	0.980	1.538	0.966	0.913	1.366	0.824

Table II.2 - continued.

K	v_0				v_0			
	$\varphi = 0$	$\varphi = 10$	$\varphi = 10^3$	$\varphi = 10^5$	$\varphi = 10^2$	$\varphi = 10^4$	$\varphi = 10^6$	
144	0.0002	0.0002	0.0002	0.0001	.16	0.002	0.0009	0.0001
196	0.018	0.017	0.012	0.003	.25	0.051	0.021	0.003
256	0.065	0.061	0.041	0.010	.36	0.133	0.059	0.010
4	0.195	0.183	0.122	0.033	.64	0.297	0.144	0.029
625	0.348	0.327	0.227	0.068	1.	0.428	0.228	0.055
9	0.471	0.443	0.317	0.111	1.96	0.604	0.379	0.116
1.6	0.638	0.604	0.459	0.249	3.24	0.708	0.491	0.162
3.6	0.804	0.772	0.640	0.346	6.25	0.810	0.625	0.251
6.4	0.879	0.852	0.742	0.447	9.	0.853	0.690	0.306
10.	0.918	0.896	0.806	0.528	16.	0.904	0.768	0.396
19.6	0.956	0.940	0.876	0.642	36.	0.948	0.854	0.520
32.4	0.973	0.960	0.913	0.715	64.	0.967	0.897	0.603
62.5	0.986	0.977	0.945	0.793	100.	0.977	0.922	0.659
90.	0.990	0.983	0.957	0.828	196.	0.986	0.948	0.732
160.	0.994	0.989	0.971	0.872	324.	0.990	0.962	0.777
360.	0.997	0.995	0.983	0.912	625.	0.994	0.975	0.826
640.	0.999	0.997	0.989	0.936	900.	0.996	0.980	0.849
1000.	0.999	0.998	0.992	0.950				

Table II.3 - Numerically calculated values of v_0 as a function of K and φ .

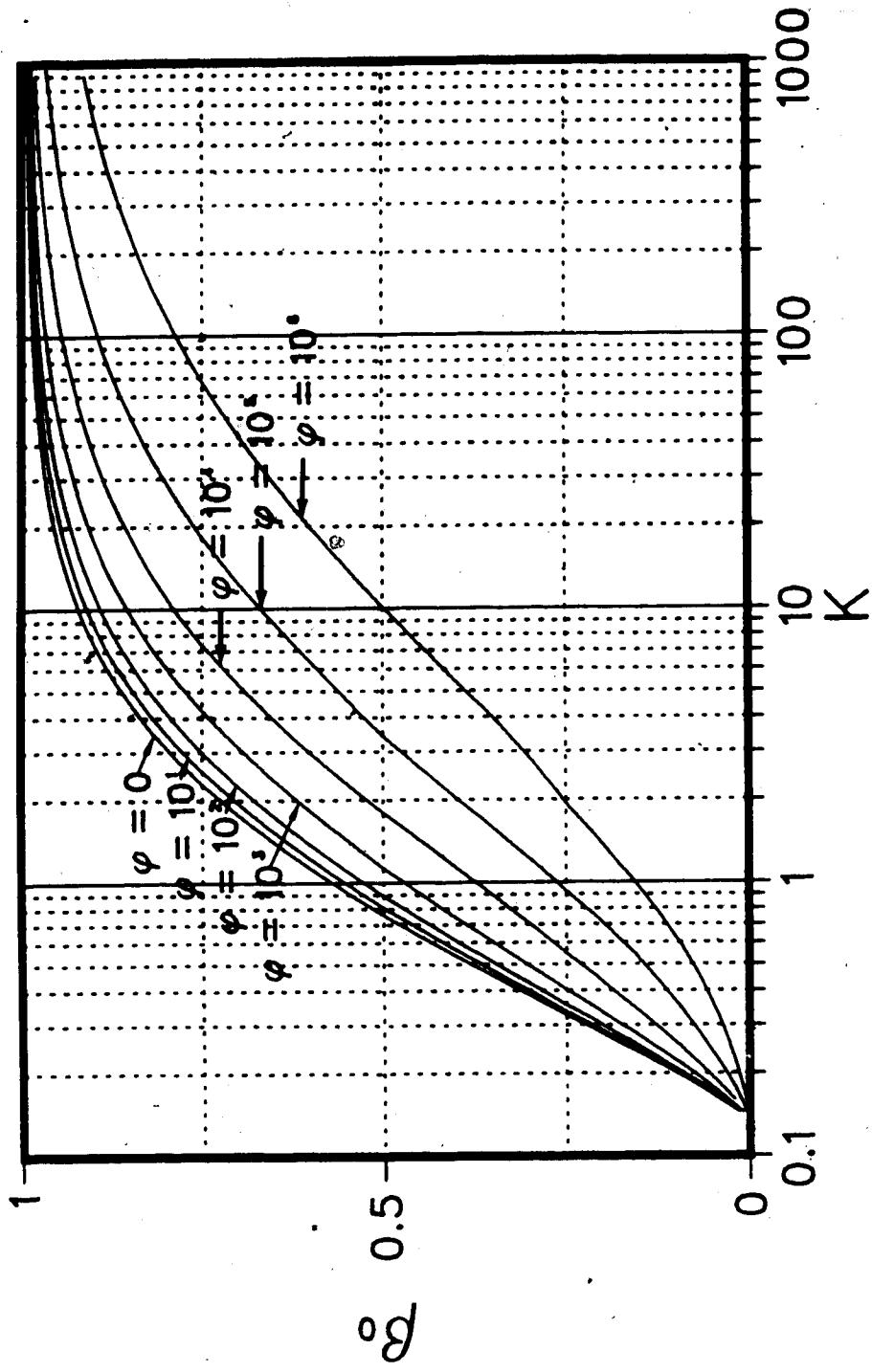


Figure II.3 - Smooth curves drawn for the numerical data of β_0 vs. K and φ .

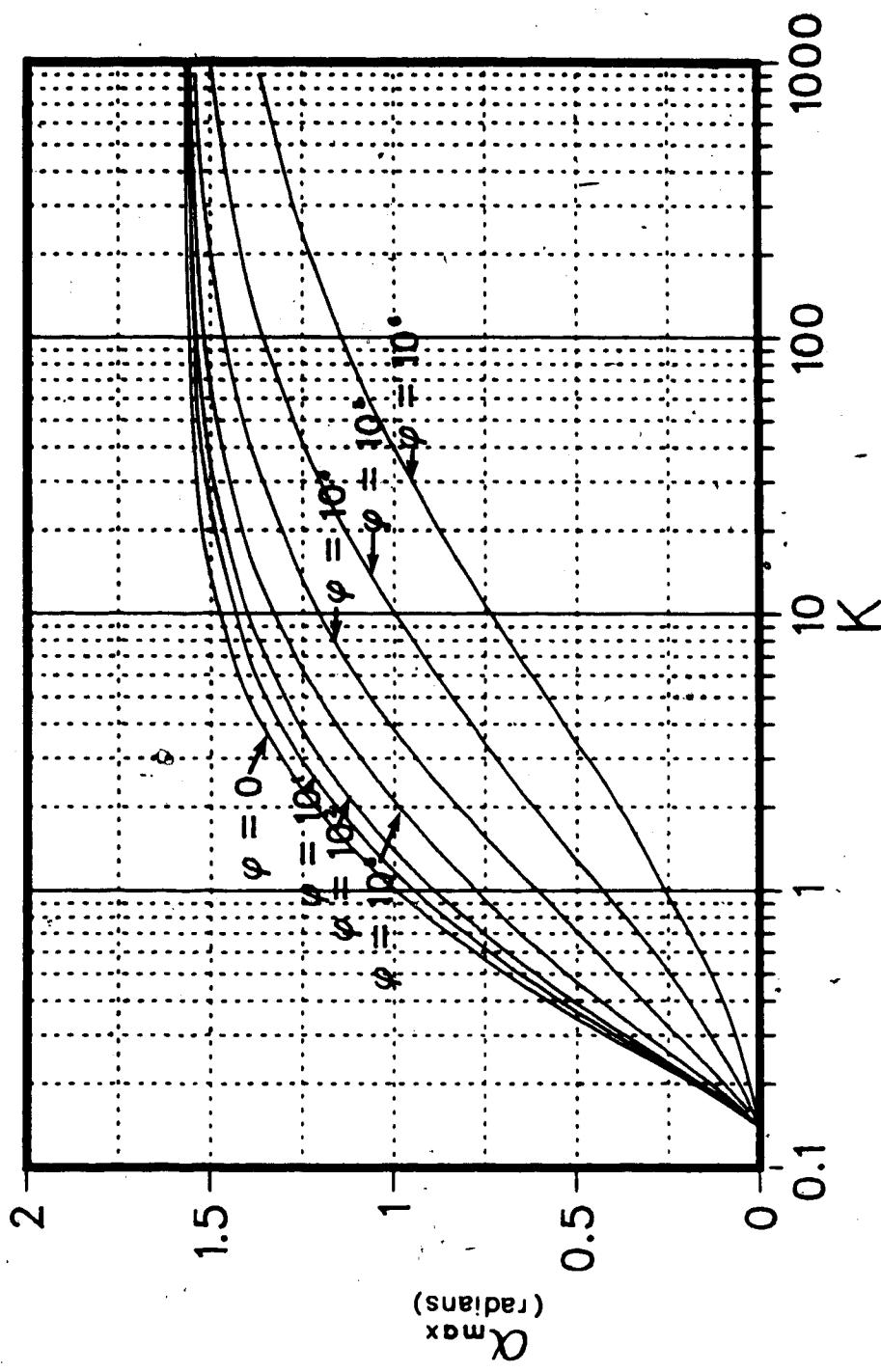


Figure II.4 - Smooth curves drawn for the numerical data of α_{\max} vs. K and φ .

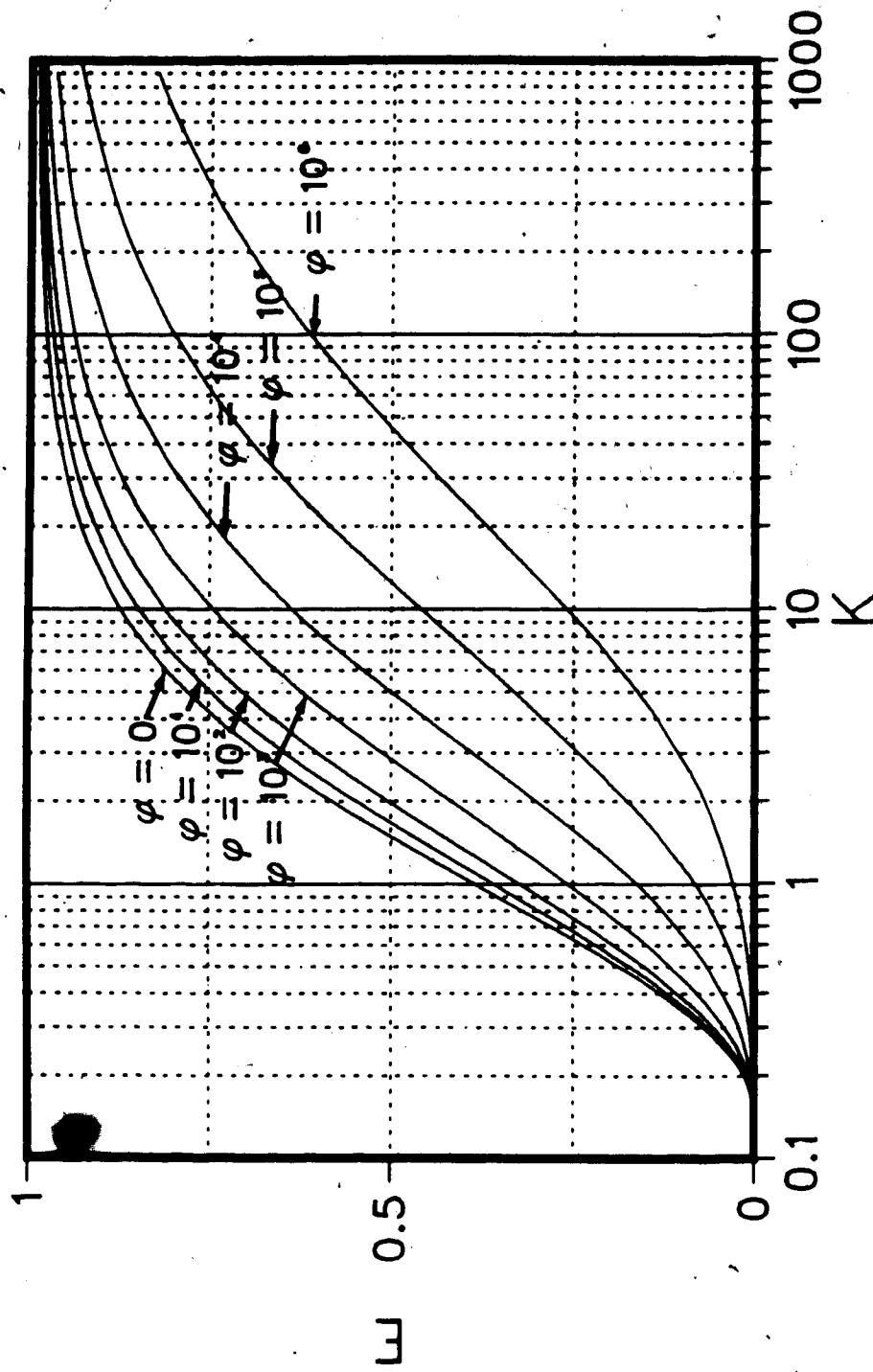


Figure II.5 - Smooth curves drawn for the numerical data of E vs. K and φ .

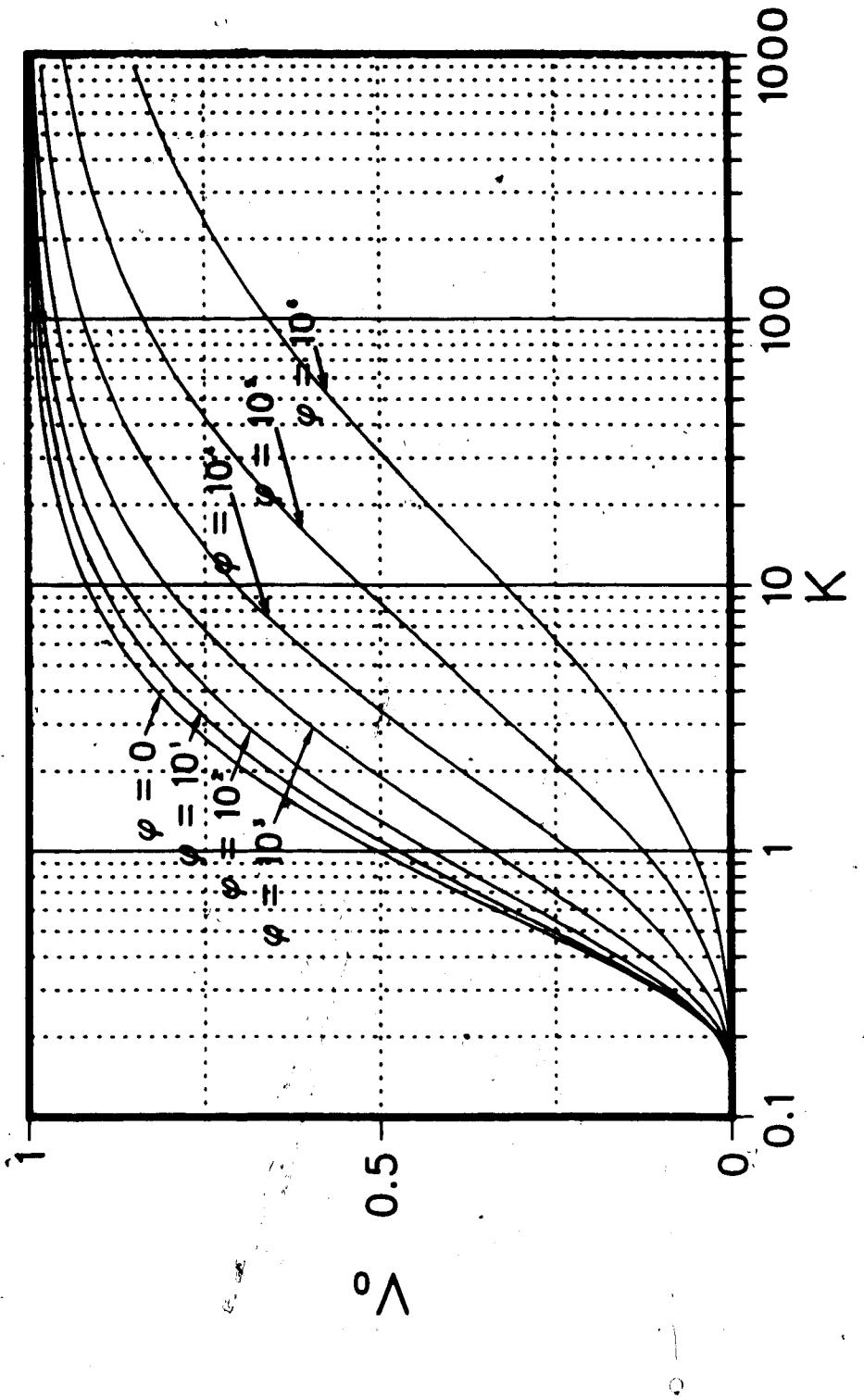


Figure III.6 - Smooth curves drawn for the numerical data of V_0 vs. K and ϕ .

Author (s)	$K_o (K, Re), \xi (U, D_g, D_c)$	$E (K_o), E (K, Re), E (\delta)$	Range
Langmuir and Blodgett	$K_o = .125 + \frac{(K - .125)}{1 + .0967 Re \cdot 6367}$	$E = .466 (\log_{10} (8K_o))^2$ $E = \frac{K_o}{K_o + \frac{\xi}{2}}$ $E = \frac{K}{K + \frac{\xi}{2} \cdot 112 Re^{.63} + .75 \cdot 10^{-4} Re^{1.38}}$.125 < $K_o < 1.1$ $K_o > 1.1$ $E > 0.5$
Lozowski, Stalabass and Hearty	$K_o = .125 + \frac{(K - .125)}{1 + .0967 Re \cdot 6367}$	$E = .489 (\log_{10} (8K_o))^{1.978}$ $E = \frac{K_o}{K_o + \frac{\xi}{2}}$.125 < $K_o < 0.9$ $K_o > 0.9$
Cansdale and McNaughton	$K_o = \frac{K}{1 + .087 Re (.76 Re^{-.027})}$	$E = .53 (\log_{10} (8K_o))^{1.8}$ $E = \frac{K_{1.1}}{K_{1.1} + 1.223}$.125 < $K_o < 1.1$ $K_o > 1.1$
Makkonen	$K_o = \frac{K}{1 + .087 Re (.76 Re^{-.027})}$	$E = .5 (\log_{10} (8K_o))^{1.6}$ $E = \frac{K_{1.1}}{K_{1.1} + 1.426}$.125 < $K_o < 0.8$ $K_o > 0.8$
Stalabass	$\xi = \frac{U^{.6} D_d^{1.6}}{D_c}$	$E = \frac{\xi - 3200}{\xi + 27000}$ $E = 0,$	$\xi > 3200$ $\xi < 3200$

Table II.4 - Approximations to the numerical data of Langmuir and Blodgett derived by various authors.

to various numerical icing models (Cansdale and McNaughtan, 1977, Stallabrass, 1980, Lozowski; Stallabrass and Hearty, 1983, Makkonen, 1984). A number of these fitted relations for $E(R, \phi)$ are listed in Table II.4.

In order to compare these relations, values computed from each are listed with a selection of the original data in Table II.5. The data all fall within a range of ϕ between 10^2 and 10^4 , since some of the fitting functions were specifically designed for similar ranges, and because this range should apply to most physical applications, corresponding to $10 \leq D_d \leq 4000$ microns, $0.01 \leq D_c \leq 1.0$ metre, $5 \leq U \leq 200$ metres per second and $-20^\circ\text{C} \leq T \leq 0^\circ\text{C}$. It is worth noting that the natural range is actually narrower than this, since the largest drops do not occur together with the highest wind speeds (e.g. marine icing), nor do the smallest droplets occur at the lowest speeds (e.g. aircraft icing).

Langmuir and Blodgett's own fit is the most accurate, since it includes a correction table for $E < 0.5$, and for larger values of R , (giving E values > 0.5), they give theoretically derived functions for β_0 , a_{\max} , and E , which assume essentially straight-line trajectories. For this more elaborate scheme, the standard error (computed with respect to the numerical data) is generally an order of magnitude smaller than it is for the other fitting schemes. Even more important, however, is the fact that the variation among the values calculated by the various empirical

formulae for a given case can exceed 10 percent. These discrepancies will add to the difficulty of making comparisons among the icing predictions made by the various models which use these relations.

To reduce any further confusion, the new data have been fitted with approximating functions of improved form and greater accuracy. These functions are described below.

The approach taken by other authors to derive the formulae of Table II.4 has been to collapse a family of curves, such as that in Figure II.3, into a single curve via the ratio K_0/K , calculated as a function of Re , or in the case of Stallabrass (1980), via the parameter ξ , a function of D_d , U , and D_c (see Table II.4). K_0 is the K value read from the $\phi = 0$ curve which yields the same value of β_0 , a_{\max} , E or V_0 as do the original K and ϕ . Extensive manipulation of the new data has revealed major problems with this approach. Specifically, the three families of curves for β , a_{\max} and E are not collapsible by the same ratio K_0/K , but require three different ratios. Only LB take this into account for small values of E , by using their correction table, and theoretical approximations for larger E . Furthermore, the K_0/R 's are not single-valued functions of Re (see Figure II.7), as has usually been assumed.

In the Stallabrass (1980) method, although ξ is used instead of K and ϕ , a similar problem arises for his curves, since they collapse to a single curve only for very large values of ξ .

φ	K	LB (numerical)	LB (approx.)	Lozowski, et al.	Makkonen	Cansdale	Stallabrass
10^2	.16	0.004	0.004	0.004	0.002	0.005	0.056
	.64	0.181	0.183	0.181	0.191	0.202	0.284
	3.24	0.591	0.588	0.568	0.630	0.594	0.634
	16.	0.855	0.854	0.836	0.884	0.867	0.867
	100.	0.966	0.965	0.958	0.996	0.972	0.966
	625.	0.992	0.992	0.989	0.995	0.994	0.992
	10 ³	0.196	0.009	0.010	0.0	0.0	0.017
	.625	0.127	0.129	0.132	0.120	0.134	0.170
	3.6	0.542	0.540	0.509	0.560	0.522	0.536
	19.6	0.830	0.825	0.796	0.852	0.831	0.827
10^4	90.	0.942	0.939	0.925	0.956	0.949	0.943
	640.	0.986	0.985	0.942	0.992	0.990	0.988
	.25	0.012	0.013	0.014	0.0	0.0	0.0
	1.	0.148	0.142	0.146	0.120	0.134	0.152
	6.25	0.546	0.539	0.499	0.557	0.518	0.531
	36.	0.820	0.811	0.778	0.850	0.829	0.831
	196.	0.937	0.934	0.921	0.960	0.954	0.951
	900.	0.974	0.974	0.972	0.988	0.987	0.985
standard error:							
0.003 0.018 0.017 0.010 0.019							

Table II.5 - Comparison of some original E values calculated numerically by Langmuir and Blodgett, and those derived from the approximations of Table II.4.

For these reasons, an accurate fit to the entire range of data given is difficult to find and awkward to use. Consequently, an approximating function is given here only for ϕ between 10^2 and 10^4 , which is the range applicable to most atmospheric, land-based and marine icing problems.

A non-linear regression analysis on this range gives the following result for $0.17 \leq K \leq 10^3$:

$$\begin{aligned} z(K, \phi) = & (C_{X,1} K^{C_{X,2}} \exp(C_{X,3} K^{C_{X,4}} + C_{X,5})) \\ & - (C_{X,6} (\phi - 100)^{C_{X,7}}) \\ & (C_{X,8} K^{C_{X,9}} \exp(C_{X,10} K^{C_{X,11}} + C_{X,12})), \end{aligned} \quad \text{II.11}$$

where X is either β_0 , a_{\max} , E or V_0 , and the constants $C_{X,n}$ are as listed in Table II.6. The basic form of this function is a standard fitting function supplied by the BMDP non-linear regression package (Dixon, et al., 1981).

Values for $K < 0.17$ and $K > 10^3$ may be read from the curves or roughly approximated as follows:

$$\beta_0 = a_{\max} = E = V_0 = 0.0, \text{ for } K < 0.17, \quad \text{II.12}$$

and

$$\beta_0 = E = V_0 = 0.99, \text{ for } K > 10^3, \quad \text{II.13}$$

$$a_{\max} = 1.56, \text{ for } K > 10^3.$$

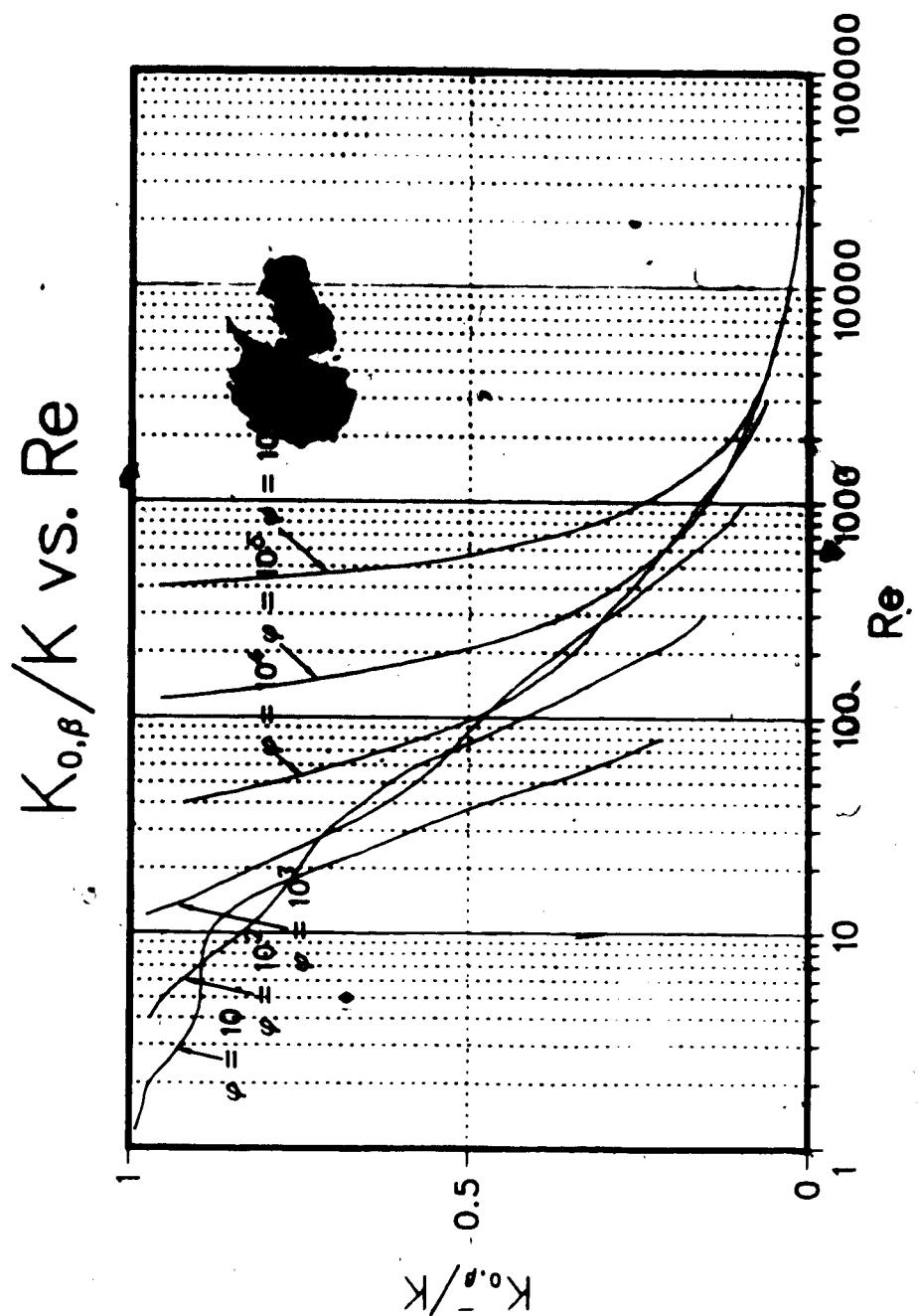


Figure 14.7 The dependence of $\frac{K_{0,\beta}}{K}$ on Reynolds Number Re . $K_{0,\beta}$ is defined by: $\beta_0(K,\varphi) = \beta_0(K_{0,\beta}, 0)$

	$X = \beta_0$	$X = \alpha_{\max}$	$X = E$	$X = V_0$
$c_{X,1}$	1.218	2.433	1.066	1.030
$c_{X,2}$	$-6.70 \cdot 10^{-3}$	$-4.70 \cdot 10^{-3}$	$-6.16 \cdot 10^{-3}$	$1.68 \cdot 10^{-3}$
$c_{X,3}$	-0.551	-0.375	-1.103	-0.796
$c_{X,4}$	-0.643	-0.576	-0.688	-0.780
$c_{X,5}$	-0.170	-0.781	-0.028	-0.040
$c_{X,6}$	$3.05 \cdot 10^{-3}$	$8.50 \cdot 10^{-3}$	$6.37 \cdot 10^{-3}$	$9.44 \cdot 10^{-3}$
$c_{X,7}$	0.430	0.383	0.381	0.344
$c_{X,8}$	2.220	1.757	3.641	2.657
$c_{X,9}$	-0.450	-0.298	-0.498	-0.519
$c_{X,10}$	-0.767	-0.420	-1.497	-1.060
$c_{X,11}$	-0.806	-0.960	-0.694	-0.842
$c_{X,12}$	-0.068	-0.179	-0.045	-0.029

Table 11.6 Coefficients for approximating impingement parameters by the expression:

$$x(K, \varphi) = (c_{X,1} K^{c_{X,2}} \exp(c_{X,3} K^{c_{X,4}}) + c_{X,5}) \\ (c_{X,6}(\varphi - 100)^{c_{X,7}}) (c_{X,8} K^{c_{X,9}} \exp(c_{X,10} K^{c_{X,11}}) + c_{X,12})$$

Within the range of $0.17 < K < 10^3$ the standard error of the fit with respect to the numerical data is given in Table II.7. It compares very favorably to the errors given in Table II.5 for the fitting schemes previously discussed.

E. Limitations of LB Theory

The collision efficiency theory of Langmuir and Blodgett assumes that the air is in potential flow about the collecting cylinder, that only steady-state drag forces are exerted on the water droplets by the flow, and that the flow itself remains unaffected by the presence of the droplets.

The potential flow field about an infinitely long, circular cylinder is as described by Equations II.7^a, and illustrated in Figure II.8(a). Figure II.8(b) illustrates the analogous situation for viscous flow at a cylinder Reynolds number of $\approx 10^4$, corresponding to $D_c = 0.0254$ m, $U = 10$ m/sec, and $T = -10^\circ$ C. The upstream flow in Figure II.8(b), which determines the accretion, deviates slightly from the potential flow case due to the presence of a sizable wake on the downstream side. The cylinder plus wake will be felt by the flow further upstream, decreasing the streamline curvature, and thus increasing the collision efficiency. In most situations, however, the effect will be small.

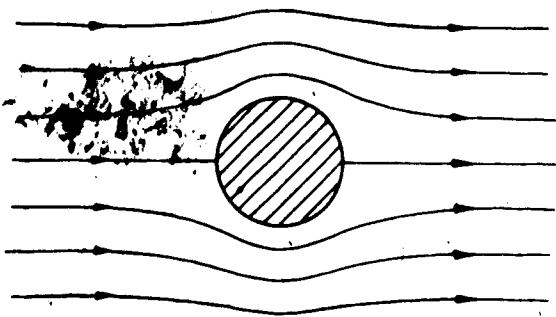
The potential flow assumption also ignores the cylinder boundary layer, which will be of negligible depth for a

numerical approximated

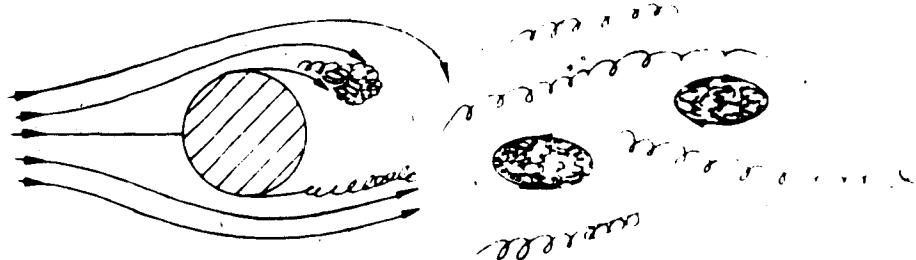
φ	β_0	α_{\max}	E	v_0^c	β_0	α_{\max}	E	v_0
10^2	.16	0.034	0.054	0.001	0.002	0.036	0.054	0.0
	.64	0.415	0.720	0.212	0.297	0.416	0.720	0.210
	3.24	0.763	1.219	0.617	0.708	0.763	1.219	0.619
	16.	0.925	1.446	0.863	0.904	0.920	1.445	0.861
	100.	0.975	1.536	0.959	0.977	0.978	1.538	0.961
	625.	0.987	1.558	0.983	0.994	0.986	1.558	0.983
10^3	.196	0.073	0.131	0.007	0.012	0.075	0.135	0.005
	.625	0.359	0.612	0.157	0.227	0.363	0.611	0.159
	3.6	0.725	1.136	0.553	0.640	0.728	1.139	0.557
	19.6	0.906	1.413	0.832	0.876	0.904	1.401	0.825
	90.	0.963	1.512	0.935	0.957	0.964	1.503	0.931
	640.	0.984	1.553	0.976	0.989	0.983	1.549	0.974
10^4	.25	0.094	0.168	0.012	0.015	0.098	0.179	0.012
	1.	0.374	0.609	0.163	0.228	0.379	0.610	0.162
	6.25	0.726	1.115	0.547	0.625	0.726	1.115	0.548
	36.	0.899	1.379	0.812	0.854	0.899	1.380	0.816
	196.	0.960	1.497	0.927	0.948	0.962	1.497	0.930
	900.	0.980	1.538	0.966	0.980	0.980	1.543	0.967

standard error: 0.002 0.005 0.003 0.006

Table II.7 - Comparisons between a selection of numerically calculated values, and approximated values calculated according to Table II.6.



(a)



(b)

Figure II.8 - (a):Potential flow streamlines about a cylinder in cross-section. **(b):**Viscous flow streamlines about a cylinder at a Reynolds number of 10,000.

smooth cylinder, but may become significant for rough surfaces. The effect, if any, of a deep, possibly turbulent boundary layer on the collision efficiency is unknown. In particular, the onset of accretion may be affected if the roughness elements are large enough to act as individual collectors.

The assumption of the airflow being undisturbed by the droplets carried in it, is a reasonable one given typical cloud liquid water contents.

In determining the deceleration of the drop by the flow, LB's steady-state viscous drag coefficient ignores buoyancy and gravitational forces, the influence of vorticity diffusion from the accelerating droplet surface, and the induced mass increase of the drop as a result of accelerating the air immediately surrounding it.

A 20 micron droplet falls in air with a terminal speed of 0.012 m/sec (Mason, 1971). Thus the vertical deviation of such a droplet approaching a 0.0254 m diameter cylinder from five diameters upstream, at a free-stream speed of 10 m/sec, is about 0.15 mm. This is only about one percent of the maximum vertical deviation due to the flow, under those conditions. At some point on the cylinder surface, the deviation due to the flow will be comparable to the gravitational drift, but since the relative difference in two adjacent trajectories will be much smaller, the collision efficiencies should not be affected.

Much larger drops, such as occur in sea spray icing, will have a correspondingly larger gravitational drift. Preliminary numerical calculations by Szilder (personal communication, 1986) indicate that there is still a very small effect on the collision efficiency, but this is a question which should be investigated further.

The other two effects on the drag are incorporated into the 'history' term, which, expressed as an acceleration, is:

$$-\frac{18\rho_a}{(2\rho_d + \rho_a) D_d} \sqrt{\frac{\mu}{\pi\rho_a}} \int_{-\infty}^t \frac{dv}{d\tau} \frac{d\tau}{\sqrt{t-\tau}} \quad II.14$$

The relative importance of the history term is often measured by the acceleration modulus (Norment, 1980):

$$N_A = D_d |\frac{dv}{dt}| / v^2, \quad II.15$$

where v is the droplet's absolute velocity. If this quantity is < 0.01 , the history term may be safely omitted from the droplet's equation of motion.

Numerical results from the trajectory integration model presented above show that N_A is well within this limit under most conditions of accretion on cylinders. However, for $K \leq 0.5$, N_A may reach values ≥ 0.01 just before impact, and for $K \leq 0.20$, the limit is exceeded as much as a few millimetres in front of the cylinder surface.

For these small K values then, the effect of the history term is to decrease the droplet's deceleration and

increase both its total velocity near the surface, and the resulting collision efficiency. Numerical integrations by Oleskiw (1982) for $K = 0.196$, $\phi = 1000$ show an increase in E of 350%, and in β_0 of 167% when the history term is included.

Thus the 'cut-off' K value, for zero collision efficiency, which arises from LB theory, is not valid for real flow. In light of these results, it may be advisable to consider Equations II.11, II.12, and II.13 as valid only for $K > 0.25$. Below this limit, the trajectories should be recalculated using the complete drag terms in the droplet equation of motion.

The last assumption listed, that of an airflow undisturbed by the presence of water droplets, is probably a reasonable one for the small liquid water contents typical of clouds, $< 5 \text{ g/m}^3$. The case for heavy spray icing may, however, be very different, and this is another question which warrants further study.

F. Summary

Earlier results for water droplet trajectories at freezing temperatures have been re-examined and improved upon. The new results differ from LB by as much as 10 percent. Tabulated data are provided to replace those of Langmuir and Blodgett for the collisional parameters of water droplets on cylinders, and a good approximating function to the data has been found which will be of use to

modellers. These results should only be applied, however, to real situations when $K > 0.25$, below which the non-steady-state drag terms have a significant effect on collision efficiencies.

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III. An Operational Model for Rime Ice Accretion

A. Introduction

Modelling of the many physical processes involved in ice accretion can be a complex and time-consuming exercise. The cost of running such models tends to limit their use as investigative tools, or for predictive purposes of an operational character. In an attempt to alleviate part of this problem, simple parameterization methods have been developed which make possible a more practical, inexpensive and realistic model. In this chapter these methods, and their application in a computer simulation of rime icing on circular cylinders will be described.

The computer model was developed with several specific purposes in mind. The first was to reduce the computational time necessary, in particular, for the calculation of impingement parameters. The quantities E , β_0 and α_{\max} for circular cylinders are available from the tabulated data and analytical fitting functions of Chapter II. Values of local collision efficiency between the stagnation line and maximum impingement angle usually have to be determined from a series of droplet trajectory pairs.¹ Oleskiw (1982) has calculated the distributions of local collision efficiency

¹ It has been suggested (McComber and Touzot, 1981) that the variation of the local radial velocity of impact be used as an approximation to the local collision efficiency. It has been previously shown (Lozowski, Finstad and Gates, 1985) that for realistic cases this can be a poor approximation at the stagnation line (compare Figures II.3 and II.6), and that the relative error increases dramatically with distance toward the accretion limit.

for several cases, from numerically integrated trajectories. From these, it was possible to derive a parameterized general form of the $\beta(\alpha)$ curve which depends only on the parameters β_0 and α_{\max} .

Use of this parameterization is the major time-saving device in the model. Several other options are provided for simulating the effect on the $\beta(\alpha)$ distribution of a droplet size spectrum which cannot be adequately modelled by the median volume size. This effect is principally in the extension of the maximum accretion angle.

A second major purpose in developing this model was to incorporate the local ice density variations into the shape of the rime ice accretions. An empirical, parameterized function is used to describe this variation; and it is shown to be reasonably successful in reproducing observed accretion shapes on non-rotating cylinders under a variety of conditions.

A third purpose was to attempt to simulate a slow rotation of the collecting cylinder about its length axis. This is meant to imitate the effects of a gravitational torque on an ice-loaded cable. Although this capability has now been built into the model, it has not yet been been possible to test it against laboratory or field data.

In order to effectively examine the validity of these methods, only dry, rime ice accretions considered. This is so that any comparisons with experimental results are not dependent on poorly known quantities, such as heat transfer,

which affect wet icing processes.

In the experimental accretion experiments used to aid in model development and verification, a crucial measurement is the droplet size spectrum, and its median volume diameter. All droplet spectra were measured by the oil slide method, an outline of which is given in Mason (1971). However, the mvd's derived from such measurements have been found to show a systematic error by Makkonen and Stallabrass (1984). We therefore apply their empirical correction formula to our data in all but a few cases, which fall outside the range of validity for the correction. The procedure followed in these cases is described in Appendix III.

To ensure rime icing under the conditions typical of the wind-tunnel experiments in which collectors with diameters of 2 to 5 cm were exposed to typical atmospheric liquid water contents and wind speeds, the air temperature in the model is assumed to be always at or below -10° C. For application to general conditions, the model should include a calculation of the critical liquid water content, or 'Ludlam limit', (Ludlam, 1951) in order to determine whether the ice will be wet or dry.

B. Parameterization for Local Collision Efficiency

For the idealized situation of straight-line droplet trajectories the local collision efficiency is given simply by $\cos \alpha$, where α is the angle of the local surface normal

with respect to the free stream direction. The collision efficiency has a maximum value of unity at the stagnation line, and a minimum of zero at $\alpha = \pi/2$.

In a more realistic situation, the trajectories curve away from the free stream direction as they approach the cylinder. The stagnation line collision efficiency is reduced from 1, and the local value reaches zero for some $\alpha = \alpha_{\max} < \pi/2$. This situation can be well approximated (especially for large droplets whose trajectory curvature is small) by directly scaling the $\cos \alpha$ distribution to these two extreme values, as follows:

$$\beta(\alpha) = \beta_0 \cos((\alpha/\alpha_{\max})(\pi/2)),$$

III.1

where the two parameters β_0 and α_{\max} are determined from tabulated values or approximating functions given in Chapter II.

A form of time dependence is built into this method since the collision efficiency at a point depends on its local surface angle, which changes with time as the ice layers are accreted on top of one another. The parameters β_0 and α_{\max} do not change with time, an approximation which becomes worse as the accretion shape grows, and departs from a circular cross-section.

For smaller droplets the simple approximation of Equation III.1 fails, since the trajectories are curved by much larger amounts. It has been found that, for a $\beta_0 <$

0.65, the $\beta(\alpha)$ curve is more closely approximated by:

$$\beta(\alpha) = \beta_0 - (\beta_0 (1 - \cos^{0.5} \alpha) / \cos^{0.5} \alpha_{\max}). \quad \text{III.2}$$

For a 0.0254 m diameter cylinder at typical ground level wind speeds of a few tens of metres per second, this corresponds to water droplets of less than 20 microns diameter.

Figure III.1 illustrates the goodness of fit of these functions against local collision efficiency curves calculated numerically by Oleskiw (1982), and approximated by Lozowski, Stallabrass and Hearty (1983).

In both Equations III.1 and III.2, a monodisperse droplet size population has been assumed, or at least a distribution for which the collision efficiency may be reasonably approximated by the collision efficiency of the median volume droplet, or mvd. The results of Appendix VI suggest that, for typical ground-level wind speeds, the single droplet size approximation holds well for $\beta_0 \geq 0.5$.

Otherwise, the effect of the size spectrum, and particularly its larger drops, should be accounted for. Droplet size distributions for icing purposes are commonly described by volume percentages in several size bins of finite width (see, for example, Figure III.2). Applying either Equation III.1 or III.2 to the median volume droplet of each size bin, a volume-weighted average distribution may be calculated. Such a weighted curve then illustrates the

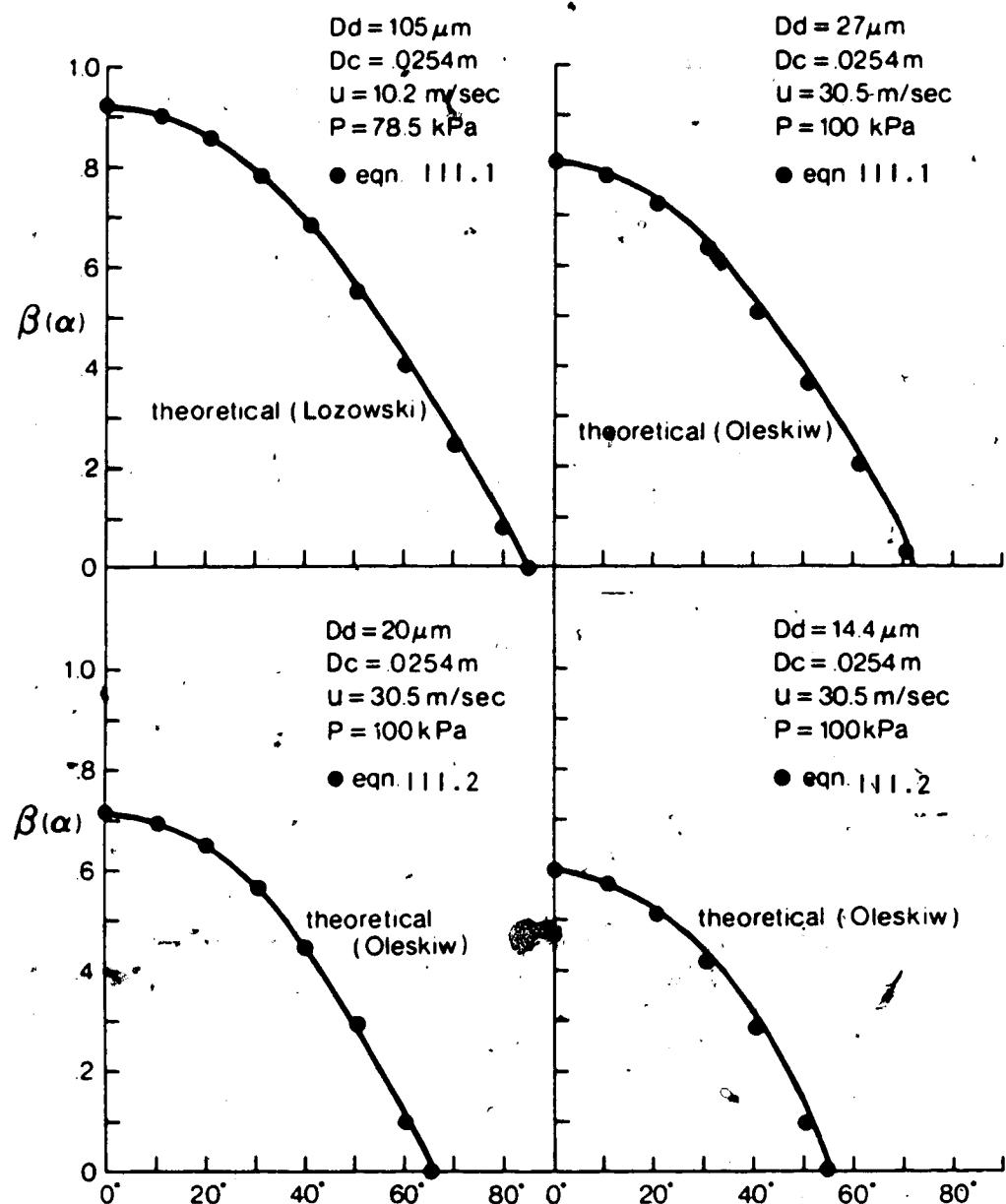


Figure III.1 - Parameterized $\beta(\alpha)$ values from Equations III.1 and III.2, compared to numerically calculated curves of Oleskiw (1982), and parameterized curve after Lozowski, Stallabrass, and Hearty (1983).

effect of the smaller droplets in extending the maximum accretion limit.

This weighting procedure, for a given input size spectrum, may be carried out within the model. This is the "weighted spectrum" option. However, the nature of the averaging procedure eventually gives rise to first-order discontinuities in the $\beta(\alpha)$ curve, and thus on the accretion surface. For this reason, the weighted spectrum option is not recommended when more than three ice layers are accreted, no matter how thin these layers are.

An alternative option is included to avoid this problem, which has been called the "simulated spectrum" option. It is another form of the $\beta(\alpha)$ curve, this time based on weighted average curves (from laboratory measured spectra) which have been scaled and fitted with polynomial functions.

for $\beta_0 < 0.65$:

$$\beta(\alpha/\alpha_m) = \beta_0 (1 + 0.22(\alpha/\alpha_m) - 3.08(\alpha/\alpha_m)^2$$

$$+ 1.86(\alpha/\alpha_m)^3),$$

III.3

and for $\beta_0 > 0.65$:

$$\beta(\alpha/\alpha_m) = \beta_0 (1 + 0.029(\alpha/\alpha_m) - 1.94(\alpha/\alpha_m)^2$$

$$+ 2.48(\alpha/\alpha_m)^3 - 4.11(\alpha/\alpha_m)^4 + 2.54(\alpha/\alpha_m)^5),$$

III.4

where $\alpha_m = \alpha_{max} + Tl$, and Tl is the "tail length", an extension to α_{max} , derived from numerical results and correlated to the mvd. For $\beta_0 < 0.65$:

$$Tl = 0.52 \exp(-mvd / 70)$$

III.5

and for $\beta_0 > 0.65$:

$$Tl = 0.37 \exp(-mvd / 70).$$

III.6

Here mvd is in μm and Tl is in radians. Equations III.3 and III.4 are compared in Figure III.2 with two weighted average curves for the cloud droplet spectra shown.

Although these methods may seem to have become very superficial, they do retain the desired economy of computation with an acceptably small loss of accuracy.

C. Variation of Local Density

A study of the local density variations in circular cylinder rime accretions will be discussed in some detail in Chapter IV. Here the result is presented briefly; an empirical curve describing local density as a function of surface slope.

Based on cross-sectional thickness measurements of several shallow rime layers grown in the laboratory under known conditions, the following function was derived for $\alpha/\alpha_m < 1$:

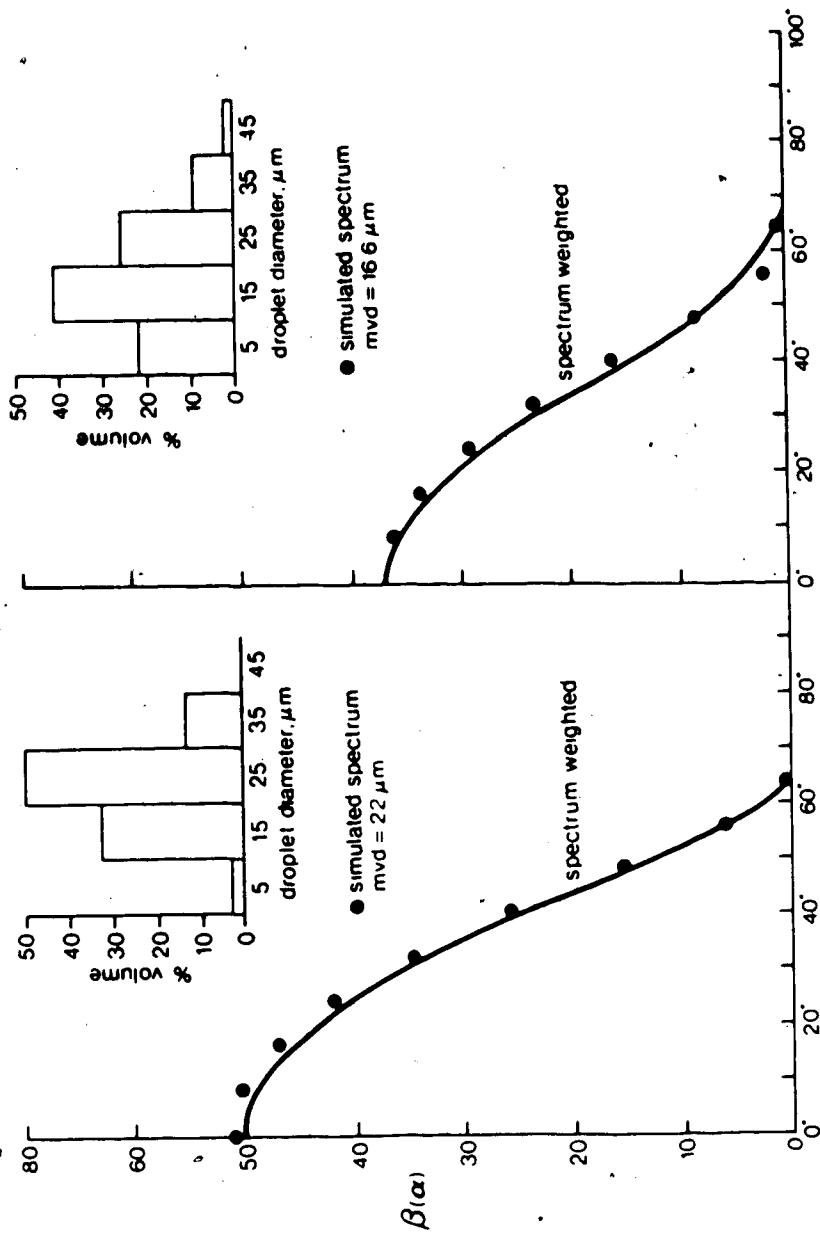


Figure III.3 - Parameterized $\beta(\alpha)$ values from Equations III.3 and III.4, compared to weighted average curves for the cloud droplet spectra shown. These spectra are derived from the distribution of Best (1951).

$$\rho/\rho_0 = (1 - 0.143(\alpha/\alpha_m) - 0.246(\alpha/\alpha_m)^2$$

$$- 0.309(\alpha/\alpha_m)^3).$$

III.7

Here α_m is defined as for Equations III.3 and III.4, and in this version of the model, $\rho_0 = 890 \text{ kg/m}^3$. This assumption is made for simplicity; in reality the rime densities may be much lower, and this will affect the accretion thickness and overall shape. Chapter IV will discuss methods for estimating more realistic values of ρ_0 .

The remainder of this section will examine the predictive success of the model under a variety of input conditions. Figure III.3 compares accretion profiles and masses from the model against ice accretions grown in wind tunnel experiments carried out in the University of Alberta's FROST tunnel (this facility and our experimental methods are described in Gates, 1981 and Lozowski and Gates, 1984). For the ranges shown of mvd, liquid water content and free-stream speed the modelled mass is always within 10% of the observed mass, and the major features of the experimental shapes are reproduced in the profiles.

However, it is not expected that the model will reproduce such features as ridges or troughs near the stagnation line. These are most likely due to changes in the flow pattern as the accretion grows, and the resulting

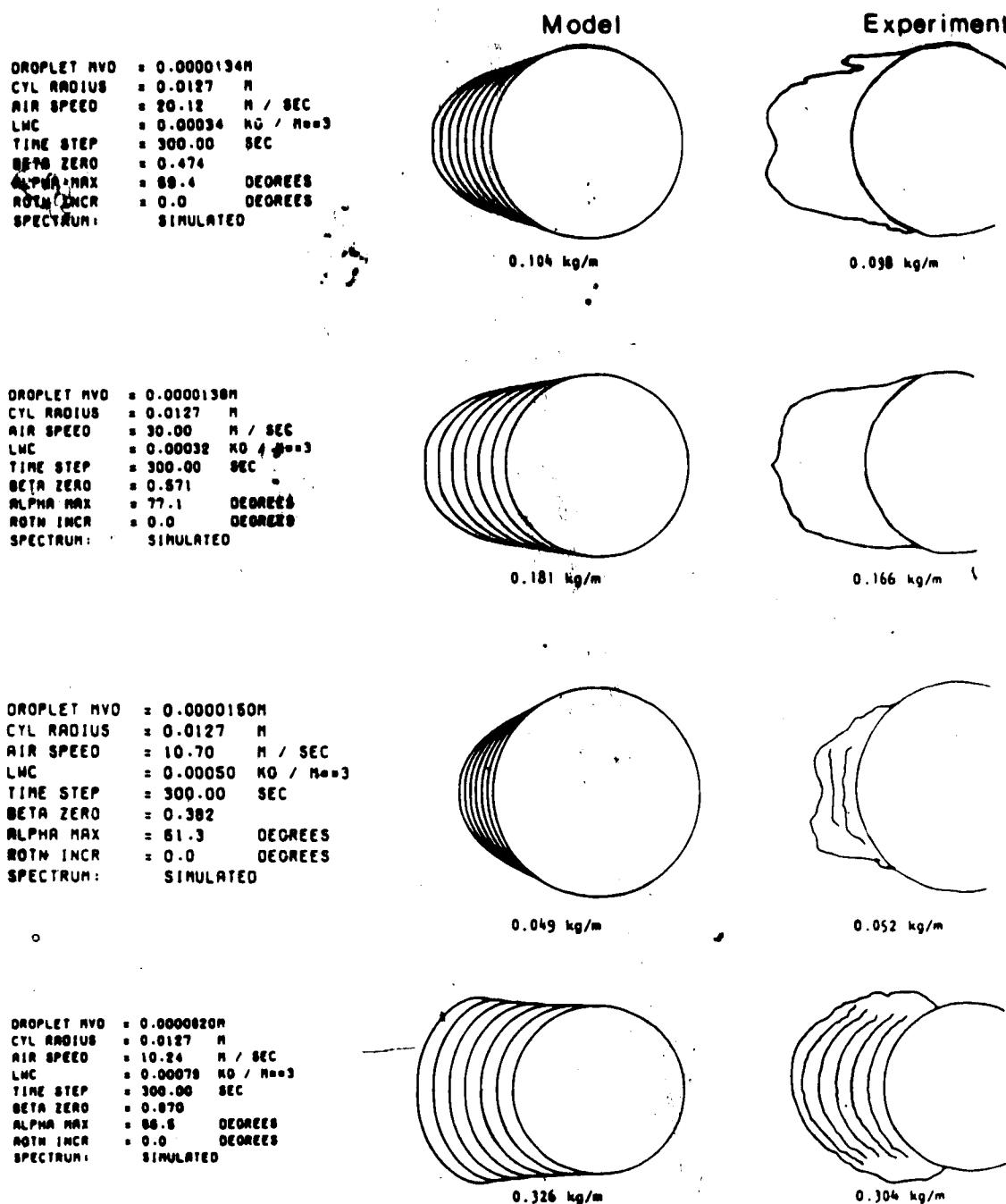


Figure III.3 - Comparisons between modelled and experimental profiles and masses per unit length for the conditions listed at left. The lower two experimental profiles have layer outlines which were marked by spray paint.

changes in the impingement characteristics.

In particular, the value of β_0 will change with time, while the model assumes a constant value. Even with this assumption, however, it is able to distinguish between widening and narrowing profiles for small differences in growth conditions, and can produce the flat front surface which often precedes the formation of ridges and troughs in the experiments.

D. Rotation

An option has been included in the model which will allow the cylinder plus accretion to be rotated with respect to the free stream by a small amount after each layer is accreted. No attempt is made to calculate the correct rotation rate due to the gravitational and aerodynamic torque; the rotation speed in degrees per layer must be specified as one of the input parameters.

In order to achieve this effect within the model algorithm, the local collision efficiency and local density must be calculated as functions of the position angle rather than the surface angle on the current surface. (The position angle at a surface point is the angle between a line joining the point to the centre of the cylinder, and the free-stream direction.) This somewhat less accurate method was adopted in order to avoid problems when accreting over top of the previous layers' accretion limit.

There are presently no laboratory or field data available with which to test this part of the model. It is presented here only as an option for future study. An example of a rotating case is shown in Figure III.4.

The Fortran77 code written for the model is included in Appendix II. The code will also run on an IBM or compatible personal computer fitted with a math co-processor and graphics capability; a six layer rotated accretion takes about 90 seconds to compute on a regular PC.

E. Model Applications

This model is intended for use in the laboratory, in conjunction with wind-tunnel experiments. There the required input data are readily available, as accurately as present equipment and methods can determine them.

However, if this model or some future version of it were to be applied in the field, the situation becomes very different. Good estimates of the input data (duration, wind speed, liquid water content, droplet mvd and type of spectrum) must somehow be obtained. In addition, a field version should be able to predict at least the transition to wet icing, and so the air temperature will also be required as input.

These data are rarely available for a remote site. Local meteorological services may provide estimates or forecasts for duration, wind speed, air temperature, and precipitation, but usually not liquid water content or

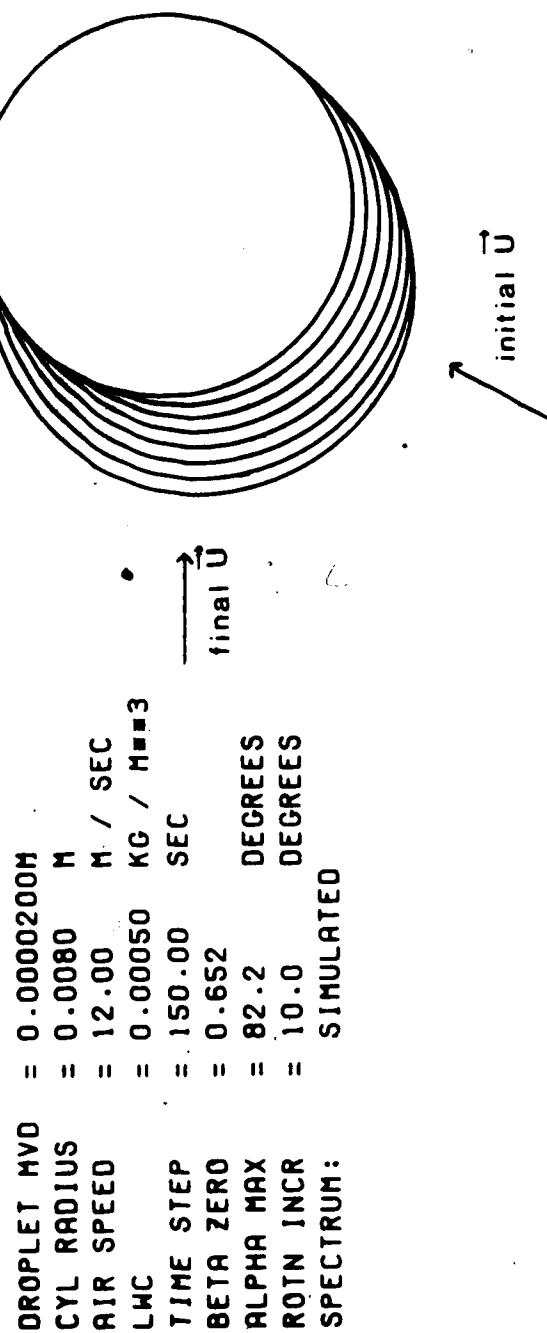


Figure III.4 - An example of the model output for a rotating cylinder.

cloud droplet size.

Clearly, real-time icing forecasting with such a model is a long way off. However, the model could conceivably be used to determine icing climatologies for different regions where adequate meteorological data have been obtained. This would mean sufficiently long-term observations of liquid water contents and droplet sizes (i.e. cloud type) to give an idea of the typical values of these quantities under the typical meteorological situations which lead to icing events in a given region.

However, given the current state of knowledge of field conditions, the best application for this model remains in furthering the understanding of what is happening under controlled rime icing conditions in the laboratory, which it is capable of doing much more efficiently and cheaply than previous models.

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IV. Local Densities of Circular Cylinder Rime Accretions

A. Introduction

A situation very commonly described for the study of rime ice accretion is that of a non-rotating, circular cylinder, enveloped in a steady cross-flow of cold, cloudy air. Rime ice is formed when all of the impinging cloud droplets are able to freeze completely at the point of impact, and before the arrival of another droplet. Rime ice is always 'dry', which means there is no liquid water on the surface or trapped within the ice. The conditions necessary for dry ice to form are described by the so-called "Ludlam limit" (Ludlam, 1951), a maximum value of cloud liquid water content computed as a function of air temperature, cylinder size and air speed. Above this limiting value, the accretion will be 'wet', or glaze ice, where the droplets may run along the surface before freezing, or are incorporated into pockets of liquid water within the ice. This results in a much higher and more uniform density of ice. This study deals only with rime ice.

Examination of a rime accretion in cross-section (Figure IV.1) reveals details of the internal structure which govern the local density. Air bubbles and channels are apparent, usually increasing in number and size toward the lateral edges, away from the stagnation line. At the extreme edges of the accretion are often found the individual elongated or fan-shaped structures known as



Figure IV.1 - A 1.5 cm thick cross-section of a circular cylinder rime accretion, viewed in transmitted light. The accretion conditions are those of Sample 1 in Table IV.1.

'rime feathers', which seem to merge into the accretion from the sides. These are seen more easily in a lateral view (Figure IV.2).

It is apparent even from such a cursory examination that there should be a trend in the local ice density, decreasing from the stagnation line out toward the edges. When averaged over intervals larger than the typical air channel size, a smooth distribution should emerge. The actual shape of this distribution and how it arises are of interest to modellers of the icing process.

Numerical models have been devised which predict the results of accretion experiments performed in wind tunnels under controlled conditions, e.g. Oleskiw (1982), Lozowski, Stallabrass and Hearty (1983), McComber and Touzot (1981). The parameters which determine the rate of ice growth in such models are the accretion object's size, the air temperature and pressure, air speed, liquid water content and the sizes and relative number distribution of the water droplets in the air stream. These quantities are needed to calculate the local and overall collision efficiencies on the collecting object, which together with the local ice density determines the mass and the shape of the modelled ice deposit.

Although collision efficiency and its local distribution have been well studied, at least for circular cylinders (see for example, Oleskiw, 1982, McComber and Touzot, 1981, Langmuir and Blodgett, 1946), the density of

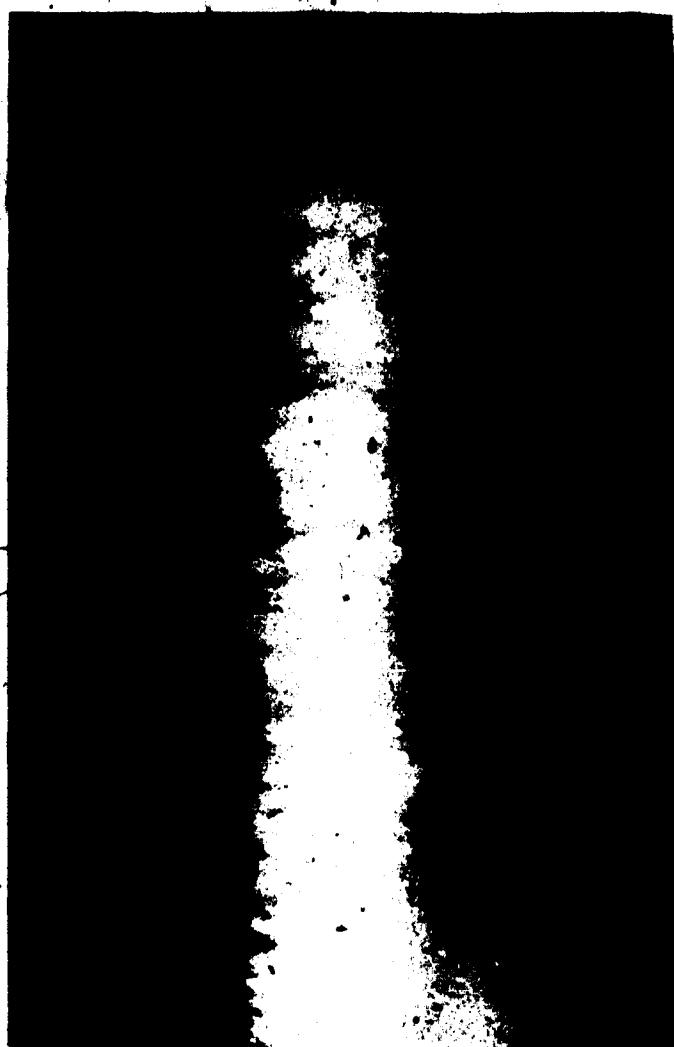


Figure IV.2 - A typical circular cylinder accretion showing
rime feathers.

rime has not received as much attention. Some studies have dealt with the densities of ice accretions on rotating cylinders or hailstones (e.g. Macklin, 1962, List, Cantin and Ferland, 1970, Makkonen and Stallabrass, 1984).

Bain and Gayet (1983) have attempted to adapt these results to the fixed cylinder case; their method will be covered in the next section. Problems with this method, very similar to those outlined by Rasmussen and Heymsfield (1985), are also discussed, and a new density distribution is derived directly from experimental accretion shapes. Both distributions are tested in the model, and also against direct laboratory measurements of local density.

B. Rotating Cylinder Densities Applied to the Non-rotating Case.

Macklin (1962) made extensive measurements of the bulk densities of rotating cylinder ice deposits. The results were found to correlate best with the parameter, called R , given by:

$$R = r_d^4 v_0 / T_s \quad IV.1$$

where r_d is the radius of the medium volume droplet, v_0 is the impact speed of that droplet at the stagnation line, and T_s is the mean surface temperature of the accretion. Similar correlations were found for the experimental results of Bain and Gayet (1983) and Makkonen and Stallabrass (1984).

on rotating cylinders, although the latter found significantly larger densities in the lower range of R .

Makkonen and Stallabrass attribute this difference to their method of measuring droplet sizes with optical probes. Macklin's droplets were measured using the oil slide method, which according to Makkonen and Stallabrass gives underestimated sizes, in comparison to the optical probe method. Makkonen and Stallabrass' correlation for the range $0.2 < R < 1.0$ is:

$$\rho(R) = 0.378 + 0.42 \log_{10}(R) - 0.0823(\log_{10}(R))^2. \quad IV.2$$

Bain and Gayet (1983) made an attempt to apply the earlier correlation of Macklin to densities on a fixed cylinder accretion, by introducing local variations into the R parameter as follows:

$$R(\theta) = r v_r(\theta) / T_s(\theta), \quad IV.3$$

where now $v_r(\theta)$ is the radial component of droplet impact velocity as a function of the radius angle θ (measured with respect to the stagnation line). Bain and Gayet derive the total droplet speed at impact from an assumed cosine distribution of speed with cylinder radius angle, scaled to the stagnation line value as calculated by Langmuir and Blodgett. $T_s(\theta)$ is the variation of local surface temperature calculated from the model of Lozowski,

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Stallabrass and Hearty (1983).

An initial objection to this approach is philosophical. Macklin's correlation is a purely empirical one; it is not a physical model and therefore should not be extended to a physical situation so different from that which provided the original data.

On the fixed cylinder, $v_r(\theta)$ is not even analogous to v_0 on the rotating cylinder, which is the maximum impact speed felt at any one point during one rotation.

Furthermore, on the rotating cylinder, the droplets arrive at a given point from a large range of angles, which should perhaps result in closer packing of the droplets and a higher density. More importantly, the shadowing effects of feather structures and the formation of air channels between them is not accounted for.

However, a more concrete objection is found by comparing this type of distribution to that found by direct measurement of the density, and by comparing the model accretion shapes and masses it produces to experimental profiles. The first comparison will be discussed in Section D of this Chapter. For the model comparison the simple rime accretion model described in Chapter III is used.

The experimental accretion profile shown on the left in Figure IV.3 was grown, under the conditions listed, in the University of Alberta FROST tunnel. Also shown are model profiles produced by three different distributions of local

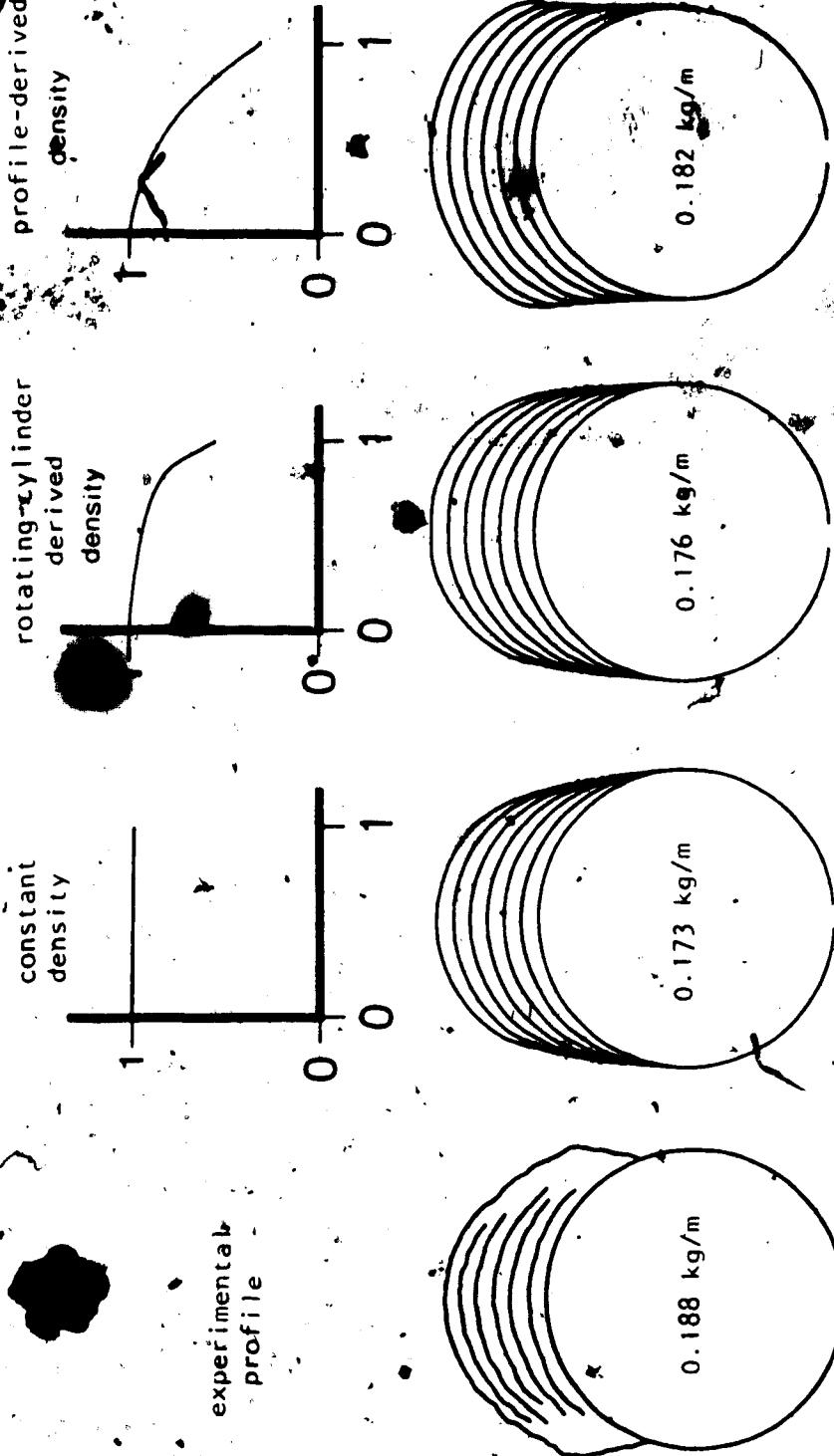


Figure IV.3 - Experimental and three model accretions showing the effects of three different density distributions (plotted above as ρ/ρ_0 vs. r/a)
 Conditions are: $D_d = 51 \mu\text{m}$, $D_c = 0.0254 \text{ m}$, $U = 10.7 \text{ m/sec}$, $1 \text{ cm} \times 0.49 \text{ g/m}$, $\tau = 300 \text{ sec}$

density. The profile at centre left results from a constant density, that at centre right from a distribution that is derived following Bain and Gayet. The profile at the far right uses a third distribution which is described below.

The three density distributions are illustrated above their respective profiles. For easy comparison, each is plotted as ρ/ρ_0 vs. α/α_m , where ρ_0 is the local density at the stagnation line, α is the local angle of the surface normal with respect to the free stream direction, and α_m is defined in Chapter III. On the bare cylinder, α is identical with θ , but not for subsequent ice layers. The angle α is used rather than θ , so that the density distribution will be time-dependent, changing as the ice shape changes.

The middle distribution is derived from Equation IV.1, using $v_r(\alpha)$ as calculated by Bain and Gayet, and $T_s(\alpha)$ calculated by the model of Lozowski, Stallabrass and Hearn. The resulting curve is fitted with a third-order polynomial:

$$\frac{\rho(\alpha/\alpha_m)}{\rho_0} = (1 - 0.303 (\alpha/\alpha_m)) \\ + 0.936 (\alpha/\alpha_m)^2 - 1.074 (\alpha/\alpha_m)^3.$$

IV.4

The second modelled profile is very similar to the first, constant density profile, which was too narrow in comparison with the experimental shape. Both of the modelled masses are less than the observed mass, by 8 and 6

percent, respectively. The result using the third distribution shown is in better agreement with the experimental shape, by having a widening profile, and a mass within 3 percent of observation. This distribution and its derivation are described next.

C. Profile-Derived Density Distribution

The thickness (Th) of ice deposited in time τ at a given point on the bare cylinder surface, which we describe by its surface angle α , is

$$Th(\alpha) = \beta(\alpha) U lwc \tau / \rho(\alpha), \quad IV.5$$

where U is the free-stream speed, and lwc is the liquid water content. This version of the dry-growth equation ignores curvature effects, e.g. as described by Lozowski, Stallabrass and Hearty (1983). This is a reasonable simplification given the very small spacing of the points which define the icing surface in the model.

Given IV.5 then, if both the local thickness and the local collision efficiency $\beta(\alpha)$ are known, and if the interval τ is assumed to be small enough that $\beta(\alpha)$ and the density $\rho(\alpha)$ remain essentially constant, IV.5 may be solved for $\rho(\alpha)$.

Figure IV.4 shows another experimental accretion grown in the FROST tunnel, and its growth conditions. The total accretion time was 30 minutes. Every five minutes, a layer



Figure IV.4 - A circular cylinder rime deposit in profile, with paint layers marking the surface at five minute intervals. Conditions are those of Case B in Table III.1.

of paint was sprayed onto the surface, allowing changes in the shape of the ice surface to be followed with time. It also allowed local thickness measurements of thin layers to be obtained, which are unaffected by the melting or sublimation losses that must occur on the top surface, between the time of the end of the test and photography of the profile.

The distribution of local collision efficiency was calculated from the median volume diameter, and an assumed droplet size distribution, scaled to the mvd, based on measurements taken in the same wind tunnel. The mvd was originally measured by the oil slide method, which is known from the work of Makkonen and Stallabrass (1984) to be an overestimate; but for Case B of Table IV, the oil slide estimate falls outside of the size range for which Makkonen and Stallabrass' empirical correction formula is valid. Therefore, an mvd has been adopted which was derived from the stagnation line thickness measurement, the liquid water content measurement, and an assumed local density for the stagnation line of 890 kg/m^3 (a reasonable density in this case based on visual inspection of the accretion). The method used to derive this estimate of the mvd is described in Appendix III.

For Case A of the same table it was possible to make use of the correction formula mentioned above to correct the oil slide measured mvd.

In this way the distribution $\rho(a)$ has been found for several layers of two different accretions. The data are listed in Table IV.1 and plotted in Figure IV.5. A cubic polynomial fit (solid curve, Figure IV.5) gives:

$$\rho(a/a_m)/\rho_0 = (1 - 0.143 (a/a_m))$$

$$- 0.246 (a/a_m)^2 + 0.309 (a/a_m)^3$$

IV.6

The relative distribution is thus assumed independent of g_{fc} and air temperature; this assumption implies that these factors only come into the determination of the scaling parameter ρ_0 . Normalization of the distribution by the two parameters ρ_0 and a_m makes the expression above easy to use in an icing model, assuming that it is generally applicable to a range of accretion conditions. Evidence for this was presented in Figure III.3 of the last Chapter, where the model results are compared to several independent wind tunnel experiments.

D. Laboratory Measurements of Local Density

For further evidence to support our proposed distribution of local density, a series of direct measurements were made from wind-tunnel accretions.

Following the techniques described by List, Cantin and Ferland (1970), local densities were measured by drilling small holes (2.5 - 3.0mm diameter, less than 1 cm depth) in

	(degrees)	Th (mm)	$\beta (\alpha)$	$\rho (\text{kg/m}^3)$
Case A	0	3.34	0.407	767
	10	3.12	0.393	792
	20	3.25	0.334	646
	30	2.9	0.249	540
	40	1.7	0.155	574
	50	0.9	0.069	482
Case B	0	1.48	0.842	890
	10	1.48	0.825	871
	20	1.48	0.776	820
	30	1.40	0.699	780
	40	1.33	0.592	696
	50	1.03	0.458	695
	60	0.81	0.302	583

Table IV.1 - Local densities as a function of surface slope
 derived from thickness measurements of shallow ice layers
 accreted at -10°C under the following conditions:

	$D_d (\mu\text{m})$	$D_c (\text{m})$	$I_{WC} (\text{g/m}^3)$	$U (\text{m/sec})$	$\tau (\text{sec})$
Case A	15.5	0.0254	0.98	10.7	600
Case B	51.0	0.0254	0.49	10.7	300

$\beta (\alpha)$ is calculated from Equations II.11, III.3 and III.4.

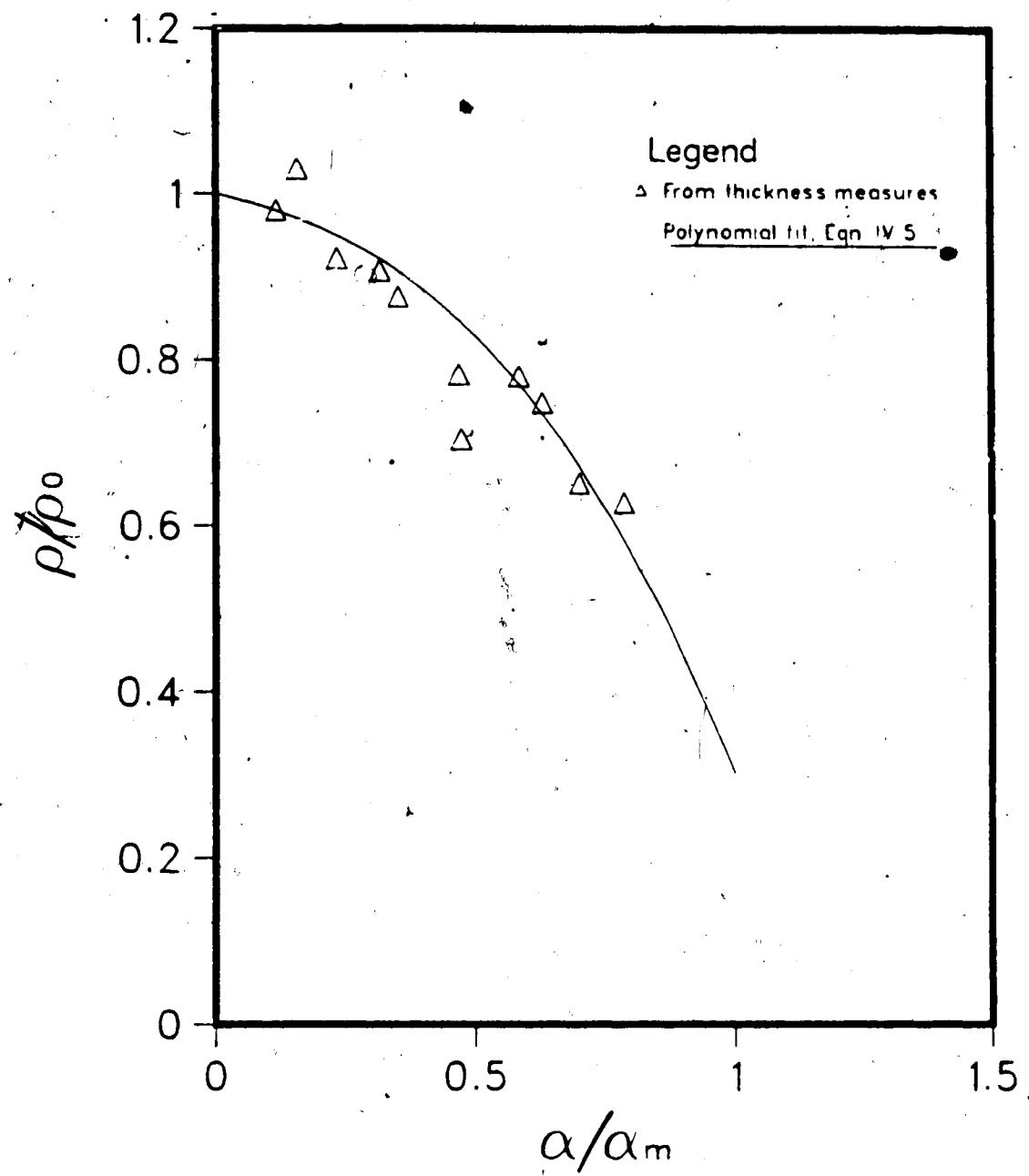


Figure IV.5 - Local densities derived from thickness measurements of shallow ice layers from Table IV.1. The solid curve is a polynomial fit to the data (Equation IV.3).

cross sectional cuts from two different rime accretions (see Figure IV.6). The location, volume and mass loss were recorded for each hole. The mass was measured to within 0.00005 g with a Mettler model H20 balance. For the data of Sample 1, Table IV.2, the hole depths and diameters from three sections of a single accretion were measured to within .05 mm with a Vernier caliper. For Sample 2 of Table IV.2, holes from two sections of accretions grown under nearly identical conditions were measured to 0.01mm in depth and 0.001mm in diameter, with the same bore and depth gauges used by List et al. (1970) to determine local densities in hailstones. The surface angles were approximated by the more easily measured radius angles, since paint layers to define the intermediate surfaces would have altered the densities. This is a good assumption for shallow layers, at any rate, and measurements have been used only from the layer immediately above the cylinder surface. The error in using the radius angle approximation is estimated to be five degrees.

Another source of error is mass loss from the entire sample through sublimation and handling during the hole drilling procedure. Several experimental estimates were made of the size of this error, and the average (0.001 g) was subtracted from the measured mass losses from the holes.

The resulting density measurements are shown in Table IV.2 and in Figures IV.7 and IV.8. The solid and dashed curves shown in each of these figures are from Equations



Figure IV.6 - An accretion section of about 1.5 cm. thickness, showing the holes drilled for the local density measurements. The condition are those of Sample 1 in Table IV.2

Sample 1		Sample 2	
α ($\pm 5^\circ$)	ρ (± 60 kg/m 3)	α ($\pm 5^\circ$)	ρ (± 30 kg/m 3)
2	786	2	907
2	667	3	902
2	660	4	723
6	719	6	747
8	620	7	770
8	792	8	909
10	714	13	788
12	645	15	815
12	688	17	766
12	685	19	724
15	629	23	823
18	519	24	683
20	715		
21	630		
27	675		
28	427		
30	681		
30	733		
30	701		
35	539		
38	580		
42	749		
50	504		

Table IV.2 - Experimentally measured local densities
as a function of surface slope taken from accretions
grown under the following conditions:

D_d (m)	D_c (m)	I_{WC} (g/m 3)	U (m/sec)	T ($^\circ$ C)
16.4	0.046	0.53	14.9	-14
14.0	0.0254	0.36	30.0	-10

IV.2 and IV.4, respectively, where the choice of ρ_0 has been based on the density measurements, and α_{\max} is calculated from the median volume diameter droplet, following the methods of Chapters II and III.

The scatter of points in both figures is considerable, and is much larger than the measurement errors represented by the error bars will allow. The most likely cause of the large scatter is in the nature of the rime itself. The internal air channels are scattered at random, and since they are often of the same size as the holes, this will introduce a natural scatter. Also, in the very fragile rime, the drilling process may cause walls between adjacent air channels to collapse, resulting in additional mass loss. This problem is especially bothersome near the edges of the accretion, where the density is lowest. Here, walls between the holes themselves will sometimes collapse, making it very difficult to obtain density values at large surface angles. Part of the problem might be alleviated by using accretions samples grown on a larger diameter cylinder, so that the holes do not have to be drilled so close together to get the same interval of surface angle. Unfortunately, the only accretions available were grown on cylinders of one and two inches in diameter.

For whatever reason, the results are too scattered to allow a clear choice between one parameterized distribution and the other. This is due, in part, to the lack of data at

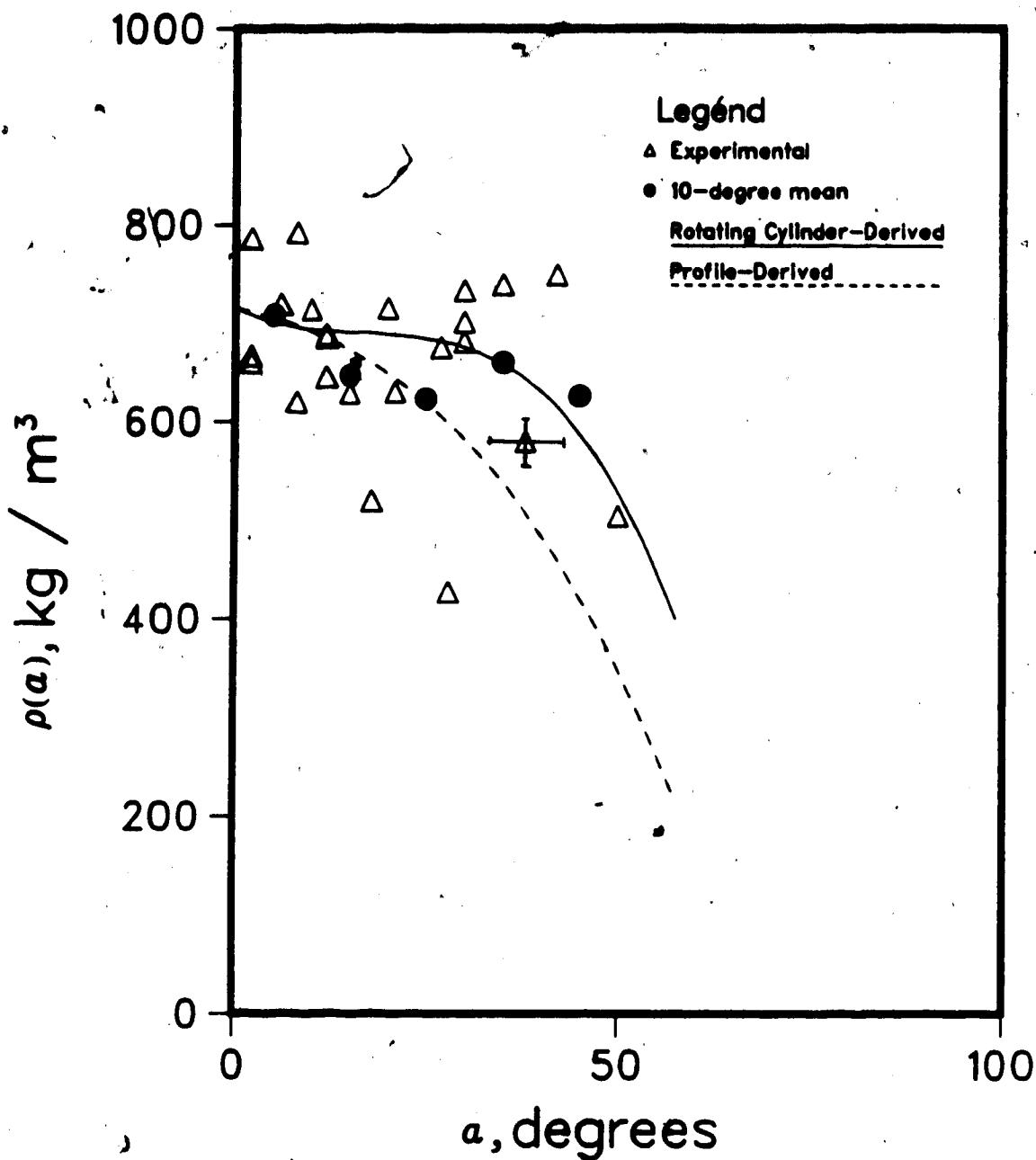


Figure IV.7 - Local density vs. surface angle; experimental measures and parameterized curves for the conditions of Sample 1 in Table IV.2. The filled symbols represent mean values for 10° intervals in surface slope.

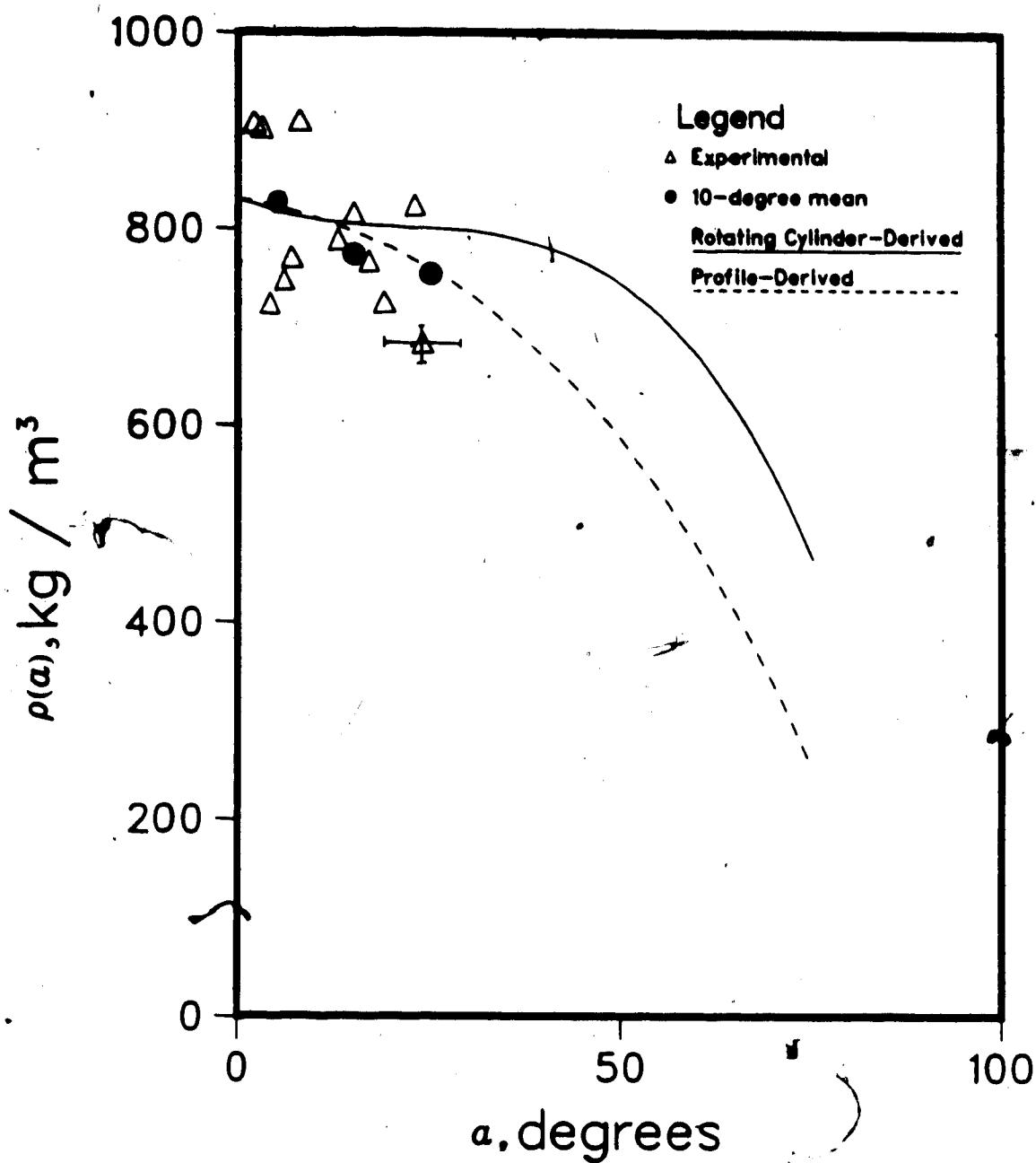


Figure IV.8 - Local density vs. surface angle; experimental measures and parameterized curves for the conditions of Sample 2 in Table IV.2. The filled symbols represent mean values for 10° intervals in surface angle.

large surface angles. When averaged over 10 degree intervals (filled symbols) the experimental data in both figures favour the profile-derived curve up to a surface angle of 30°, but for larger angles Figure IV.7 they lie closer to the rotating cylinder curve. However, there are far fewer points defining the averages for larger angles, and none at all for Figure IV.8.

Although these direct density measurements have not been conclusive, the model results and the practical advantages of the profile-derived distribution justify its adoption over the other. The application of this method is efficient and inexpensive, assuming that the scaling parameters may be easily determined. One of these parameters, the maximum accretion angle, α_{\max} , may be quickly approximated according to the analytical function given in Chapter II, but ρ_0 remains to be determined.

E. Local Density at the Stagnation Line

Qualitative observations of many rime accretions show that the stagnation line density also varies with the environmental conditions. Since there are, at present, insufficient quantitative data on this variation, the correlation for rotating cylinder densities, Equation IV.1, may perhaps be applied. The justification for doing so here, but not for the entire density distribution, is that there is no shadowing at the stagnation line, and the conditions are more nearly analogous to the rotating

cylinder case.

- However, some important differences must be noted between the two situations. In the rotating cylinder correlation, the parameter v_0 is the impact speed of the median volume droplet at the stagnation line only. Over the remainder of the accretion surface, droplets impact at speeds ranging between zero and v_0 , so the average impact speed, at least for cases with large a_{max} , is roughly one half v_0 , and perhaps somewhat less if a_{max} is small. Thus, v_0 for the rotating cylinder is \leq twice the average speed of impact. On the stagnation line of the non-rotating cylinder, the impact speed is constant and equal to v_0 , so in order to substitute into the R parameter the correct analogy for twice the average speed, we use twice v_0 . v_0 is estimated from Equation II.11.

The R parameter is further altered by substituting for the mean surface temperature, the value at the stagnation line ($T_{s,0}$) which is determined using the model of Lozowski, Stallabrass and Hearty (1983). The new correlation parameter is then:

$$R' = 2 r_d v_0 / T_{s,0} \quad IV.7$$

The results are shown in Figure IV.9, for five accretions for which ρ_0 has been experimentally measured by the methods described in the last section. The fit to Equation IV.1 is quite good, considering the size of the

measurement errors in ρ_0 (i.e., the same size as for Figure IV.8). Until more experimental data are available, this is an acceptable method for estimating ρ_0 .

F. Summary

Previous methods for estimating the local variation of rime ice density are examined, and found to be inadequate. A new method is derived empirically, and tested in model accretions, and against direct laboratory measurements of local density. The laboratory measured distribution is inconclusive, but model results show the new method to be more successful than the old for predicting the shapes and masses of circular cylinder rime ice accretions.

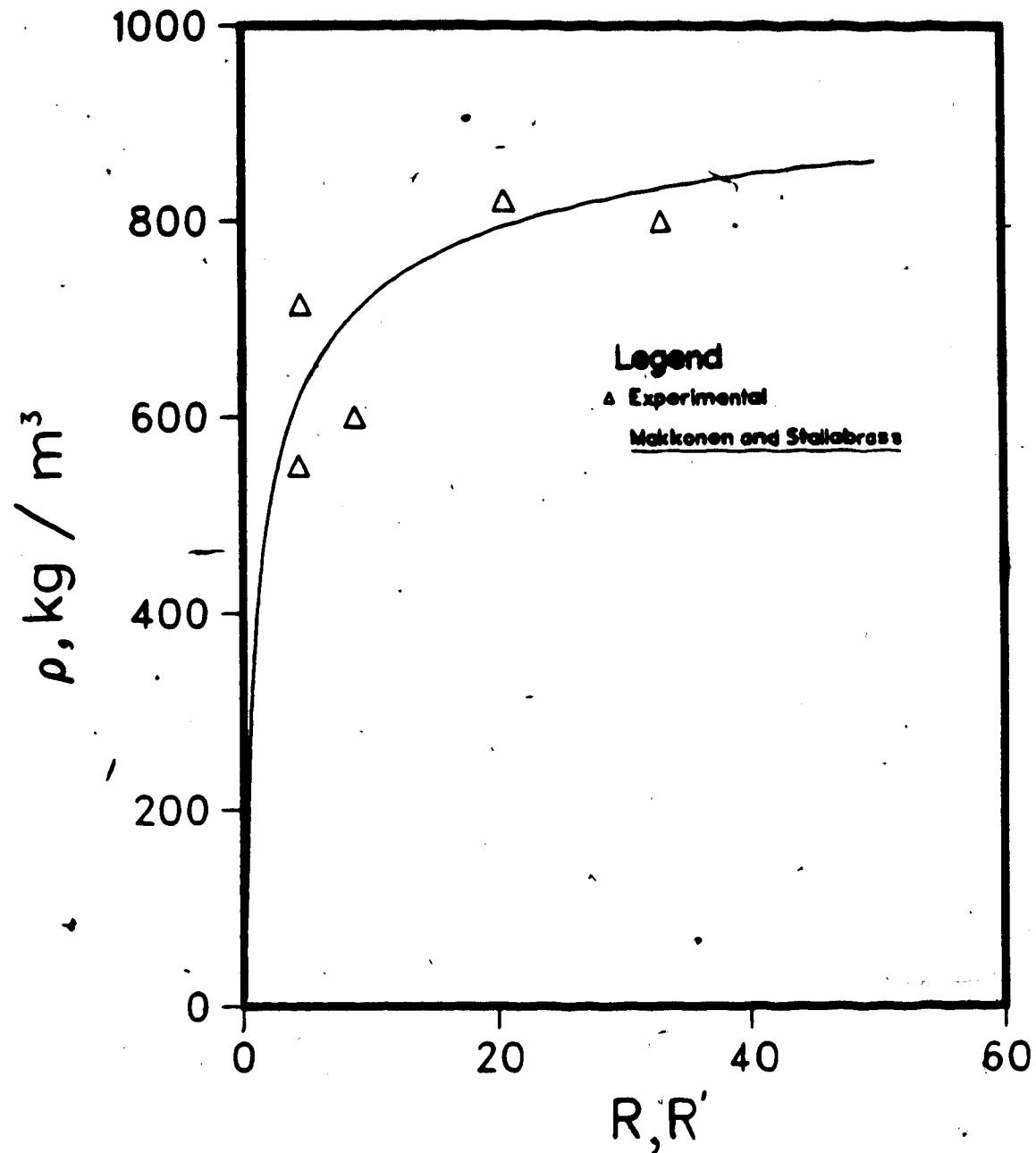


Figure IV.9 - Stagnation line density measurements vs. R' (symbols), and the rotating cylinder density correlation of Makkonen and Stallabrass, Equation IV.2, vs. R (solid line). R and R' are defined in Equations IV.1 and IV.7, respectively.

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V. Rime Icing on Airfoils

A. Introduction

A numerical simulation of ice accretion on an arbitrary airfoil section involves greater computational difficulty than does the corresponding problem of icing on circular cylinders. The major reason is the requirement for numerical integration of the impinging droplet trajectories in order to determine collision efficiencies. Unlike the cylinder case, there is (except for the Joukowski airfoil) no analytical solution of the potential flow to simplify the task, nor are there convenient tables of previously calculated collision efficiencies as functions of the flow and inertia parameters.

This requirement adds considerably to the model's consumption of computing time and resources. Previous airfoil accretion models (e.g. Oleskiw, 1982, MacArthur et al., 1982, Cansdale and Gent, 1983) have thus tended to be expensive to run, and may be prone to numerical instabilities.

However, from the results of one such model (Oleskiw 1982), simple parameterizations have been derived for the local collision efficiency distribution with surface slope on NACA 0015 and 0012 airfoils. Three separate analytical functions approximate this distribution for monodisperse or average droplet spectra, or for any spectrum when the angle of attack is non-zero. Use of these functions requires the

determination of only one value of local collision efficiency, that at the stagnation line; thus considerably reducing the computation time. This single value may then be re-determined at intervals during the accretion to give the model a true time-dependence.

These parameterizations have been incorporated into a new, more efficient numerical icing model. Other computational simplifications include the representation of the airfoil and ice surfaces by straight line segments rather than by a spline curve, and a straightforward algebraic method to determine the droplet's point of collision with the surface. An approximate distribution function for the local density has also been introduced. Again, rime icing only is considered, i.e. the air temperature is assumed to be below -10° C, and the impinging water is assumed to freeze immediately and completely. Most of the parameterizations and simplifications could also be used in a model for wet accretion.

The model results are presented in comparison with those of Oleskiw (1982) and with real accretions grown in wind tunnel experiments.

B. Model Outline

A brief outline of the model algorithm is as follows:

Step 1. Specify the x, y coordinates of the original airfoil surface. The coordinates are dimensional,

with the origin at the nose of the airfoil.

Step 2. Find the local collision efficiency at the stagnation line, β_0 , by the following procedure:

1. Solve for the potential flow field about the airfoil or accretion shape by the method of Kennedy and Marsden (1976).
2. Integrate two droplet trajectories near the stagnation line and find their collision points.
3. Calculate β_0 from $\Delta y / \Delta l$, the ratio of initial separation normal to the free-stream of the trajectories to their final separation in arc length on the surface.

Step 3. For each surface point, calculate the thickness accreted perpendicular to the surface during one layer time step τ . The local collision efficiency at each point is approximated from one of three parameterized functions of β_0 and the local surface slope, α , depending on the type of droplet size spectrum and the angle of attack, Φ (defined as the angle between the free-stream direction and the airfoil chord). The local density distribution is also approximated from a parameterized function of surface slope, α .

Step 4. Calculate the mass and mean density of the accreted ice layer.

Step 5. Repeat Steps 3 and 4 for the number of layers requested. Repeat Step 2 only for every fifth layer. The layer time step, τ , should be chosen so that the combined stagnation line thickness of five layers is not more than 1/10 of the original chord length.

Step 7. Plot the original airfoil and all accreted layers.

The sections below will discuss each of Steps 1 through 4 in detail.

C. Specifying the Airfoil Shape

This version of the model employs the standard airfoil section known as the NACA (United States National Advisory Committee on Aerodynamics) four-digit wing section, for which Abbott and von Doenhoff (1959) give the thickness distribution in non-dimensional x,y-coordinates normalized to the chord length, as:

$$y = 0.05 h_0 (0.2969 x^{0.5} - 0.126 x - 0.3516 x^2)$$

$$+ 0.2843 x^3 - 0.1015 x^4)$$

V.1

The number h_0 gives the maximum airfoil thickness as a

percentage of the chord length, as shown in Figure V.1, and also makes up the last two digits of the "four-digit" designation. The first two digits represent c_m , the maximum camber as a percentage of the chord, and c_p , the abscissa of the camber ordinate in tenths of the chord. All of the parameterizations used in the model are derived assuming that both c_m and c_p are zero. The maximum thickness h_0 is entered as model input, but expressed as a fraction of the chord, rather than a percentage.

Although defined above for non-dimensional coordinates, the model carries the airfoil and accretion profiles in dimensional coordinates throughout most of the calculations, because the chord length (initial chord plus accumulated ice thickness) is changing with every layer step. Where a non-dimensional profile is required for the potential flow calculation, the coordinates are normalized using the current chord length.

Since the profile is always represented by a vector of coordinate points, the points must be of sufficiently high density to ensure that the curved surfaces are well approximated by straight line segments joining them. This avoids the expensive computational process of fitting a few surface points with a spline curve, which may be prone to instabilities. It also simplifies the determination of trajectory impact points, but it does require more data storage. However, even this can be minimized by using a point distribution scheme which concentrates points in the

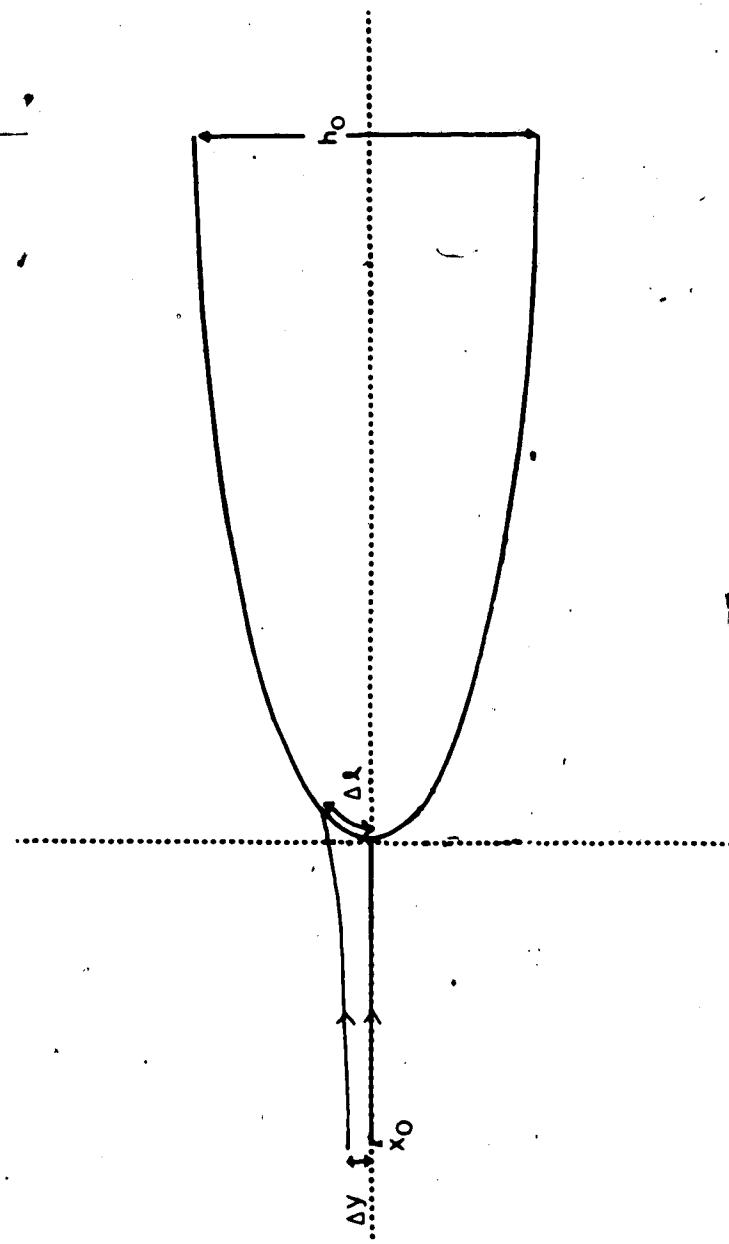


Figure V.1 - Determination of collision efficiency on an airfoil. $\beta = \frac{dy}{dx} \approx \frac{\Delta y}{\Delta x}$. Also illustrated is the definition of Φ , the angle of attack.

regions of strong curvature at the nose and tail.

One such scheme is that suggested by Kennedy and Marsden (1976), in which the distribution of non-dimensional x-coordinates is given by:

$$x_i = 0.5 (1 - \cos(2\pi i / N)) \quad V.2$$

where i lies between 1 and N , an even integer. This scheme was found to give good results with $N = 600$ to specify the complete airfoil section for plotting purposes, although only half this number of points, defining the front portion of the airfoil, is allowed to accrete ice.

In the potential flow solution, a profile consisting of a much smaller number of points is desired to reduce the computational requirements; Kennedy and Marsden suggest that 40 points is an optimum number. This can be simply constructed from either the original airfoil profile, or from subsequent accretion shapes by taking every 15th point of the denser profile. The new sparse profile will also retain the desired distribution pattern of points.

Accretion thicknesses are calculated for each surface point in the direction perpendicular to the local surface slope. This causes the point separation to increase near the nose with each succeeding layer. After several layers, the separation may have become too large for a reasonable representation of the surface by straight line segments, and has destroyed the original point distribution.

A way to remedy the situation is to shift each coordinate point, after each layer is accreted, closer to the nose along the straight line segment joining it and its nearest neighbour (whose position has already been shifted). The amount of the shift must be carefully chosen to preserve the desired distribution of points, and an accurate profile. Because the original distribution concentrated points near the nose, and because points near the nose diverge faster than elsewhere, those points must be shifted by larger amounts (relative to the segment length) than points further away.

A shift of half the segment length at the nose, decreasing linearly to one eighth at about half the chord length gives the desired results. To test that profiles produced by this scheme are accurate, model profiles were also made with four times the optimum point density. There was no noticeable difference in the resulting plotted profiles.

D. Local Collision Efficiency

Parameterization schemes were derived in Chapter III for local collision efficiency β on circular cylinders as a scaled cosine function of local surface slope α . If the numerical results of Oleskiw (1982) for local collision efficiency on a NACA airfoil at 0° angle of attack are plotted against surface slope, they also closely follow a cosine curve, but one which is displaced negatively along

the abscissa. The form holds for several cases computed by Oleskiw, as seen in Figures V.2, V.3 and V.4, and is described by:

$$\beta(\alpha) = \cos(\alpha) - (1 - \beta_0), \quad V.3$$

where β_0 is the value of β at the stagnation line. This is the only parameter required to approximate the entire distribution, since the maximum accretion angle α_{\max} , in radians, follows from:

$$\alpha_{\max} = \arccos(1 - \beta_0). \quad V.4$$

Appendix IV gives a geometrical derivation of Equation V.3.

So far, a monodisperse droplet population has been assumed. In practice, of course, a spectrum of droplet sizes is present. Often the collision efficiency of the entire population is modelled by a single droplet size at the median volume diameter of the spectrum. This is generally a good approximation to the weighted average $\beta(\alpha)$ distribution calculated according to the relative volumes of the droplet sizes in the spectrum, as discussed in Appendix VI.

In real populations, when relatively large numbers of very small droplets are present in the population, they can change the shape of the β distribution, as well as force the mvd-based maximum accretion limit to a smaller value.

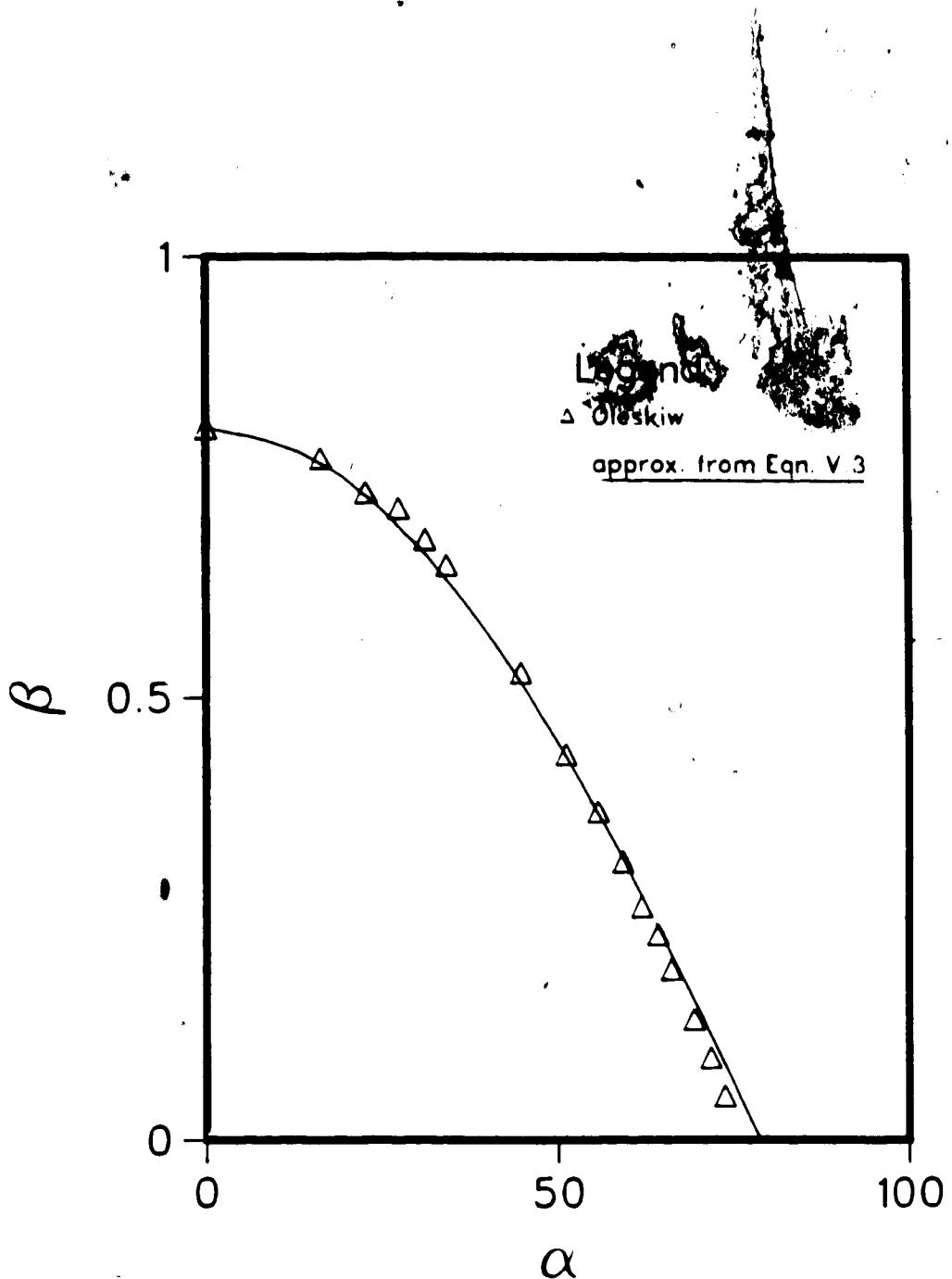


Figure V.2 - Comparison of Oleskiw's numerical results and Equation V.3, for collision efficiency of a single drop size. Airfoil is a NACA 0015, $U = 61 \text{ m/sec}$, Chord = 0.213 m, $D_d = 20 \mu\text{m}$.

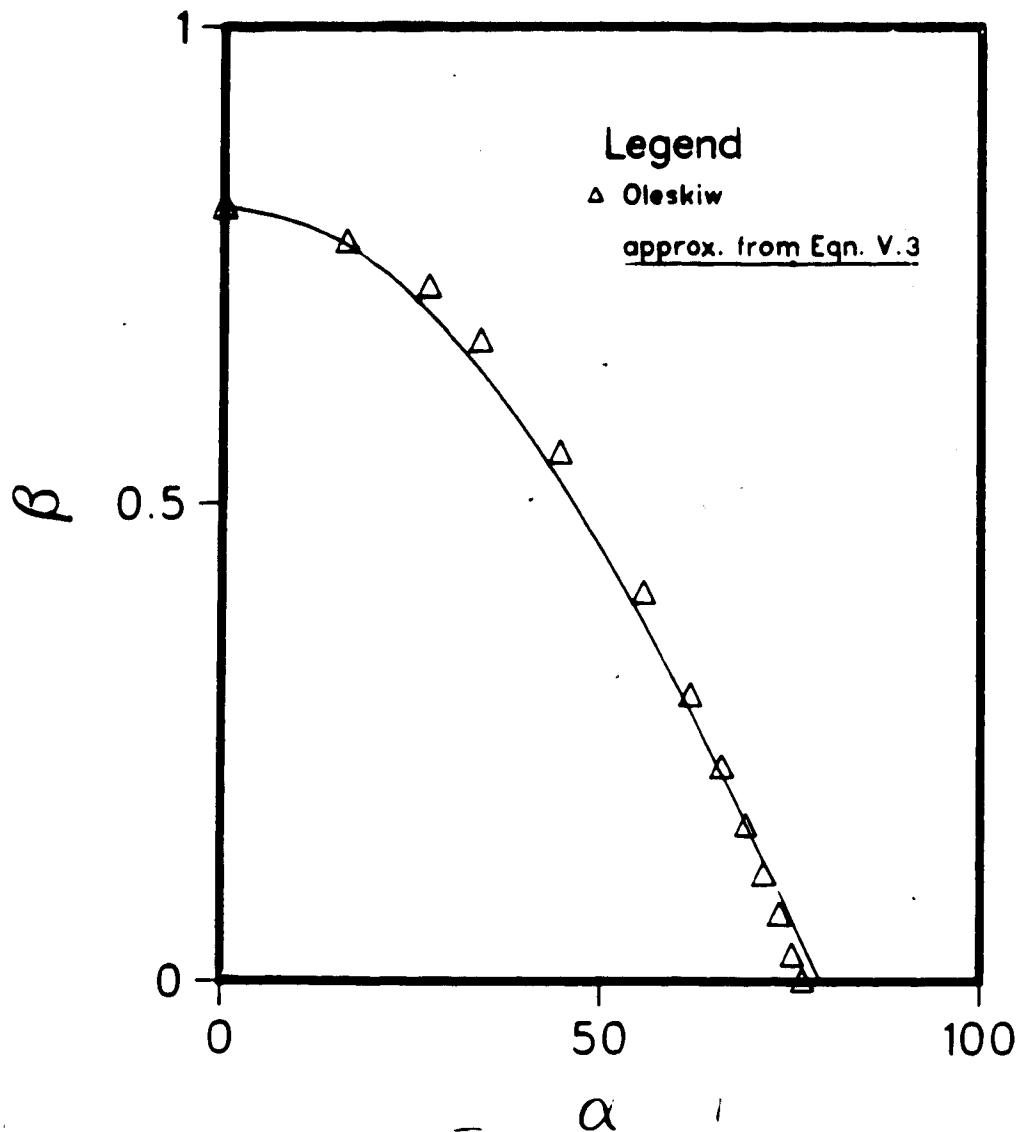


Figure V.3 - Comparison of Oleskiw's numerical results and Equation V.3, for collision efficiency of a single drop size. Airfoil is a Joukowski 15%, $U = 78.2$ m/sec, Chord = 0.33 m, $D_d = 25.5 \mu\text{m}$.

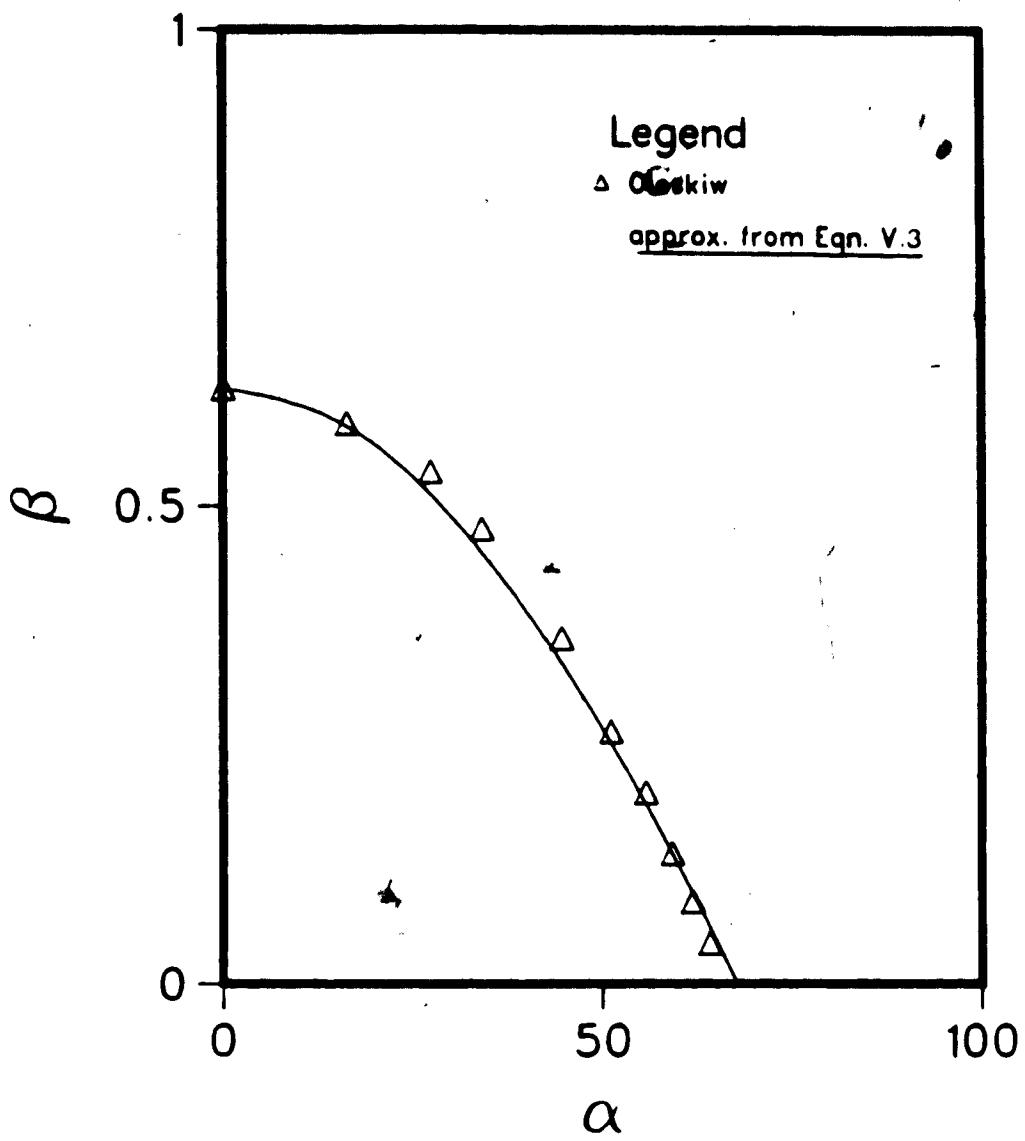


Figure V.4 - Comparison of Oleskiw's numerical results and Equation V.3, for collision efficiency of a single drop size. Airfoil is a Joukowski 15%, $U = 78.2$ m/sec, Chord = 0.33 m, $D_d = 13.2 \mu\text{m}$.

than actually occurs. For airfoils, the effect becomes important for β_0 less than about 0.8. For these cases, an alternative parameterization is proposed. It is derived from a polynomial fit to an average of several spectrum-weighted $\beta(\alpha)$ curves calculated by the author for droplet spectra observed in the FROST tunnel at the University of Alberta. The form of the parameterization is:

$$\beta(\alpha/\alpha_m) = \beta_0 (1 + 0.039(\alpha/\alpha_m) - 1.842(\alpha/\alpha_m)^2$$

$$- 0.543(\alpha/\alpha_m)^3 + 1.792(\alpha/\alpha_m)^4$$

V.5

$$- 0.444(\alpha/\alpha_m)^5,$$

where $\beta_0 = \beta_{0,mvd}$, and

$$\alpha_m = \alpha_{\max,mvd} + 0.13 \exp(-mvd / 70).$$

V.6

Here α_m is again in radians, and mvd is in μm . In Figures V.5 and V.6, the approximation of Equation V.5 is compared with spectrum-weighted distributions for two other observed droplet spectra, which appear in Table A.1 as "Battan and Reitan stratus" (Battan and Reitan, 1957), and "FROST 3", a spectrum measured in the FROST tunnel. These figures also show the difference in β_0 between the spectrum-weighted value and that derived from the mvd for these cases.

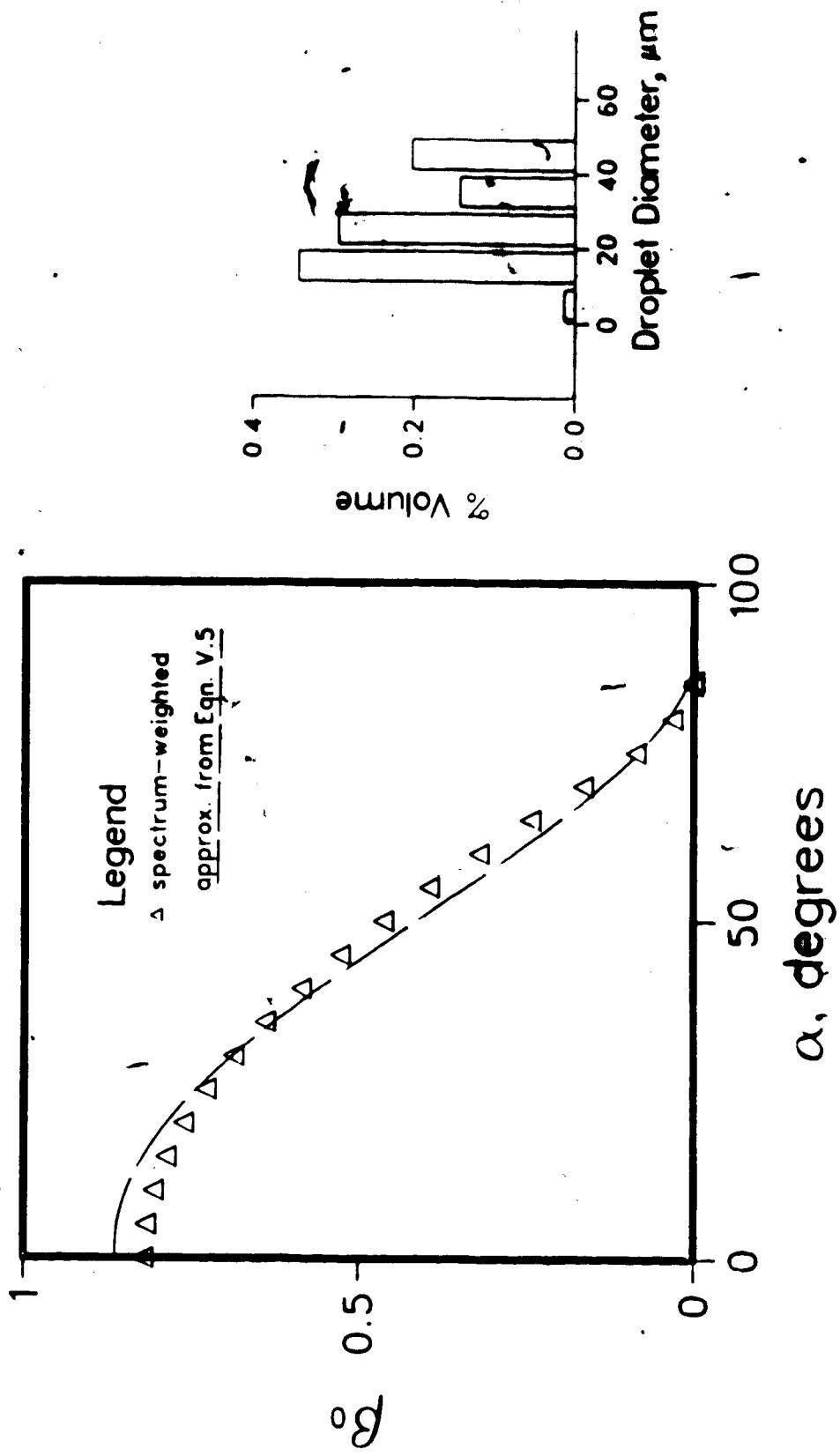


Figure V.5 - Comparison of spectrum-weighted collision efficiencies and Equation V.5 for the spectrum shown ($m_{vd} = 26.2 \mu\text{m}$) on a NACA 0015, $U = 50 \text{ m/sec.}$, chord = 0.2 m.

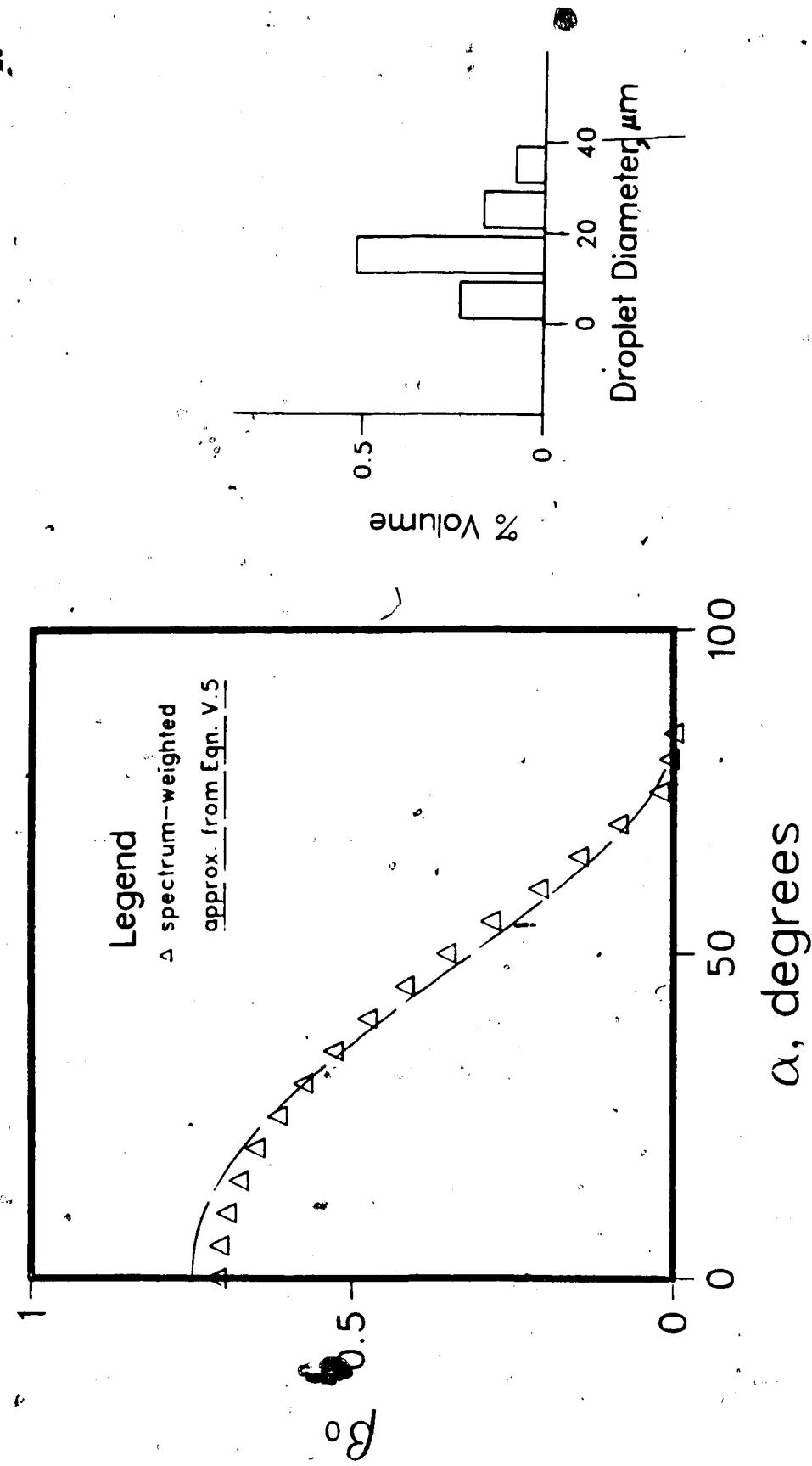


Figure V.6 - Comparison of spectrum-weighted collision efficiencies and Equation V.5 for the spectrum shown ($m_{vd} = 17.2 \mu\text{m}$) on a NACA 0015, $U = 50 \text{ m/sec}$, chord = 0.2 m.

When the angle of attack (Φ) is non-zero, Oleskiw's numerical results have been fitted with the following function:

$$\beta(\alpha') = 0.95 \beta_0 \cos(\alpha'), \quad V.7$$

where β_0 is calculated as for the $\Phi = 0$ case, and

$$\alpha' = ((\pi/2) (\alpha_0 - \alpha)) / (\pi/2 - \alpha). \quad V.8$$

The position of maximum collision efficiency has been designated by α_0 , and may be approximated by:

$$\alpha_0 = \pi \log_{10} ((180 \Phi/\pi) + 1) / (0.180 \exp(30.5 h_0)), \quad V.9$$

where α_0 and Φ are in radians.

Unfortunately, there are only two numerical cases currently available (from Oleskiw, 1982) to test this fit, as illustrated in Figures V.7 and V.8. Further calculations at some time in the future may help to refine the approximation. In the meantime, it does give a rough estimate for the $\beta(\alpha)$ distribution when Φ is non-zero, which is otherwise very costly to compute.

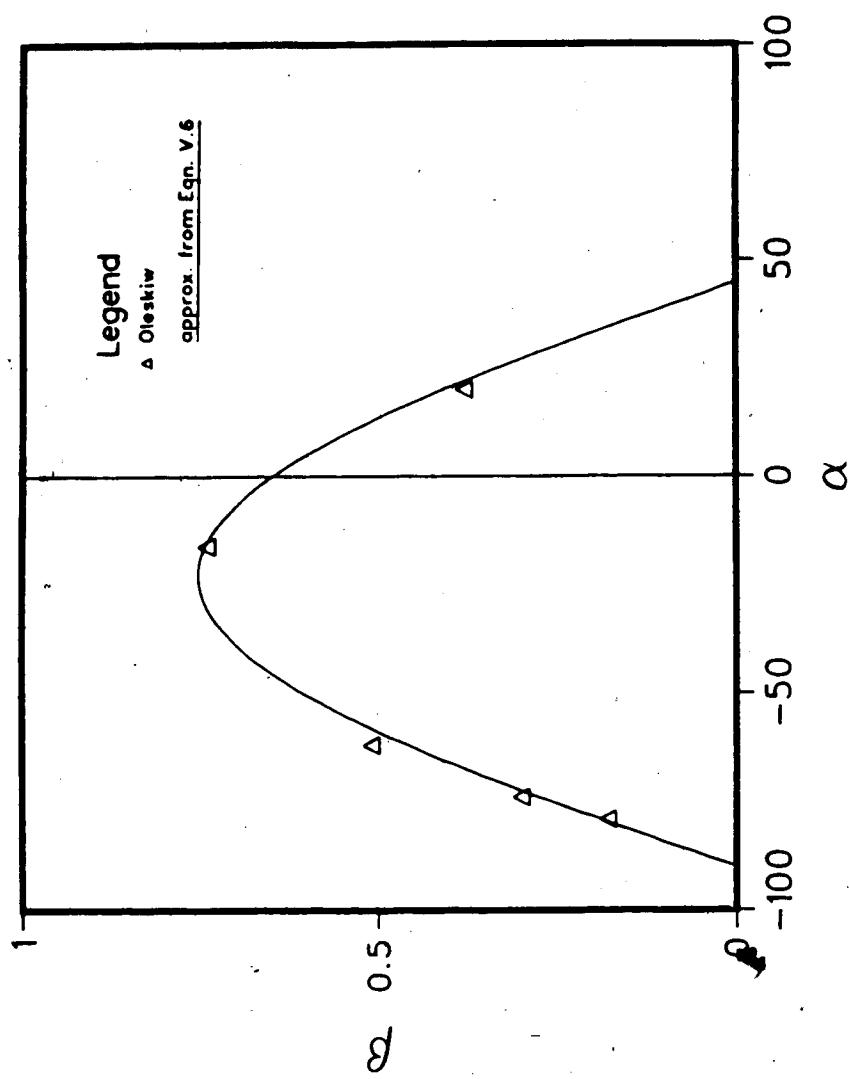


Figure V.7 - Comparison of Oleskiw's numerical results and Equation V.7, for collision efficiency on a NACA 0015 at an angle of attack = 8° , $U = 30.5$ m/sec, Chord = 0.169 m, $D_d = 20 \mu\text{m}$.

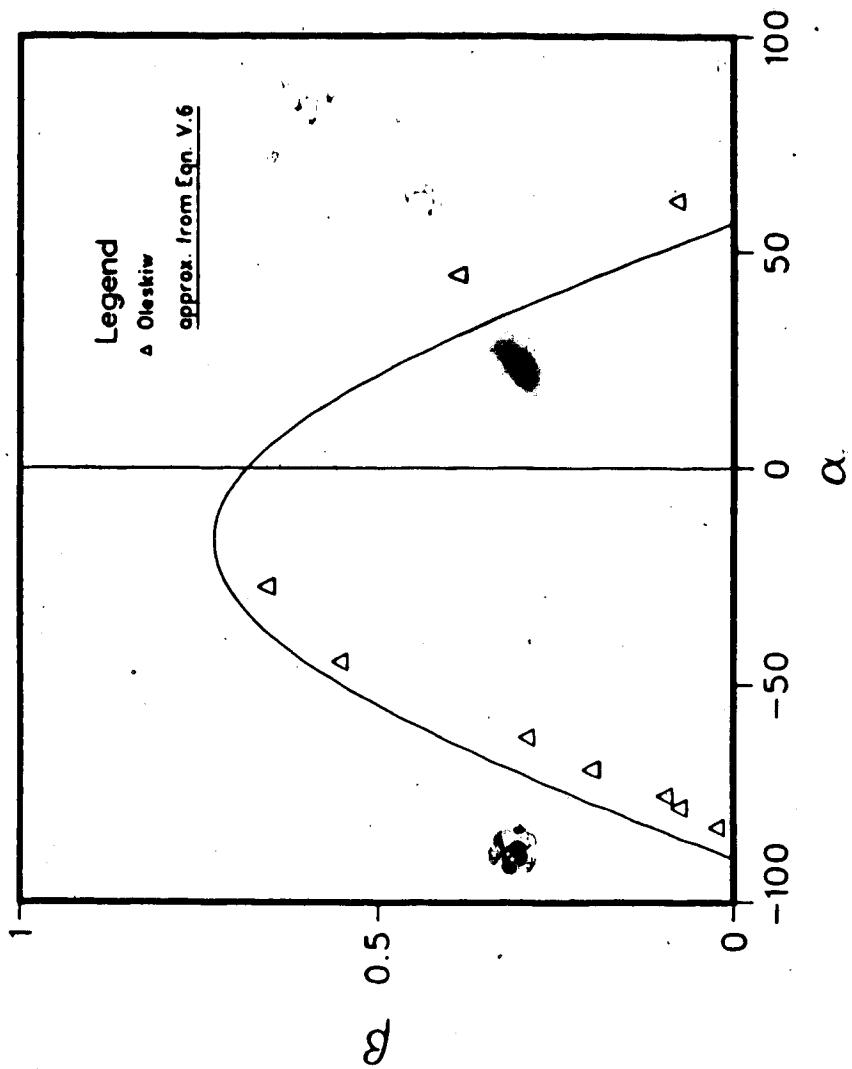


Figure V.8 - Comparison of Oleskiw's numerical results and Equation V.7, for collision efficiency on a Joukowski 15% at an angle of attack = 4° , $U = 78.2 \text{ m/sec}$, Chord = 0.33 m, $D_d = 18.6 \mu\text{m}$.

E. Integration of Droplet Trajectories

The equations of motion of a cloud droplet carried by the air stream, in a more general form than for the cylinder problem (Equation II.6), are:

$$\frac{dv_x}{dt} = -3/4 (\rho_a / \rho_d) (C_d/D_d) v (v_x - u_x) \quad V.10$$

$$\frac{dv_y}{dt} = -3/4 (\rho_a / \rho_d) (C_d/D_d) v (v_y - u_y)$$

In this steady-state formulation the additional term due to gravity has been neglected, as has the "history" term, arising from the droplet's acceleration with respect to the flow. Neglect of these terms is justified by the results of Oleskiw (1982), and also by examination of the "acceleration modulus", (Norment, 1980):

$$N_A = D_d |\frac{dv}{dt}| / v^2 \quad V.11$$

If this quantity is less than about 0.01, the history term may be safely neglected. The results of Oleskiw and of Norment showed that this limit is not exceeded in most aeronautical modelling situations.

The drag coefficient is again estimated from the expressions of Beard and Pruppacher (1969), Equation II.9. The droplet trajectories are integrated by the Heun method (Mesinger and Arakawa, 1976), similar to the calculations in

Chapter II, except for the determination of the air speeds u_x and u_y .

The streamfunction ψ of the potential flow field about the current accretion shape is numerically calculated following the method of Kennedy and Marsden (1976). This is a version of the surface singularity method originated by Hess and Smith (1967), in which the airfoil surface is represented by a set of point vortices. Kennedy and Marsden have slightly altered, at the trailing edge, the usual boundary condition of a constant streamfunction along the airfoil surface. They claim this gives a very accurate result with a minimum of computing time. Full details are available in their paper. The air speeds may then be approximated from a finite-difference calculation of the streamfunction derivatives, with the spatial differences equal to the droplet diameter.

The droplet's initial x-coordinate is five chord lengths upstream of the nose, and the initial separation of the two trajectories, Δy , is 10^{-8} metres. An optimum time step was found to be $\Delta t = 0.0025$ (U/chord). The justification for these choices is given by Table V.1, which examines the sensitivity of the model results to changes in these parameters.

As the droplet approaches the airfoil, the model begins testing for its position with respect to the current accretion surface. At each time step Δt , the array of accreting surface points is searched for the pair of

Δt (chord/U)	β_0 ($x_0 = 10 \cdot \text{chord}$)	β_0 ($x_0 = 5 \cdot \text{chord}$)
0.005	0.787	0.812
0.0033	0.794	0.804
0.0025	0.796	0.801
0.002	0.798	0.799

Table V.1 - Sensitivity of the integration results to the time step size Δt , and the initial x-position x_0 . Trajectory is for a 20 micron drop impinging on a NACA 0015 airfoil of 0.213 m chord, in a 61 m/sec airstream.

adjacent points which bracket the current y-coordinate of the droplet. The droplet's x-coordinate is then compared to the x-coordinates of these two points. If either or both of them are greater than the droplet's x-coordinate, the current droplet position lies inside the accretion surface.

The collision point is then estimated by solving for the intersection of the straight line segments, one joining the two bracketing surface points, and one joining the current and immediately previous droplet positions (see Figure V.9). The arc length $\Delta\ell$ separating the two collision points on the surface is also approximated as a straight line. β_0 is then just $\Delta y / \Delta\ell$.

All of these straight line approximations, and the simple integration method, are sufficiently accurate because of the very small distances between surface points near the nose, the small Δt and initial Δy of the trajectories.

A comparison of the integration results with those of Oleskiw for his Case 55 shows very good agreement. Oleskiw calculates $\beta_0 = 0.805$ for a 20 μm drop impinging on a chord NACA 0015, at 0° angle of attack and 61 m/s. The present model result is 0.801 for the same initial conditions. The small difference is easily accounted for by the use of a slightly different formulation for drag coefficient and by Oleskiw's method of approximating β_0 . This is from a spline fit to several $\beta(\alpha)$ points that are distributed along the surface, away from the stagnation line.

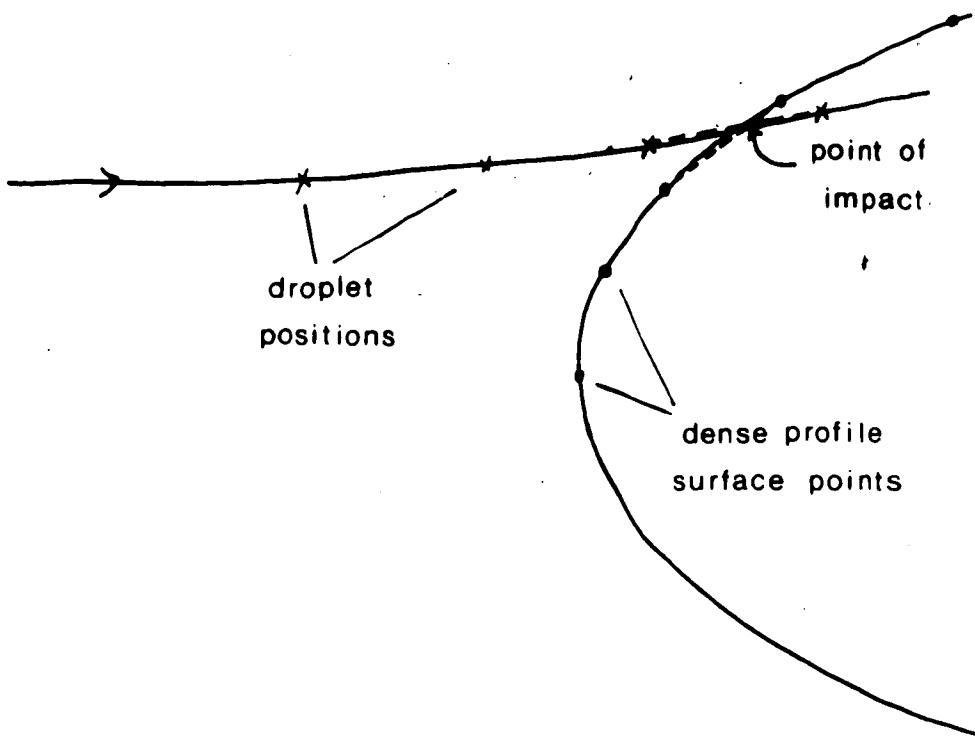


Figure V.9 - Determination of droplet impact point on the surface.

F. Local Density Variations

Once β_0 has been determined, α_{\max} and the distribution of local collision efficiency may be approximated from Equations V.3, V.5 or V.7. The local ice thickness (Th) accreted over a layer time step τ at each surface point is then given by the dry growth equation:

$$Th = \beta(\alpha) U_{lwc} \tau / \rho(\alpha), \quad V.12$$

where now a suitable distribution for $\rho(\alpha)$ must be found.

This would ideally be done using the same techniques that were used in Chapter IV, to investigate circular cylinder accretion densities by direct measurement of local densities and of layer thicknesses. However, the wind tunnel experiments on airfoils which were performed in the FROST tunnel for this study, were grown on airfoils with thicknesses less than one inch. This is too small to allow effective measurements by the hole drilling technique, and the layers outlined by paint spray for some of the experiments were too thick to allow good results using the layer method.

Therefore, an approximate distribution was derived by plugging the above-mentioned circular cylinder curve into the model, and adjusting it until it could produce reasonable profiles. The result works best when used with the 'simulated spectrum' parameterization for local

collision efficiency. We emphasize that this distribution is not yet supported by any empirical evidence, and that it should be adjusted further when such evidence is available. For the present however, a good estimate for the density distribution is:

$$\rho(a/a_m) = \rho_0 (1 - 0.163(a/a_m) + 0.235(a/a_m)^2$$

V.13

$$- 2.357(a/a_m)^3 + 1.374(a/a_m)^4 + 0.019(a/a_m)^5),$$

where ρ_0 is assumed, for simplicity, to be always 890 kg/m^3 . In certain conditions, the maximum density may be less than this value, but the model has no capability to determine this, at present.

The mass and mean density of each layer are calculated in the same way as for the circular cylinder described in Chapter III.

G. Model Results

Available data for model validation include the numerical results of Oleskiw, and a number of wind tunnel experiments carried out in the University of Alberta FROST tunnel, at NASA/Lewis and described in Olsekiw (1982b), at the National Aeronautical Establishment by Stallabrass (1958), and at the National Research Council Low Temperature Laboratory by Lozowski and Stallabrass (1978). In all of the experimental cases, the droplet sizes were determined by the

oil slide method, and have been corrected by us according to the calibration of Makkonen and Stallabrass (1984) discussed in Appendix III. An exception is the case taken from Stallabrass (1958), for which the liquid water content given by Stallabrass was derived from the observed ice thickness and the oil slide measurement of mvd. Stallabrass' numbers have thus been retained for both the droplet size and lwc. In all other cases, however, the accretion conditions were measured independently of the resulting thicknesses.

Figure V.10 illustrates the modelled profiles of Oleskiw and of the present model for identical accretion conditions. Since a constant local density was used, the agreement in shape of the layers between the two models reflects the accuracy of the parameterized collision efficiency distribution, for this case Equation V.3. The present model also lacks the surface instabilities which are already appearing in the Oleskiw model profile, arising from the spline curve approximation to the layer surface. Most importantly, the bottom profile of Figure V.10 represents a reduction in computational costs of a factor of ten over the Oleskiw model.

The conditions of Figure V.10 were chosen to agree with those of an experimental case grown in the NASA/Lewis tunnel. This experimental profile appears in Figure V.11, together with a corresponding model simulation using a constant local density, and one using the 'simulated spectrum' version of the local collision efficiency

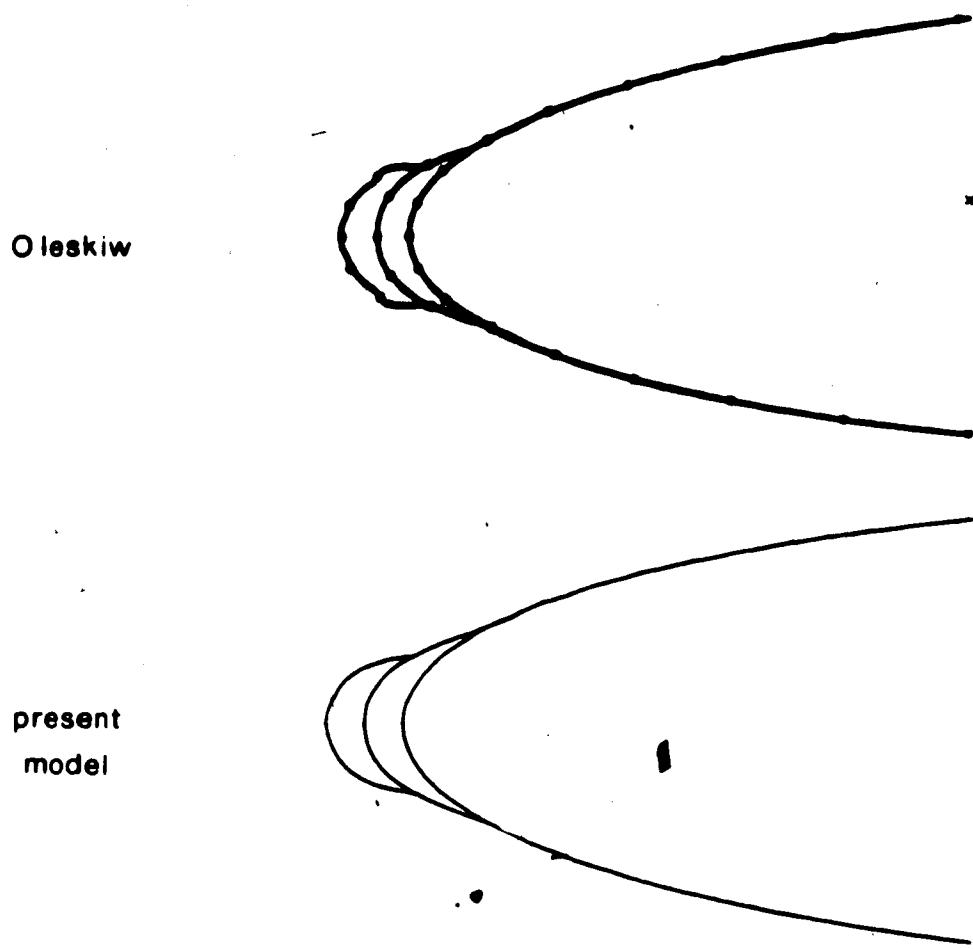


Figure V.10 - Comparison of present model results with those of Oleskiw (1982b) for a constant density accretion on a NACA 0015 at $\Phi = 0^\circ$, $U = 58 \text{ m/sec}$, $lwc = 1.02 \text{ g/m}^3$, chord = 0.533 m, $t = 5 \text{ min}$. A monodisperse droplet population at 12 μm diameter is assumed, and an air temperature of -15° C .

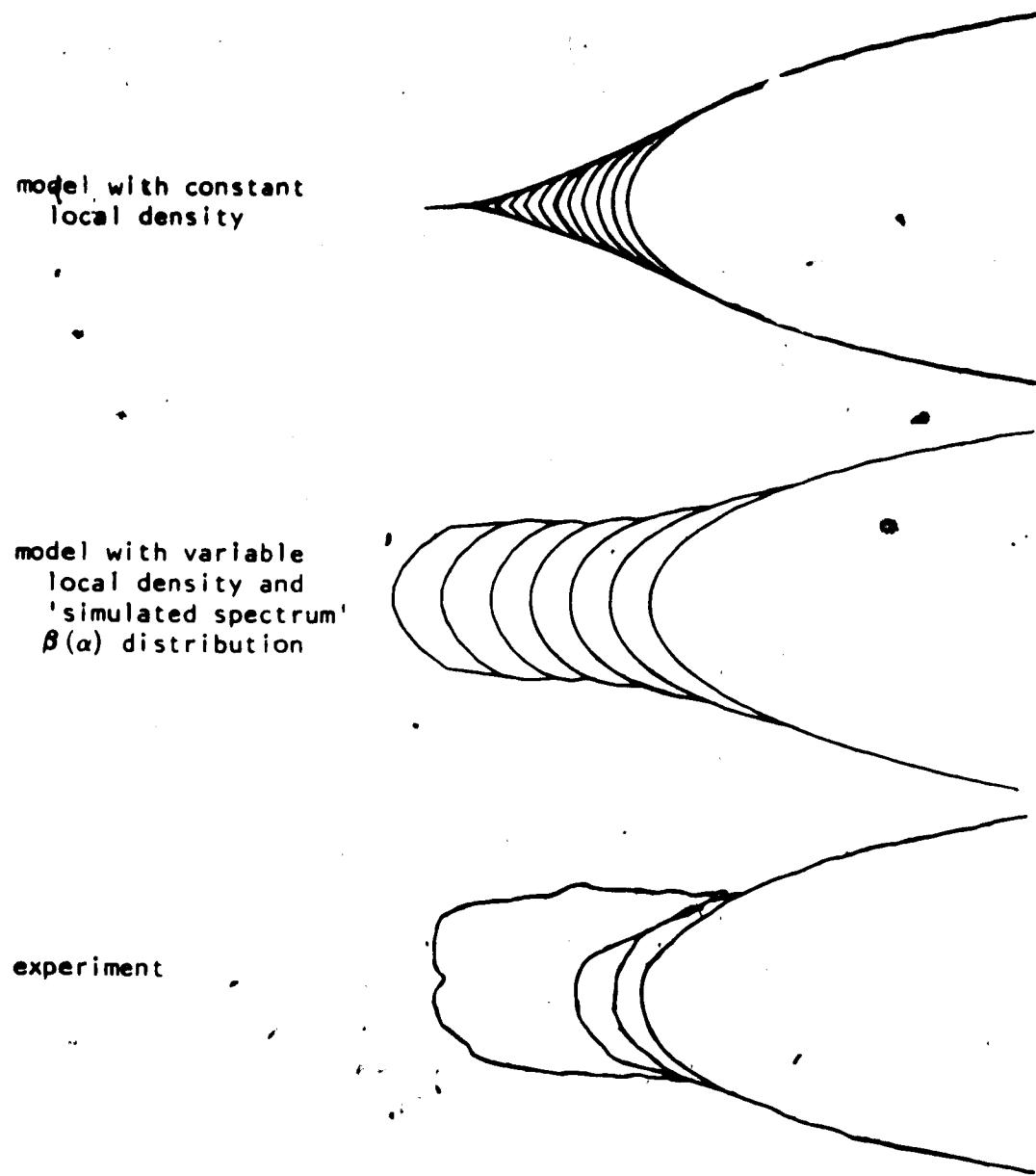


Figure V.11 - Comparison of present model results with an experimental accretion grown in the NASA/Lewis tunnel and described in Oleskiw (1982b), for the same conditions as in Figure V.10, but for a NACA 0012 airfoil, $t = 15$ min. Intermediate profiles at $t = 2$ and 5 minutes are also marked.

distribution, Equation V.5, together with the local density distribution of Equation V.12.

The two modelled profiles demonstrate a great improvement, in terms of a realistic accretion shape, in using the new local density and 'simulated spectrum' distributions over the old constant density and monodisperse spectrum versions. The model profiles are also seen to be free of surface instabilities and thus capable of accreting many more layers than the Oleskiw model.

Except for the forward surface of the experimental accretion, which seems to contain a small central trough, the agreement in shape is very good. Central troughs and ridges are features which are often seen to form on wind-tunnel accretions. They most probably arise from disturbances in the laminar flow pattern near the nose, as the ice layer alters its shape. The simple, potential flow model does not include such effects, and such features are not expected to show up in these simulations.

The experimental results of Lozowski and Stallabrass (1978) and of Stallabrass (1958) allow a test of Equation V.7, for local collision efficiency at non-zero angles of attack. Figure V.12 shows four experimental accretions and the corresponding model profiles, all including density variation and local collision efficiencies estimated for the $\Phi \neq 0$ case from Equation V.7. Since the method of approximating accretion limits for $\Phi \neq 0$ is somewhat crude,

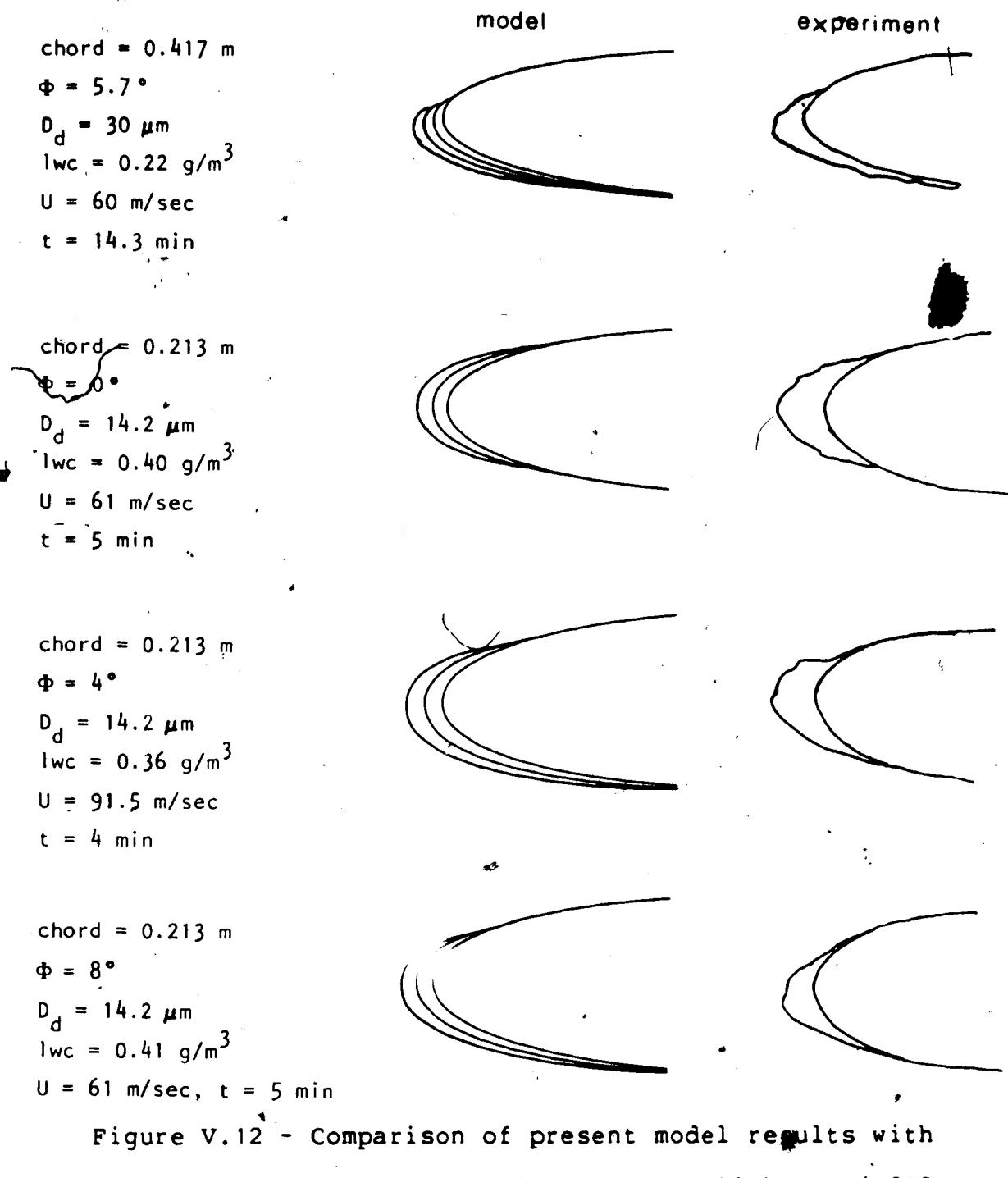


Figure V.12 - Comparison of present model results with experimental profiles from Stallabrass (1958, top profile, NACA 0012) and Stallabrass and Lozowski (1978, lower three profiles, Air temperature is -15° C.)

comparison in this region shows only rough agreement. However, the general agreement in shape and stagnation line thicknesses is encouraging. These cases at least illustrate that the approach taken in Equation V.9, to approximate the position of maximum collision efficiency α_0 , is worth pursuing and refining, especially given the large computational costs of determining this position from a series of droplet trajectory integrations.

For all of the cases mentioned so far, accretion masses have not been available. However, a series of wind-tunnel experiments were carried out in the University of Alberta FROST tunnel, for which was measured the mass of ice accreted on a 15 cm length of a NACA 0012 airfoil under various conditions. These results, with intermediate layers as marked by paint spray, are shown in Figure V.13 together with the modelled masses and profiles.

The differences between the model masses and the experimental masses for the two cases are 11 and 16 percent. This is of the same order as typical measurement errors for input parameters such as droplet size and liquid water content.

These examples have shown the usefulness of a simple approach toward numerical methods for ice accretion modelling. The present results are comparable to and much less costly than those of the Oleskiw model, and are of

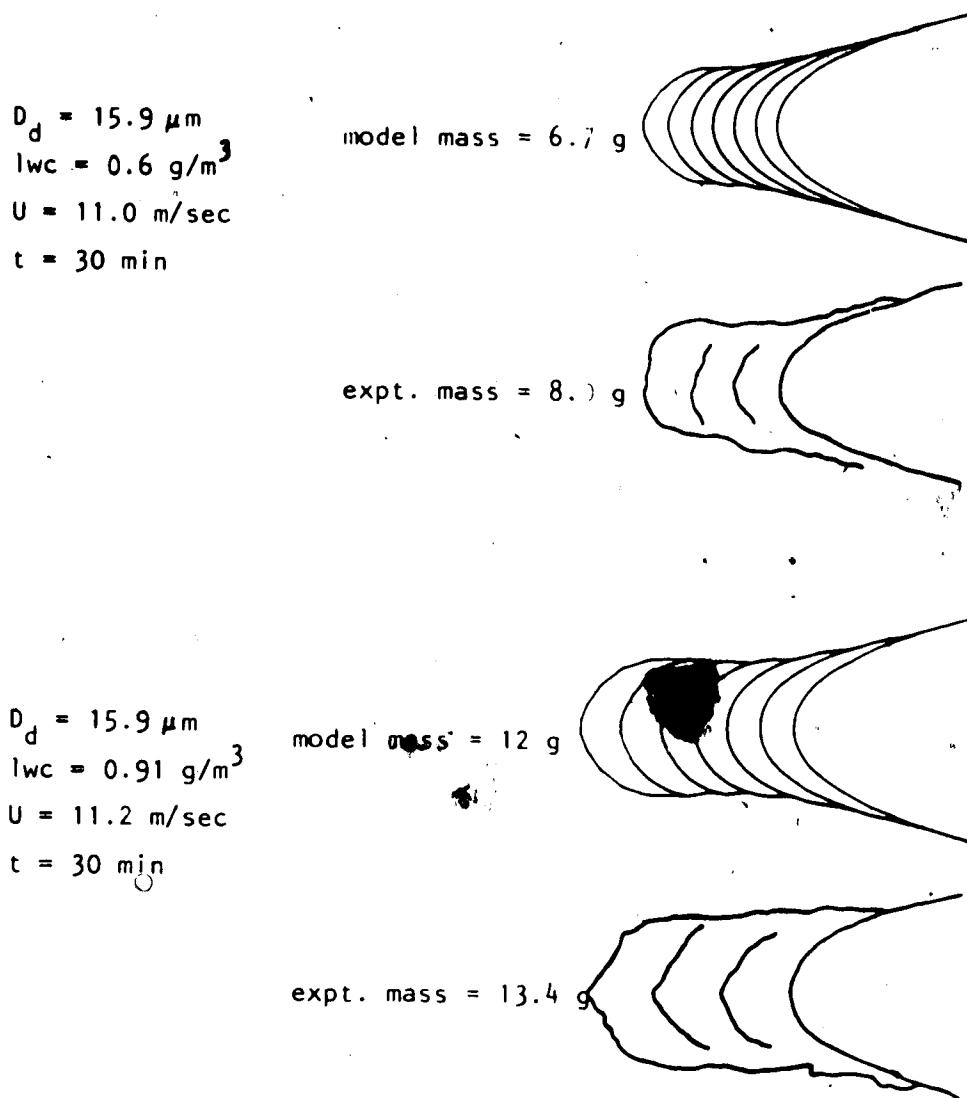


Figure V.13 - Comparison of present model results with experimental profiles and masses of accretions grown in The University of Alberta FROST tunnel on a NACA 0012 at $\Phi = 0^\circ$, chord = 0.176 m, air temperature = -10° C .

sufficient accuracy to allow practical and realistic comparisons with rime experiments. Such methods will also help to make the modelling problem for wet icing more tractable.

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VI. Concluding Remarks

In the four separate, but interrelated studies of this thesis several improvements have been put forward for physical models of ice accretion. Since conclusions were given within each chapter, these remarks will summarise the results, and give suggestions for future research.

Chapter II presented a re-calculation of the long-established work of Langmuir and Blodgett on collision efficiencies and other impingement parameters of circular cylinders. High-speed digital computer results for droplet trajectory integrations have provided a small improvement over the old analog computer results, but a larger improvement has been made in the application of a non-linear regression fit to the tabulated numerical values of collision efficiency, impact speed and accretion limit. This new analytical approximation could be very useful to modellers, who presently use a variety of functions to approximate the same thing, none of which is as accurate as the new expression.

These results are equally applicable to all types of icing, wet or dry, from transmission lines to marine structures, as long as it occurs on circular cylinders. A possibility for future work here is to extend the re-calculations to the spheres and flat ribbons also considered by Langmuir and Blodgett.

The circular cylinder rime model of Chapter III makes use of these new collision efficiencies, as well as of new

parameterizations for the distribution of local collision efficiency under various conditions. The efficiency and accuracy of these methods was well demonstrated in that chapter, by model results which reproduced realistic profiles, and accreted masses to within 10 percent of observation. The speed of the model; a few seconds on the University of Alberta Amdahl 580, and about 90 seconds on an IBM PC; make it highly practical.

The cylinder model also represents an improvement on older models in its version of local ice density distribution, derived in Chapter IV. Based on empirical data, this distribution is more realistic than previous versions, based on rotating cylinder correlations, have been. Since an attempt to validate the density distribution by direct measurement was inconclusive, further work, perhaps using different measurement techniques, would be valuable.

A logical extensions to the circular cylinder model would be the addition of gravitational and aerodynamic torque calculations in order to determine the correct rate of rotation. Another would be the determination of stagnation line density, following the method described in Chapter IV. The stagnation line density is required as a scaling parameter of the empirical density distribution. This would require a simple heat balance calculation at the stagnation line in order to determine the surface temperature there.

The airfoil model of Chapter V shows that parameterization methods similar to those of Chapter III may also be found for more difficult shapes, where the potential saving in computational costs is much larger. Here the parameterization for $\rho(\alpha)$ is in particular need of empirical validation, and the collision efficiency approximations would benefit from further verification and refinement using numerical results for a wider range of conditions.

The airfoil model has the capability to update the potential flow approximation, from which the β_0 parameter is derived, at intervals throughout the accretion period. Although a suggestion was made for the length of this interval, it was only an estimate. More numerical experiments should be carried out to discover the optimum interval, since this procedure is an expensive one.

Both models fail in the prediction of the troughs and ridges which sometimes form along the stagnation line of experimental accretions, often after the formation of a flat forward surface. An interesting exercise would be to investigate the flow patterns and droplet trajectories about these flat surfaces, in an attempt to understand how such features are formed. Experimental data should be gathered on the conditions under which troughs and ridges form, since they are not always present. Such studies may discover ways to model their formation.

Appendix I

Program Listing: GRAZE

LEVEL 1.1.1 (DEC 81)

VS FORTRAN DATE: JUL 24, 1986 TIME: 13:13:40

OPTIONS IN EFFECT: NOLIST NOXREF NOODSTANT MODECK SOURCE /TERM OBJECT FIXED
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This program calculates trajectory pairs for water droplets in an air flow field around a circular cylinder, and determines the stagnation line collision efficiency.

maximum accretion angle and overall collision efficiency. The air is assumed to be at a temperature of -10 degrees C, and a pressure of 100 kPa.

The equation of motion of the droplet is integrated using the Heun scheme (Ressinger, F. and Arakawa, A. 1976. Numerical Methods Used in Atmospheric Models, Vol. 1. W.M.O. GARP Publications Series No 17, 6pp.) The time step is set to 1/400 of the time to travel one cylinder radius at the free stream speed. A maximum of 7000 steps are allowed, by which time the droplet is assumed to have collided with the cylinder or to have passed it.

Note that the droplet's initial x-coordinate is set to be 10 cylinder radii upstream of the cylinder axis. The initial x-component speed is set equal to the air's x-component, and the y-component is set to half the air's y-component.

A first trajectory is integrated beginning at the input initial x-coordinate, which is then incremented by a very small amount (DELTAY) for the second trajectory. If both trajectories intercept the cylinder, the normal impact velocity, impact angle, impact position, and local collision efficiency are calculated.

The grazing trajectory is determined by successively halving the distance between a pair of initial y-coordinates which bracket the initial y-coordinate of the grazing trajectory. The initial values of y-min and y-max are input by the user; a judicious choice will decrease the number of trajectories needed to determine the grazing one.

If a Stokes flow calculation is requested (i.e. Langmuir phi parameter = 0), the drag coefficient is set to 24 / Reynolds number. Otherwise, the formulation of Beard and Pruppacher is used for the drag coefficient.

All operations are carried out in double precision. Output is written to a file named SOURCE.

To compile on enter:

\$RUN *FORTRAN5 SCARDS=GRAZERUN

To run, enter \$RUN GRAZERUN. You will be prompted for the following input parameters:

LEVEL 1.1.1 (DEC 81)

VS FORTRAN

DATE: JUL 24. 1986 TIME: 13:13:40

NAME: MAIN PAGE: 2

52 C Droplet DIAMETER in microns
53 C
54 C Cylinder RADIUS in metres
55 C
56 C Free stream air speed in metres per second
57 C
58 C Initial droplet y position in metres
59 C
60 C Initial value for y-min
61 C
62 C Initial value for y-max
63 C
64 C y(es) or n(o) for Stokes flow option
65 C
66 C Subprograms
67 C
68 C
69 C SUBROUTINE TRAJEC - calculates droplet trajectories
70 C by integrating the droplet equations of motion by
71 C the trapezoidal rule with one iteration step.
72 C
73 C SUBROUTINE ITERAT - calculates the air speeds (from
74 C the analytical flow solution) and droplet
75 C accelerations (from the equations of motion) for
76 C the current iteration and time step.
77 C
78 C SUBROUTINE COLLID - calculates final impact position
79 C if the droplet collides with the surface.
80 C
81 C
82 C Variable Dictionary
83 C
84 C ARC1, - arc lengths along cylinder surface
85 C ARC2, from stagnation line to the impact points
86 C of the two trajectories
87 C
88 C BETA, - local collision efficiency
89 C
90 C CYLRAD, - cylinder radius in metres (INPUT)
91 C
92 C DELTAY, - time step in seconds
93 C
94 C DELTAY, - difference in initial y coordinate of
95 C the two trajectories, in metres
96 C DROPO1, - droplet diameter, (input in microns, but
97 C converted to metres (INPUT))
98 C FINVX, - average values of the final velocities
99 C FINVY
100 C
101 C FINVX1, - x-components of droplet velocity at
102 C impact
103 C FINVX2
104 C FINVY1, - as above, for y-components
105 C FINVY2

LEVEL 1.1.1 (DEC 81)

VS FORTRAN

DATE: JUL 24. 1986 TIME: 13:13:40

NAME: MAIN PAGE:

107 C HALFP1 - PI / 2 (FORTRAN77 constant parameter
 108 C
 109 C HIT1, - logical flags for each trajectory
 110 C which indicate whether a collision
 111 C HIT2 with the cylinder has occurred
 112.1 C
 113. C NOTRAJ - no. of trajectories used to find grazing traj
 114. C
 115. C NSTEPS - maximum number of time steps to be
 116. C integrated
 116.1 C
 117. C PHI - angle of trajectory to the surface
 118. C normal at impact
 118.1 C
 119. C STK - input character variable to set Stokes
 120. C flow flag
 120.08 C
 120.5 C STOKES - logical flag for Stokes flow option
 120.7 C
 120.71 C THETA - average of THETA1 and THETA2
 120.8 C
 120.81 C THETA1, - position angles of the impact points
 120.87 C THETA2 of the two trajectories
 120.93 C
 121. C VEL - free stream air speed in metres/sec (INPUT)
 122. C
 122.1 C VNORM - droplet velocity normal to the cylinder
 122.4 C surface at impact
 122.7 C
 123. C XINIT - initial droplet x-coordinate in metres
 124. C
 125. C YINIT - initial droplet y-coordinate in metres (INPUT)
 126. C
 127. C YMIN, - minimum and maximum initial y-coordinates
 128. C YMAX which bracket the graze coordinate
 129. C
 130. C
 131. C
 132. C
 133. C
 134. C
 135. C
 136. C
 137. C
 138. C
 139. C DOUBLE PRECISION DROPO1,VEL,CYLRAD,XINIT,YINIT,FINVX,
 140. + FINVY,VNORM,PHI,THETA,THETA1,THETA2,
 141. + BETA,FINVX1,FINVX2,FINVY1,FINVY2,
 142. + ARC1,ARC2,DELTAT,DELTAY,YMIN,YMAX
 143. C
 144. C
 145. C
 146. C
 147. C
 148. C
 149. C
 150. C
 151. C
 152. C
 153. C
 ISN 2
 ISN 3
 ISN 4
 ISN 5
 ISN 6
 ISN 7
 154. C INTEGER
 155. C
 156. C LOGICAL
 157. C
 158. C CHARACTER*1
 159. C
 160. C
 161. C
 162. C Prompt user for input parameters
 163. C
 164. C WRITE (6,10)
 165. C FORMAT (//,***** Droplet trajectories about a'

+ Cylinder ***** //)

```

166.      C
167.      C      WRITE (6,*) 'Droplet DIAMETER to microns?'
168.      C      READ (5,*) DROPD1
169.      C      DROPD1 = DROPD1 * 1.0E-6
170.      C
171.      C      WRITE (6,*) 'Cylinder RADIUS in metres?'
172.      C      WRITE (6,*) 'Free stream air speed in m/sec?'
173.      C      READ (5,*) CYLRAD
174.      C
175.      C      WRITE (6,*) 'Initial droplet y-coordinate in metres?'
176.      C      WRITE (6,*) 'Initial value of ymin?'
177.      C      READ (5,*) VEL
178.      C
179.      C      WRITE (6,*) 'Initial value of ymax?'
180.      C      READ (5,*) YINIT
181.      C
182.      C      WRITE (6,*) 'Initial value of ymin?'
183.      C      READ (5,*) YMINT
184.      C      WRITE (6,*) 'Initial value of ymax?'
185.      C      READ (5,*) YMAX
186.      C
187.      C      WRITE (6,*) 'Do you wish a Stokes flow calculation? y/n?'
188.      C      READ (5,15) STK
189.      C      15 FORMAT (A1)
190.      C
191.      C      Set flag for Stokes flow case
192.      C
193.      C      IF ((STK .EQ. 'Y') .OR. (STK .EQ. 'Y')) THEN
194.      C          STOKES = TRUE
195.      C          READ (5,15) STK
196.      C          15 FORMAT (A1)
197.      C
198.      C      Set maximum number of time steps, initial x-coordinate, and
199.      C      the difference in initial y-coordinates increment number
200.      C      of trajectories
201.      C
202.      C      STOKES = FALSE
203.      C
204.      C      Initialize the total number of trajectory pairs
205.      C
206.      C      NOTRAJ = 0
207.      C      100 CONTINUE
208.      C
209.      C
210.      C      Set maximum number of time steps, initial x-coordinate, and
211.      C      the difference in initial y-coordinates increment number
212.      C      of trajectories
213.      C
214.      C      NOTRAJ = NOTRAJ + 1
215.      C      NSTEPS = 7000
216.      C      DELTAY = 1.00E-8
217.      C      XINIT = 10 * CYLRAD
218.      C
219.      C      Calculate time step
220.      C
221.      C

```

LEVEL 1.1.1 (DEC 81)

VS FORTRAN DATE: JUL 24, 1986 TIME: 13:13:40

NAME: MAIN PAGE: 5

```

222. C Integrate first trajectory
223. C
224. C CALL TRAJEC (XINIT,YINIT,DROPDI,VEL,CYLRAD,DELTAT,NSTEPS,
225. + STOKES,THETA1,ARC1,HIT1,FINVX1,FINVY1)
226. C
227. C Correct initial y position and integrate second trajectory
228. C
229. C YINIT = YINIT + DELTAY
230. C CALL TRAJEC (XINIT,YINIT,DROPDI,VEL,CYLRAD,DELTAT,NSTEPS,
231. + STOKES,THETA2,ARC2,HIT2,FINVX2,FINVY2)
232. C
233. C Test for result of trajectory (miss, graze or collision)
234. C
235. C IF((HIT1) .AND. (HIT2)) THEN
236. C
237. C Calculate impact parameters and local collision efficiency
238. C
239. C
240. C IF ((YINIT-DELTAY .NE. 0.0) YMIN = YINIT - DELTAY
241. C BETA = DELTAY / (ARC2 - ARC1)
242. C THETA = (THETA1 + THETA2) / 2.0
243. C FINVX = (FINVX1 + FINVX2) / 2.0
244. C FINVY = (FINVY1 + FINVY2) / 2.0
245. C PHI = HALFPI - DATAN(FINVX/FINVY) + THETA
246. C VNORM = DSQRT((FINVX * FINVX) + (FINVY * FINVY))
247. C
248. C Write results to screen
249. C
250. C
251. C WRITE (6,20) THETA,VNORM,PHI,BETA
252. C 20 FORMAT (//, Both trajectories collided with cylinder //,
253. C //, Position angle at impact point: ,F7.3
254. C //, Velocity normal to surface at impact: ,F7.3
255. C //, Angle of trajectory to surface normal: ,F7.3
256. C //, Local collision efficiency: ,F5.3 //)
257. C ELSE IF (HIT1) THEN
258. C WRITE (6,30) THETA1
259. C 30 FORMAT (//, First trajectory grazed cylinder //,
260. C //, Surface angle of max accretion point: ,F7.3 //)
261. C
262. C WRITE (6,*), overall coll. eff.: ,(YINIT - DELTAY)/CYLRAD
263. C WRITE (6,*), no of trajectories: ,NOTRAJ
264. C GO TO 200
265. C
266. C YMAX = YINIT -DELTAY
267. C 40 WRITE (6,40)
268. C FORMAT (//, Both trajectories missed the cylinder //),
269. C
270. C Determine new initial position and start next trajectory pair
271. C
272. C 67 YINIT = (YMAX + YMIN) / 2.0
273. C 68 GO TO 100
274. C
275. C 69 200 STOP
276. C END

```

LEVEL 1.1.1 (OEC 81) VS FORTRAN

DATE: JUL 24, 1986 13:13:40

NAME: MAIN PAGE: 6

STATISTICS SOURCE STATEMENTS = 69. PROGRAM SIZE = 2732 BYTES. PROGRAM NAME = MAIN PAGE: 1.

STATISTICS NO DIAGNOSTICS GENERATED.

***** END OF COMPILATION 1 *****

LEVEL 1.1.1 (DEC) 81)

DATE: JUL 24. 1986 TIME: 13:13:41

NAME: MAIN PAGE: 7

OPTIONS IN EFFECT: NOLIST NOMAP NOXREF NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(0) LANGVL(77) NOTIPS FLAG(') NAME(MAIN) LINECOUNT(60)

277 C *
278 C *
279 C *
280 C *
281 C * Integrates trajectory of droplet whose initial
282 C * position is specified in the parameter list. Calls
283 C * ITERAT for iterative calculations, and COLLID to
284 C * determine impact parameters for colliding trajectories.
285 C *
286 C *
287 C * INPUT VARIABLES:
288 C * XINIT - initial x position
289 C * YINIT - initial y position
290 C * DROPD1 - droplet diameter
291 C * CYLRAD - cylinder radius
292 C * VEL - free stream ~~VE~~ speed
293 C * DELTAT - time step
294 C * NSTEPS - maximum number of time steps
295 C * STOKES - logical flag for Stokes flow
296 C *
297 C * OUTPUT VARIABLES:
298 C * THETA - surface angle of impact
299 C * ARC - arc length along cylinder surface
300 C * from stagnation line to impact point
301 C * HIT - logical flag set to TRUE if
302 C * a collision occurs
303 C * FINVX - droplet x velocity on impact
304 C * FINVY - droplet y velocity on impact
305 C *
306 C * OTHER VARIABLES:
307 C * X,Y - arrays containing x,y positions
308 C * for each time step
309 C * VX,VY - arrays containing x,y components
310 C * of droplet speed for each step
311 C * DVX, - arrays containing x,y components
312 C * of droplet accelerations
313 C *
314 C *
315 C *
316 C *
317 C *
318 C *
ISN 1 319 C *
ISN 2 320 C *
ISN 3 321 C *
ISN 4 322 C *
ISN 5 323 C *
324 C * INTEGER
325 C * LOGICAL
326 C *
327 C *
328 C *
329 C *
SUBROUTINE TRAJEC(XINIT,YINIT,DROPD1,VEL,CYLRAD,DELTAT,
NSTEPS,STOKES,THETA,ARC,HIT,FINVX,FINVY)
DOUBLE PRECISION X(7000),Y(7000),
VX(7000),VY(7000),DVX(7000),DVY(7000)
DOUBLE PRECISION XINIT,YINIT,DROPD1,CYLRAD,VEL,DELTAT,
THETA,ARC,FINVX,FINVY
NSTEPS
HIT,STOKES
Initialize values for integration

LEVEL 1.1.1 (DEC 81)

VS FORTRAN DATE: JUL 24 P 1986 TIME: 13:13:41 NAME: TRAJEC PAGE: 8

```

      C   X(1) = -XINIT
      C   Y(1) = YINIT
      C   VY(1) = (DABS(2 * VEL * CYLRAD * CYLRAD * X(1) * Y(1)) /
      C   + ((X(1) * X(1)) + (Y(1) * Y(1))) * 2.0) / 2.0
      C   VX(1) = VEL
      C   I = 1
      C   HIT = .FALSE.

 338.  C   Begin integration
 339.  C
 340.  C   100 CONTINUE
 341.  C
 342.  C   Calculate current step air speeds and accelerations
 343.  C
 344.  C   CALL ITERAT(X(I),Y(I),VX(I),VY(I),VEL,CYLRAD,DROPDI,
 345.  C   + STOKES,DVX(I),DVY(I))
 346.  C
 347.  C   Calculate next step droplet speeds and positions
 348.  C
 349.  C   VX(I+1) = VX(I) - (DELTAT * DVX(I))
 350.  C   VY(I+1) = VY(I) + (DELTAT * DVY(I))
 351.  C   X(I+1) = X(I) + (VX(I) * DELTAT) - (0.5 * DVX(I))
 352.  C   + (VY(I) * DELTAT)
 353.  C   Y(I+1) = Y(I) + (VY(I) * DELTAT) + (0.5 * DVY(I))
 354.  C   + (VX(I) * DELTAT)
 355.  C
 356.  C   Calculate next step air speeds and accelerations
 357.  C
 358.  C   CALL ITERAT(X(I+1),Y(I+1),VX(I+1),VY(I+1),VEL,
 359.  C   + CYLRAD,DROPDI,STOKES,DVX(I+1),DVY(I+1))
 360.  C
 361.  C   Iterate calculations of next step droplet speed and
 362.  C   position using averaged values of accelerations
 363.  C
 364.  C   VX(I+1) = VX(I) - (DELTAT / 2.0) * (DVX(I) +
 365.  C   + DVX(I+1))
 366.  C   VY(I+1) = VY(I) + (DELTAT / 2.0) * (DVY(I) +
 367.  C   + DVY(I+1))
 368.  C   X(I+1) = X(I) + (VX(I) * DELTAT) - (0.25 *
 369.  C   + (DVX(I) + DVX(I+1)) * (DELTAT * DELTAT))
 370.  C   Y(I+1) = Y(I) + (VY(I) * DELTAT) + (0.25 *
 371.  C   + (DVY(I) + DVY(I+1)) * (DELTAT * DELTAT))
 372.  C
 373.  C   Test whether trajectory has intersected the cylinder
 374.  C
 375.  C   IF (DSQRT((X(I+1)**2 + (Y(I+1)**2) LE CYLRAD) THEN
 376.  C     HIT = .TRUE.
 377.  C     FINVX = (VX(I) + VX(I+1)) / 2.0
 378.  C     FINVY = (VY(I) + VY(I+1)) / 2.0
 379.  C     CALL COLLIDIX(I,X(I+1),Y(I),Y(I+1),CYLRAD,
 380.  C                   THEITA,ARC)
 381.  C
 382.  C   ELSE IF (I .LT. (NSTEPS - 1)) THEN
 383.  C     I = I + 1
 384.  C     GO TO 100
 385.  C

```

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:13:41 NAME: TRAJEC PAGE: 9
*.....1.....2.....3.....4.....5.....6.....7.....8.....9.....

ISN 366 C RETURN
ISN 32 367
ISN 33 368 END

STATISTICS SOURCE STATEMENTS = 33, PROGRAM SIZE = 338354 BYTES, PROGRAM NAME = TRAJEC PAGE: 7.

STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPIRATION 2 *****

LEVEL 1.1.1 (DEC 81)

VS FORTRAN

DATE: AUG 07, 1986 TIME: 10:44:54

OPTIONS IN EFFECT:

NOLIST NOMAP NOXREF NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(O) LANG,LVL(77) NOFIPS FLAG(I) NAME(MAIN) LINECOUNT(60)

NAME: MAIN PAGE: 10

```

389. C *****
390. C *****
391. C * SUBROUTINE ITERAT - calculates for the current step:
392. C * air speed components from analytical solution to
393. C * flow around a cylinder, the droplet Reynolds
394. C * number, drag coefficient according to the formula-
395. C * tion of Beard and Pruppacher, and the droplet
396. C * acceleration components from the equations of motion.
397. C *****
398. C * INPUT VARIABLES :
399. C   X,Y      - current droplet positions
400. C   VX,VY.   - current droplet speeds
401. C   VEL      - free stream air speed
402. C   CYLRAD   - cylinder radius
403. C   DROPD1   - droplet diameter
404. C   STOKES   - logical flag for Stokes flow
405. C *****
406. C * OUTPUT VARIABLES :
407. C   DVX,DVY - current droplet accelerations
408. C *****
409. C * OTHER VARIABLES :
410. C   UX,UY    - current air speed components
411. C   DROPRE   - droplet Reynolds number
412. C   DRAGCO   - drag coefficient
413. C   COEFF    - combined coefficients of the
414. C   equations of motion
415. C   DENOM    - denominator of air speed
416. C   EQUATIONS - equations
417. C *****
418. C *****
419. C   SUBROUTINE ITERAT(X,Y,VX,VY,VEL,CYLRAD,DROPD1,STOKES,DVX,DVY)
420. C   ISN      1
421. C   421. DOUBLE PRECISION X,Y,VX,VY,DVX,DVY,VEL,CYLRAD,DROPD1,
422. C   422. DROPRE,DRAGCO,COEFF,UX,UY,DENOM
423. C   423. C
424. C   424. LOGICAL
425. C   425. STOKES
426. C   426. C
427. C   427. C Calculate air speeds
428. C   428. C
429. C   429. DENOM = ((X * X) + (Y * Y)) ** 2
430. C   430. UX   = VEL * ((CYLRAD**2) * (((Y * Y) - (X * X)) / DENOM))
431. C   431. UV   = DABS(2 * VEL * ((CYLRAD**2) * ((X * Y) / DENOM)))
432. C   432. C
433. C   433. C Calculate Reynolds number and drag coefficient
434. C   434. DRPRE = DROPD1 * DSQRT(((VX - UX) ** 2) + ((VY - UY)
435. C   435. * 2)) / 1.262E-5
436. C   436. +
437. C   437. IF (STOKES) THEN
438. C   438. DRAGCO = 24.0 / DRPRE
439. C   439. ELSE IF (DROPRE LT 2.0) THEN
440. C   440. DRAGCO = (24.0 / DRPRE) + (2.45 / (DRPRE ** 0.045))
441. C   441. ELSE IF (DROPRE LT 21.0) THEN

```

```

LEVEL 1.1.1 (DEC 81)          VS FORTRAN          DATE: AUG 07. 1986   TIME: 10:44:54    NAME: ITERAT PAGE: 11
                                         1.....2.....3.....4.....5.....6.....7.....8.....9

ISN   13     442.          DRAGCO = (24.0 / DROPRE) + (2.76 / (DROPRE ** 0.198))
ISN   14     443.          ELSE IF (DROPRE .GT. 200.0) THEN
ISN   15     444.          DRAGCO = (24.0 / DROPRE) + (4.73 / (DROPRE ** 0.371))
ISN   16     445.          + (6.24E-3 * (DROPRE ** 0.38))
ISN   17     446.          ELSE
ISN   18     447.          DRAGCO = (24.0 / DROPRE) + (4.536 / (DROPRE ** 0.368))
ISN   19     448.          ENDIF
ISN   20     449.          C calculate accelerations
ISN   21     450.          C
ISN   22     451.          COEFF = 9.92E-4 * (DRAGCO / DROPRE)
ISN   23     452.          DVX = COEFF * DSQRT(((VX - UX) ** 2) + ((VY - UY) ** 2))
ISN   24     453.          + (VX - UX)
ISN   25     454.          DVY = -COEFF * DSQRT(((VX - UX) ** 2) + ((VY - UY) ** 2))
ISN   26     455.          + (VY - UY)
ISN   27     456.          RETURN
ISN   28     457.          END
ISN   29     458.          C
ISN   30     459.          C
ISN   31     460.          C
ISN   32     461.          C

*STATISTICS* SOURCE STATEMENTS = 23. PROGRAM SIZE = 1380 BYTES. PROGRAM NAME = ITERAT PAGE: 10.
*STATISTICS* NO DIAGNOSTICS GENERATED.
***** END OF COMPILEATION 3 *****

```

LEVEL 1.1.1 (DEC 81)

DATE: JUL 24, 1986 TIME: 13:13:41

NAME: MAIN PAGE: 12

OPTIONS IN EFFECT: NOLIST NOMAP NOXREF NODECK
OPTIMIZE(O) LANGVL(77) NOTLIPS FLAG(I) NAME(MAIN) LINECOUNT(60)

NOTEST SEQ

462 C *
463 C *
464 C * SUBROUTINE COLLID - Calculates the x,y position
465 C * of impact on the cylinder surface, position angle
466 C * of the impact point, and arc length measured from
467 C * the stagnation line.
468 C *
469 C * INPUT VARIABLES:
470 C * XOUT, - last droplet position outside
471 C * YOUT the cylinder
472 C * XIN, - first droplet position inside
473 C * YIN the cylinder
474 C * CYLRAD - cylinder radius
475 C *
476 C * OUTPUT VARIABLES :
477 C * THETA - position angle of impact point
478 C * on the cylinder measured from
479 C * the stagnation line
480 C * ARC - arc length from stagnation line
481 C * to the impact point
482 C *
483 C * OTHER VARIABLES :
484 C * SLOPE - slope of the trajectory segment
485 C * which crosses the cylinder surface
486 C * i.e., between XOUT,YOUT and
487 C * XIN,YIN
488 C * A,B,C - coefficients of the quadratic
489 C * equation to solve for the intercept
490 C * of the trajectory and cylinder
491 C * surface
492 C * IMPX, - impact position on the surface
493 C * IMPY
494 C *
495 C *
496 C *
497 C *
498 C *
499 C * DOUBLE PRECISION XOUT,XIN,YOUT,YIN,CYLRAD,THETA,ARC,
500 C *
501 C *
502 C * Calculate slope of trajectory at impact point
503 C *
504 C * SLOPE = (YIN - YOUT) / (XIN - XOUT)
505 C * XOUT = XOUT / CYLRAD
506 C * XIN = XIN / CYLRAD
507 C * YOUT = YOUT / CYLRAD
508 C * YIN = YIN / CYLRAD
509 C *
510 C * Find impact point by solving quadratic equation with
511 C * coefficients as follows
512 C *
513 C * A = (SLOPE * SLOPE) + 1.0
514 C * B = (2.0 * SLOPE * YOUT) - (2.0 * (SLOPE * SLOPE) * XOUT)
ISN 1
ISN 2
ISN 3
ISN 4
ISN 5
ISN 6
ISN 7
ISN 8
ISN 9

LEVEL 1 1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:13:41 NAME: COLLID PAGE: 13

ISN 10 515 C -(2.0 * SLOPE * XOUT * YOUT) + ((SLOPE * SLOPE) *
 516. + (XOUT * XOUT) + (YOUT * YOUT) - 1.0
 ISN 11 517 IMPX = (-B - DSQRT((B * B) - (A * A * C))) / (2 * A)
 ISN 12 518 IMPY = (SLOPE * (IMPX - XOUT)) + YOUT
 ISN 519 C calculate impact point angle and arc length
 520. C
 ISN 521 C
 ISN 13 522 THETA = DATAN (IMPY / DABS(IMPX))
 ISN 14 523 ARC = CYLRAD * THETA
 ISN 524 C
 ISN 15 525 C
 ISN 16 526 C
 RETURN
 END

STATISTICS SOURCE STATEMENTS = 16, PROGRAM SIZE = 840 BYTES, PROGRAM NAME = COLLID PAGE: 12.

STATISTICS NO DIAGNOSTICS GENERATED.

***** END OF COMPILEATION 4 *****

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:13:42 NAME: MAIN PAGE: 14

SUMMARY OF MESSAGES AND STATISTICS FOR ALL COMPILENTIONS

STATISTICS SOURCE STATEMENTS = 69. PROGRAM SIZE = 2732 BYTES. PROGRAM NAME = MAIN PAGE: 1

STATISTICS NO DIAGNOSTICS GENERATED.

***** END OF COMPILATION 1 *****

STATISTICS SOURCE STATEMENTS = 33. PROGRAM SIZE = 338354 BYTES. PROGRAM NAME = TRAJEC PAGE: 7

STATISTICS NO DIAGNOSTICS GENERATED.

***** END OF COMPILATION 2 *****

STATISTICS SOURCE STATEMENTS = 23. PROGRAM SIZE = 1380 BYTES. PROGRAM NAME = ITERAT PAGE: 10

STATISTICS NO DIAGNOSTICS GENERATED.

***** END OF COMPILATION 3 *****

STATISTICS SOURCE STATEMENTS = 16. PROGRAM SIZE = 840 BYTES. PROGRAM NAME = COLLID PAGE: 12

STATISTICS NO DIAGNOSTICS GENERATED.

***** END OF COMPILATION 4 *****

***** SUMMARY, STATISTICS ***** 0 DIAGNOSTICS GENERATED. HIGHEST SEVERITY CODE IS 0

Appendix II

Program Listing: OMNICYL

OPTIONS IN EFFECT: MOLIST NOXREF NOGOSTMT NODECK

VS FORTRAN

DATE: JUL 24, 1986 TIME: 13:15:27

PAGE: 1

PAGE: 1

OPTIMIZE(0) LANGLV(77) NOFIPS FLAG(I) NAME(MAIN) LINCOUNT(50)

NOTE#

SEQ

1 1 2 3 4 5 6 7 8 9

***** Cylinder Rime Icing Model *****

Karen J. Finstad

June, 1986

This program will calculate and plot rime ice accretion layers on a cylinder, as well as calculate the mass and mean density of each layer. Local collision efficiencies for each point on the surface are calculated as a cosine function of the current local surface angle and scaled to the value of stagnation line collision efficiency (Beta zero), and maximum accretion angle (Alpha max).

Two forms of this function are used for droplet diameters above and below 20 microns. Beta zero and Alpha max are estimated from a non-linear regression fit to data computed by droplet trajectory program GRAZE.

Two optional procedures are available for treating the effects of droplet size spectra, when the assumption of a monodisperse population at the median-volume diameter (mvd) becomes inaccurate (roughly, when the mvd is less than 50 micron, for ground level wind speeds):
1. the user may enter weighting factors for each of 16 size bins of width 10 microns; or
2. use a "simulated spectrum" curve for the local collision efficiency (l.c.e.), which differs principally from the cosine approximation by the addition of a "tail" onto the maximum accretion limit. The length of this tail is a function of the mvd. Again, there is a slightly different form of the curve for mvd above and below 20 microns.

The first of these options is recommended only when three or less layers are to be accreted, since the weighted averaging procedure eventually produces first-order discontinuities in the l.c.e. curve, and thus on the modelled ice surface.

Local density is calculated from an empirically-derived function of local surface slope

The cylinder may be rotated clockwise, after each layer is accreted. The rotation increment is included in the input parameter list. When this option is chosen, (i.e., if the input rotation increment is greater than zero), the l.c.e. and local density are calculated as a function of position angle, instead of surface angle. This somewhat less accurate method is necessary to avoid problems when accreting over top of the previous layers.

Also please note that rotation of more than 90 degrees in total is not possible in this version

51
50

52. C The output consists of a plot description file written into
 53. C "pdf" attached to unit 8, and a mass and density table written
 54. C into -mass attached to unit 9. (These are temporary files.)

55. C
 56. C The plot description file is produced using the system
 57. C subroutine library *PLOTLIB. All of the plotting is done
 58. C in a single program subroutine, so that other available
 59. C plotting routines may be substituted if desired.
 60. C Please note: In order to run *PLOTLIB subroutines (or any
 61. C subroutines which have not been compiled in Fortran77 and
 62. C which contain character strings or variables in their
 63. C argument lists), in conjunction with this program, (or any
 64. C Fortran77 program) the command *PROCESS (list of subroutines)
 65. C MUST appear in ROW 1, COLUMN 1 if the subroutines are
 66. C called directly from MAIN. If they are called from
 67. C another subroutine, the *PROCESS command must be placed
 68. C in COLUMN 1 OF THE LINE IMMEDIATELY FOLLOWING THE END
 69. C STATEMENT OF THE PREVIOUS ROUTINE.

70. C
 71. C This code will also run on the IBM Professional Fortran
 72. C compiler for PC's and PC compatibles. A math co-processor
 73. C and graphics capability is required. (Graphics routines
 74. C must be supplied.) A 6-layer accretion with rotation
 75. C takes about 90 seconds on a standard PC. Storage needed
 76. C is about 370,000 bytes.

77. C
 78. C To compile (on MTS), first empty -mass and -pdf. Then
 79. C \$RUN *FORTRANS SCARDS=OMNIIC SPUNCH=CYLRUN.

80. C
 81. C To run, enter \$RUN CYLRUN *PLOTLIB. You will be prompted
 82. C for the following input parameters:
 83. C median volume droplet DIAMETER in microns
 84. C
 85. C or n to enter full spectrum, followed by weights,
 86. C entered individually
 87. C
 88. C y or n to simulated spectrum option
 89. C cylinder RADIUS in metres
 90. C
 91. C
 92. C free stream air speed in metres per second
 93. C
 94. C liquid water content in kg per cubic metre
 95. C
 96. C time step per layer in seconds
 97. C
 98. C total number of layers
 99. C
 100. C rotation increment in degrees
 101. C
 102. C
 103. C
 104. C
 105. C
 106. C
 107. C INTERNAL:

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..... 1..... 2..... 3..... 4..... 5..... 6..... 7..... 8..... 9.....

108. C SUBROUTINE DRAW - Plots cylinder profile and all
109. C accreted ice layers with appropriate legend.

110. C
111. C DOUBLE PRECISION FUNCTION MONBET - Estimates local
112. C collision efficiency as a function of surface slope for
113. C a monodisperse droplet population at the mvd. The
114. C function has two forms with mvd smaller or
115. C larger than 20 microns.

116. C
117. C
118. C SUBROUTINE PARAM - Estimates the stagnation line
119. C collision efficiency and maximum accretion angle
120. C for bare cylinder at atmospheric pressure of 100 kPa
121. C from a non-linear regression fit to data derived from
122. C droplet trajectory calculations.

123. C
124. C SUBROUTINE ROTAT1 - When the rotation increment
125. C is non-zero, this subroutine calculates the array
126. C of position angles for the current surface, to replace
127. C surface angles in the array ALPHA.

128. C
129. C SUBROUTINE ROTAT2 - shifts the array indices of
130. C the current LAYX and LAYR arrays to face the new
131. C free stream direction, which has been rotated
132. C according to the rotation increment.

133. C
134. C DOUBLE PRECISION FUNCTION SPCBET - Estimates local
135. C collision efficiency as a function of surface slope
136. C which simulates the effect of a typical droplet size
137. C spectrum centred on the mvd. The function has two forms
138. C for user with mvd smaller or larger than 20 microns.

139. C
140. C SUBROUTINE WTBET - Calculates weighted average of
141. C local collision efficiency estimates, from FUNCTION
142. C MONBET, for droplet spectrum weights in 18 10-micron
143. C size bins.

144. C
145. C
146. C EXTERNAL:

147. C
148. C *PLOTLIB - The subroutines PLOTS, METRIC, PLOT, AXIS2,
149. C LINE, SYMBOL and NUMBER (called from SUBROUTINE DRAW) are
150. C found in the public system subroutine library *PLOTLIB.
151. C (See note above regarding external subroutines in Fortran77).

152. C
153. C ***** Variable Dictionary *****
154. C
155. C ALPHA - array of angles between surface normal
156. C and the free stream direction
157. C
158. C ALPHA - Alpha max., the maximum impingement angle
159. C
160. C AM - array of maximum impingement angles for
161. C each size category of the droplet size
162. C spectrum
163. C

164 C	AREA - cross sectional area between adjacent points in the current layer (dimensioned in metres)
165 C	BETA - local collision efficiency
166 C	BZERO - stagnation line value of the local collision efficiency
167 C	BZ - array of Bzero values for each size category of the droplet size spectrum
168 C	CYLRAD - radius of cylinder in metres (INPUT)
169 C	CYLX,CYLY - arrays containing non-dimensional x and y coordinates of the cylinder surface
170 C	D1 - array of 18.10-micron wide droplet diameter size bins
171 C	DIAG - cross diagonal of the quadrilateral area used in mass calculations
172 C	DROPOI - droplet diameter in microns (INPUT)
173 C	HALFPI - $\pi/2$ (Fortran77 constant PARAMETER)
174 C	LAREA - cross sectional area of the current layer in sq. metres
175 C	LAYERS - integer number of layers
176 C	LAYR,X,LAYRY - array containing non-dimensional x and y coordinates of the accretion surfaces for up to 30 accretion layers
177 C	LDENS - mean density of the current layer in kg/cubic m
178 C	LMASS - ρ - mass of the current layer in kg per unit length (INPUT)
179 C	LWC - liquid water content in kg per cubic metre (INPUT)
180 C	L1,L2,L3 - dimensioned sides of the quadrilateral area defined by two adjacent points on the current surface and two on the surface below
181 C	MASS - array of masses per unit length for areas defined by adjacent points, in kg
182 C	P1 (Fortran77 constant PARAMETER)
183 C	R - temporary variable for storing real result during conversion of the rotation increment an integer number of surface points
184 C	
185 C	
186 C	
187 C	
188 C	
189 C	
190 C	
191 C	
192 C	
193 C	
194 C	
195 C	
196 C	
197 C	
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217 C	
218 C	
219 C	

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220.	C	RHO	- array of local densities in kg/cubic metre
221.	C	ROTN	- no. of surface points in a rotation increment
222.	C	RSPEED	- rotation increment per layer in radians (INPUT)
223.	C	SIMSPC	- logical variable, true if simulated spectrum option is chosen
224.	C	SIMOPT	- character variable for reading in response to simulated spectrum option query (INPUT)
225.	C	SLOPE	- array of local surface slopes calculated at each surface point
226.	C	SMDROP	- logical flag for droplet size, which determines function used to calculate local c.e.
227.	C	SPCOPT	- character variable for reading in response to spectrum weighting option query (INPUT)
228.	C	SPEC	- logical variable, true if spectrum weighting option is chosen
229.	C	S1,S2	- quantities used to determine the quadrilaterals for the layer area calculation
230.	C	TAILEN	- filter length for optional smoothing filter applied to c.e. curve
231.	C	TAU	- accretion time step in seconds (INPUT)
232.	C	THICK	- array of ice thicknesses normal to surface (non - dimensional)
233.	C	TMASS	- total mass of the accretion
234.	C	VEL	- free stream air speed in m per sec (INPUT)
235.	C	WGHT	- array of droplet spectrum volume percentages for each size bin (INPUT)
236.	C		
237.	C		
238.	C		
239.	C		
240.	C		
241.	C		
242.	C		
243.	C		
244.	C		
245.	C		
246.	C		
247.	C		
248.	C		
249.	C		
250.	C		
251.	C		
252.	C		
253.	C		
254.	C		
255.	C		
256.	C		
257.	C		
258.	C		
259.	C		
260.	C		
261.	C		
262.	C		
263.	C		
264.	1		
265.			DOUBLE PRECISION ALPHA(1600),CYLX(1202),CYLY(1202), LAYRX(30,600),LAYRY(30,600),
266.			RHO(600),SLOPE(600),THICK(600), DI(18),BZT(18),BZ(18),AM(18)
267.			
268.			C DOUBLE PRECISION ALPH,AREA,BETA,BZERO,CYLRAD,DIAG, DROPD,LAREA,LDENS,LMASS,LWC,L1,L2,
269.			L3,L4,MASS,RSPEED,S1,S2,TAILEN,TAU, TMASS,VFL
270.	2		
271.			
272.			
273.			
274.			
275.	3	REAL	R

```

276      C      INTEGER          LAYERS,ROTN
        ISN    4     277      C
        ISN    5     278      C      CHARACTER*1 SIMOPT,SPCOPT
        ISN    6     279      C      LOGICAL      SIMSPC,SMDROP,SPEC
        ISN    7     280      C      PARAMETER (PI = 3.1415926535897932,HALFPI = PI/2)
        ISN    8     281      C      COMMON CYLX,CYLY,LAYRX,LAYRY
        ISN    9     282      C      Fill array of droplet diameter bins
        ISN   10    283      C      DATA (DI(I),I=1,18)/5.E-6,15.E-6,25.E-6,35.E-6,45.E-6,
        ISN   11    284      C      +      55.E-6,65.E-6,75.E-6,85.E-6,95.E-6,105.E-6,
        ISN   12    285      C      +      115.E-6,125.E-6,135.E-6,145.E-6,155.E-6,
        ISN   13    286      C      +      165.E-6,175.E-6/
        ISN   14    287      C      Prepare mass table and plot description files.
        ISN   15    288      C      OPEN (UNIT = 8, FILE = '-MASS', STATUS = 'OLD')
        ISN   16    289      C      OPEN (UNIT = 9, FILE = '-PDF', STATUS = 'OLD')
        ISN   17    290      C      Set up headings for mass and density table in -MASS.
        ISN   18    291      C      WRITE (8,110)
        ISN   19    292      C      110 FORMAT (1//,15X,'Mass of Ice Accreted',//,5X,'Layer no.',/
        ISN   20    293      C      +      7X,'Mean Density',//,
        ISN   21    294      C      +      20X,'(kg/m)',9X,'(kg/cu.m)',/
        ISN   22    295      C      +      '*****'
        ISN   23    296      C      +      '*****')
        ISN   24    297      C
        ISN   25    298      C
        ISN   26    299      C
        ISN   27    300      C
        ISN   28    301      C      Set write program title and intro to screen.
        ISN   29    302      C      WRITE (8,110)
        ISN   30    303      C      110 FORMAT (1//,15X,'CYLINDER RIME ICING MODEL',/,
        ISN   31    304      C      +      '*****',/,'*****',//,20X,'Version Feb. 17, 1986',//,
        ISN   32    305      C      +      '22X,'K. J. Finstad',/,'19X',
        ISN   33    306      C      +      'Meteorology Division',/,'15X,'The University of Alberta',
        ISN   34    307      C      +      '15X,'Please enter parameters to begin accretion: ',/,'')
        ISN   35    308      C      Prompt user for input parameters.
        ISN   36    309      C      WRITE (6,120)
        ISN   37    310      C      120 FORMAT (1//,15X,'Drop size flag',/,
        ISN   38    311      C      +      '*****',/,'*****',//,20X,'',
        ISN   39    312      C      +      '22X,'K. J. Finstad',/,'19X',
        ISN   40    313      C      +      '*****',/,'*****',/,'*****',/,'*****',
        ISN   41    314      C      +      '*****',/,'*****',/,'*****',/,'*****',
        ISN   42    315      C
        ISN   43    316      C      Set drop size flag
        ISN   44    317      C      WRITE (6,*)
        ISN   45    318      C      'Median Volume Droplet Diameter in microns? '
        ISN   46    319      C      READ (5,*)
        ISN   47    320      C
        ISN   48    321      C      Set drop size flag
        ISN   49    322      C
        ISN   50    323      C      SMDROP = .FALSE.
        ISN   51    324      C      IF (DROPDI .LT. 20) SMDROP = .TRUE.
        ISN   52    325      C      DROPDI = DROPD1 * 1.OE-6
        ISN   53    326      C      Set flags for requested treatment of droplet spectrum.
        ISN   54    327      C      If full spectrum weighting is requested, input volume
        ISN   55    328      C      weights for each size bin.
        ISN   56    329      C
        ISN   57    330      C
        ISN   58    331      C      WRITE (6,*)

```

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```
      1.....2.....3.....4.....5.....6.....7.....8.....9.....  
1SN 23   332.  WRITE (6,*) 'Do you wish to enter full droplet spectrum?'  
1SN 24   333.  READ (5,9) SPCOPT  
1SN 25   334.  FORMAT (1A1)  
1SN 26   335.  SPEC = FALSE  
1SN 27   336.  IF ((SPCOPT .EQ. 'Y') .OR. (SPCOPT .EQ. 'y')) THEN  
1SN 28   337.    SPEC = TRUE  
1SN 29   338.    WRITE (6,*)  
1SN 30   339.    WRITE (6,*) 'Enter weight for the following size ranges:'  
1SN 31   340.    DO 13 I=1,18  
1SN 32   341.    WRITE (6,*) DI(I)  
1SN 33   342.    WRITE (6,*) DI(I)/1.0E-6, 'microns'  
1SN 34   343.    READ (5,*) WGHT(I)  
1SN 35   344.    CONTINUE  
1SN 36   345.  ELSE  
1SN 37   346.    WRITE (6,*)  
1SN 38   347.    'Do you wish simulated spectrum weighting?'  
1SN 39   348.    READ (5,8) SIMOPT  
1SN 40   349.    FORMAT (1A1)  
1SN 41   350.  ENDIF  
1SN 42   351.  C For simulated spectrum, calculate 'tail' length from  
1SN 43   352.  C median volume diameter.  
1SN 44   353.  IF ((SIMOPT .EQ. 'Y') .OR. (SIMOPT .EQ. 'y')) THEN  
1SN 45   354.    SIMSPC = TRUE  
1SN 46   355.    IF (SMDROP) THEN  
1SN 47   356.      TAILEN = 0.52 * DDEXP(-DROPDI / 7.0E-5)  
1SN 48   357.    ELSE  
1SN 49   358.      TAILEN = 0.37 * DDEXP(-DROPDI / 7.0E-5)  
1SN 50   359.    ENDIF  
1SN 51   360.    SIMSPC = FALSE  
1SN 52   361.    TAILEN = 0.0  
1SN 53   362.  ENDIF  
1SN 54   363.  C  
1SN 55   364.  C  
1SN 56   365.  C  
1SN 57   366.  C  
1SN 58   367.  C  
1SN 59   368.  C  
1SN 60   369.  C  
1SN 61   370.  C  
1SN 62   371.  C  
1SN 63   372.  C  
1SN 64   373.  C  
1SN 65   374.  C  
1SN 66   375.  C  
1SN 67   376.  C  
1SN 68   377.  C  
1SN 69   378.  C  
1SN 70   379.  C  
1SN 71   380.  C  
1SN 72   381.  C  
1SN 73   382.  C  
1SN 74   383.  C  
1SN 75   384.  C  
1SN 76   385.  C  
1SN 77   386.  C  
1SN 78   387.  C  
      1.....2.....3.....4.....5.....6.....7.....8.....9.....
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```
388. ISN      68      C LAYERS = LAYERS * Q1
      ISN      69      C WRITE (6,*) 'Rotation speed in degrees per layer? '
      ISN      70      C WRITE (6,*) 'Rotation speed in degrees per layer? '
      ISN      71      C READ (5,*) RSPEED
      ISN      72      C Convert rotation increment to radians, and calculate
      ISN      73      C approx. Number of surface points in rotation increment
      ISN      74      C
      ISN      75      RSPEED := (RSPEED * 2.0 * PI) / 360.0
      ISN      76      R := (RSPEED / 5.235987E-3)
      ISN      77      RDTM = NINT(R)
      ISN      78      C Call Subroutine Param to calculate Beta zero and alpha max
      ISN      79      C for the svd. If spectrum weighting option has been chosen,
      ISN      80      C call Param for each size bin.
      ISN      81      C
      ISN      82      C CALL PARAM (DROPD1,VEL,CYLRAD,BZERO,ALPHM)
      ISN      83      C IF (SPEC) THEN
      ISN      84      DO 12 I=1,18
      ISN      85      IF (WGHT(I) .NE. 0.0) THEN
      ISN      86      CALL PARAM(DI(I),VEL,CYLRAD,BZ(I),AM(I))
      ISN      87      ELSE
      ISN      88      BZ(I) = 0.0
      ISN      89      AM(I) = 0.0
      ISN      90      ENDIF
      ISN      91      C Make sure accretion limit is revised to reflect the
      ISN      92      C largest droplets present.
      ISN      93      C
      ISN      94      C IF (AM(I) .GT. AM(I-1)) ALPHM = AM(I)
      ISN      95      C
      ISN      96      C 12 CONTINUE
      ISN      97      C
      ISN      98      C
      ISN      99      C Fill array of points defining the initial surface of the
      ISN      100     C cylinder in non-dimensional coordinates. The complete
      ISN      101     C surface contains 1200 points spaced equally in x.
      ISN      102     C
      ISN      103     C DO 100 L=601,1200
      ISN      104     CYLX(I) = (DBLE(I-601) * 3.33333E-3)
      ISN      105     CYLY(I) = DSQRT(1.0 - ((1.0 - CYLX(I))*(1.0 - CYLX(I))))
      ISN      106     CONTINUE
      ISN      107     DO 150 I = 1,600
      ISN      108     CYLX(I) = 2.0 - (DBLE(I) * 3.33333E-3)
      ISN      109     CYLY(I) = -DSQRT(1.0 - ((1.0 - CYLX(I)) * (1.0 - CYLX(I))))
      ISN      110     CONTINUE
      ISN      111     C
      ISN      112     C Copy front half of surface to array which will contain
      ISN      113     C all of the layers.
      ISN      114     DO 200 I = 1,600
      ISN      115     LAYRX(1,I) = CYLX(299 + I)
      ISN      116     LAYRY(1,I) = CYLY(299 + I)
      ISN      117     ISN      96
      ISN      97
      ISN      98
```

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```

      1.....2.....3.....4.....5.....6.....7.....8.....9
      200 CONTINUE
      C Initialize total mass and begin loop to accrete ice layers.
      C
      TMASS = 0.0
      DO 300 J = 2,LAYERS
      C
      C Calculate slope and surface angle from finite differences for
      C each point on current surface. If rotation increment is
      C non-zero, call subroutine to calculate slopes and
      C position angles instead.
      C
      IF (ROTN .GT. 0) THEN
        CALL ROTAT1(J,RSPEED,SLOPE,ALPHA)
      ELSE
        DO 400 I = 2,599
          SLOPE(I) = (LAYRX(J-1,I+1) - LAYRX(J-1,I-1)) /
          DABS(LAYRY(J-1,I+1) - LAYRY(J-1,I-1))
          IF (LAYRY(J-1,I+1) .LT. LAYRY(J-1,I-1)) THEN
            ALPHA(I) = DABS(P1 - DATAN(SLOPE(I)))
          ELSE
            ALPHA(I) = DABS(DATAN(SLOPE(I)))
          ENDIF
        400 CONTINUE
      ENDIF
      C
      C Calculate for each point local density, collision efficiency and
      C ice thickness. The method of calculation for Beta depends on
      C the values of SMDROP, SPEC and SIMSPC.
      C
      DO 500 I = 2,599
        RHO(I) = 890. *
        (1.0 - 143 * (ALPHA(I)/(ALPHM+TAILEN)) -
        246 * ((ALPHA(I)/(ALPHM+TAILEN)) ** 2) -
        309 * ((ALPHA(I)/(ALPHM+TAILEN)) ** 3))
      C
      C Make sure density does not become too small
      C
      IF (RHO(I) .LT. 175.) RHO(I) = 175.
      C
      C Calculate local collision efficiency
      C
      IF (SPEC) THEN
        CALL WTBET(ALPHA(I),SMDROP,BZ,AM,WGHT,BETA)
      ELSE IF (SIMSPC) THEN
        BETA = SPCBET(ALPHA(I),SMDROP,BZERO,ALPHM,TAILEN)
      ELSE
        BETA = MBET(ALPHA(I),SMDROP,BZERO,ALPHM)
      ENDIF
      C
      C Calculate local ice thickness perpendicular to surface
      C (in non-dimensional units)
      C
      THICK(I) = BETA * VEL * 1AU * LWC / (RHO(I) * CYLRAD)
      C
      Define new layer surface coordinates
      C

```

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```

      C
      500.          C
      501.          IF (LAYRX(J-1,I) .LT. LAYRY(J-1,I-1)) THEN
      502.          LAYRX(J,I) = LAYRX(J-1,I) + (THICK(I) *
      503.          DCOS(DATAN(SLOPE(I))))
      504.          ELSE
      505.          LAYRX(J,I) = LAYRX(J-1,I) - (THICK(I) *
      506.          DCOS(DATAN(SLOPE(I))))
      507.          ENDIF
      508.          LAYRY(J,I) = LAYRY(J-1,I) + (THICK(I) *
      509.          DSIN(DATAN(SLOPE(I))))
      510.          CONTINUE
      511.          LAYRX(J,1) = LAYRX(J-1,1)
      512.          LAYRX(J,600) = LAYRX(J-1,600)
      513.          LAYRY(J,1) = LAYRY(J-1,1)
      514.          LAYRY(J,600) = LAYRY(J-1,600)
      515.          C calculate mass and mean density of the layer from
      516.          C the sum of areas between surface points on successive
      517.          C layers. First, initialize layer mass and area.
      518.          C
      519.          C
      520.          LMASS = 0.0
      521.          LAREA = 0.0
      522.          DO 560 I = 2,598
      523.          IF ((THICK(I) .GT. 0.0) .AND. (THICK(I+1) .GT. 0.0)) THEN
      524.          L4 = THICK(I) * CYLRAD
      525.          L2 = THICK(I+1) * CYLRAD
      526.          L1 = DSQRT((LAYRX(J-1,I+1) - LAYRY(J-1,I))**2 + (LAYRY
      527.          (J-1,I+1) - LAYR(J-1,I))**2) * CYLRAD
      528.          L3 = DSQRT((LAYRX(J,I+1) - LAYRY(J,I))**2 + (LAYRY
      529.          (J,I+1) - LAYR(J,I))**2) * CYLRAD
      530.          DIAG = DSQRT((LAYRX(J,I+1)-LAYRY(J-1,I))**2 + (LAYRY
      531.          (J,I+1) - LAYR(J-1,I))**2) * CYLRAD
      532.          S1 = 0.5 * (DIAG+L1+L2)
      533.          S2 = 0.5 * (DIAG+L3+L4)
      534.          AREA = DSQRT(S1 * (S1-DIAG) * (S1-L1) * (S1-L2)) +
      535.          DSQRT(S2 * (S2-DIAG) * (S2-L3) * (S2-L4))
      536.          MASS = ((RHO(I) + RHO(I+1)) / 2.0) * AREA
      537.          LMASS = LMASS + MASS
      538.          LAREA = LAREA + AREA
      539.          ENDIF
      540.          CONTINUE
      541.          C calculate mean density.
      542.          C
      543.          C
      544.          LDENS = LMASS/LAREA
      545.          C write results to mass table.
      546.          C
      547.          WRITE (8,570) J-1,LMASS,LDENS
      548.          570 FORMAT (//,.8X,12.7X,9.5,7X,F9.5)
      549.          C
      550.          C add mass of current layer to total mass.
      551.          C
      552.          C
      553.          TMASS = TMASS + LMASS
      554.          C
      555.          C call subroutine to rotate array indices if required.

```

```

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24. 1986 TIME: 13:15:27 NAME: MAIN PAGE: 11
      . . . . . 6 . . . . . 9
      . . . . . 5 . . . . . 8
      . . . . . 4 . . . . . 7
      . . . . . 3 . . . . . 6
      . . . . . 2 . . . . . 5
      . . . . . 1 . . . . . 4
      . . . . . 0 . . . . . 3
      . . . . . 1 . . . . . 2
      . . . . . 0 . . . . . 1
      . . . . . 9 . . . . . 8
      . . . . . 7 . . . . . 6
      . . . . . 6 . . . . . 5
      . . . . . 5 . . . . . 4
      . . . . . 4 . . . . . 3
      . . . . . 3 . . . . . 2
      . . . . . 2 . . . . . 1
      . . . . . 1 . . . . . 0

ISN 158      557      C      IF (ROTN .GT. 0) CALL ROTAT2(J.ROTN)
ISN 160      558      C
ISN 161      559      C      300 CONTINUE
ISN 162      560      C
ISN 163      561      C      write total mass accreted to table
ISN 164      562      C
ISN 165      563      C      WRITE (8,590) TMASS
ISN 166      564      C      590 FORMAT (8X,//, 'Total mass accreted per meter', F9.5)
ISN 167      565      C      call subroutine to plot cylinder and accreted layers
ISN 168      566      C
ISN 169      567      C      CALL DRAW(ALPHM,BZERO,CYLRAD,DROPDI,LAYERS,LWC,RSPEED,
ISN 170      568      C              SIMSPC,SPEC,TALEN,TAU,VEL)
ISN 171      569      C
ISN 172      570      C      999 STOP
ISN 173      571      C      END

*STATISTICS* SOURCE STATEMENTS = 161. PROGRAM SIZE = 27174 BYTES. PROGRAM NAME = MAIN PAGE: 1
*STATISTICS* NO DIAGNOSTICS GENERATED.
***** END OF COMPIRATION 1 *****

```

LEVEL 1.1.1 (DEC 81)

VS FORTRAN

DATE: JUL 24, 1986 TIME: 13:15:29 NAME: MAIN PAGE: 12

REQUESTED OPTIONS (PROCESS): SC(Axis2.Symbol)

OPTIONS IN EFFECT: NOLIST NOMAP NOXREF NOGOSTMT NODECK OPTIMIZE(0) LANGLVL(77) NOFIPS FLAG(1) NAME(MAIN) LINECOUNT(60)

NOTEST SEQ

```

C
572
573
574      SUBROUTINE DRAW - to plot cylinder and all accretion
575      layers. With appropriate information
576      in the legend. This subroutine uses
577      the plotting library SPOTLIB
578
579      Input Variables : as defined for Main program
580      ALPHA,BZERO,CYLRAD,DROPDI,LAYERS,LWC,
581      RSPEED,SIMSPC,SPEC,TAILEN,TAU,VEL
582
583      Common Variables : CYLX,CYLY,LAYRX,LAYRY
584
585      Other Variables : PLOTX,PLOTY - arrays containing x,y
586      plotting data and scaling information
587      for accretion layer currently being
588      plotted
589
590
591
592      C      SUBROUTINE DRAW(ALPHA,BZERO,CYLRAD,DROPDI,LAYERS,
593      LWC,RSPEED,SIMSPC,SPEC,TAILEN,TAU,VEL)
594
595      C      DOUBLE PRECISION CYLX(1202),CYLY(1202),LAYRX(30,600),
596      LAYRY(30,600).PLOTX(602).PLOTY(602)
597
598      C      DOUBLE PRECISION ALPHM,BZERO,CYLRAD,DROPDI,LWC,RSPEED,
599      TAILEN,TAU,VEL
600
601      C      INTEGER      LAYERS
602      C      LOGICAL      SIMSPC,SPEC
603
604
605      C      COMMON CYLX,CYLY,LAYRX,LAYRY
606
607      C      To plot cylinder surface, form lower surface array.
608      C      and place scaling data at the end of coordinate arrays.
609
610      C
611      7      CYLX(1201) = -6.0
612      8      CYLX(1202) = 1.0
613      9      CYL(1201) = -3.0
614      10     CYL(1202) = 1.0
615      11     PLOTX(601) = -6.0
616      12     PLOTX(602) = 1.0
617      13     PLOTY(601) = -3.0
618      14     PLOTY(602) = 1.0
619      15     C      Open plotting library
620      16     CALL PLOTS
621      17     CALL METRIC(0)
622
623

```

```

624      C Draw 6 by 9 inch box
625      C
626      CALL PLOT(2.0,2.0,-3)
627      CALL PLOT(0.0,0.6,0.3)
628      CALL PLOT(0.0,0.6,0.3)
629      CALL PLOT(0.0,0.6,0.2)
630      CALL PLOT(0.0,0.0,0.2)
631      C Label axes and draw tics
632      C
633      CALL AXIS2(0.0,0.0,'X/CYLRAD',-8.9,0.0,0,-6.0,1.0,1.0)
634      CALL AXIS2(0.0,0.0,'Y/CYLRAD',-8.6,0.9,0,-3.0,1.0,0.5)
635      C
636      C write legends to top left corner
637      C
638      C
639      CALL SYMBOL(1,1.5,4.0,10,'DROPLET MVD'   - .0,0,15)
640      CALL NUMBER(2,6.5,4.0,10,'DROPDI',0.0,0.07)
641      CALL SYMBOL(3,5.5,4.0,10,'N' ,0,0.01)
642      CALL SYMBOL(1,1.5,2.0,10,'CYL RADIUS' - .0,0,15)
643      CALL NUMBER(2,6.5,2.0,10,'CYLRAD',0.0,0.04)
644      CALL SYMBOL(3,5.5,2.0,10,'N' ,0,0.01)
645      CALL SYMBOL(1,1.5,0.0,10,'AIR SPEED' - .0,0,15)
646      CALL NUMBER(2,6.5,0.0,10,'VEL',0.0,0.02)
647      CALL SYMBOL(3,5.5,0.0,10,'N' / SEC',0.0,0.07)
648      CALL SYMBOL(1,1.4,8.0,10,'LWC'    - .0,0,15)
649      CALL NUMBER(2,6.4,8.0,10,'LWC',0,0.05)
650      CALL SYMBOL(3,5.4,8.0,10,'KG / M^0.3' ,0,0,0.09)
651      CALL SYMBOL(1,1.4,6.0,10,'TIME STEP' - .0,0,15)
652      CALL NUMBER(2,6.4,6.0,10,'TAU',0,0.02)
653      CALL SYMBOL(3,5.4,6.0,10,'SEC',0,0.03)
654      CALL SYMBOL(1,1.4,4.0,10,'BETA ZERO' - .0,0,15)
655      CALL NUMBER(2,6.4,4.0,10,'BZERO',0,0.03)
656      CALL SYMBOL(1,1.4,2.0,10,'ALPHA MAX' - .0,0,15)
657      CALL NUMBER(2,6.4,2.0,10,'(ALPHM+TAILEN)',360./6,283.0,0.01)
658      CALL SYMBOL(3,5.4,2.0,10,'DEGREES',0,0.07)
659      CALL SYMBOL(1,1.4,0.0,10,'JOIN INCR' - .0,0,15)
660      CALL NUMBER(2,6.4,0.0,10,'RSPEED',360./6,283.0,0.01)
661      CALL SYMBOL(3,5.4,0.0,10,'DEGREES',0,0.07)
662      CALL SYMBOL(1,1.3,8.0,10,'SPECTRUM',0,0.09)
663      IF (SPEC) THEN
664          CALL SYMBOL(2,6.3,8.0,10,'YES',0,0.03)
665          ELSE IF (SIMSPC) THEN
666              CALL SYMBOL(2,6.3,8.0,10,'SIMULATED',0,0.09)
667          ELSE
668              CALL SYMBOL(2,6.3,8.0,10,'NO',0,0.02)
669          ENDIF
670      C Draw initial cylinder surface
671      C
672      CALL LINE(CYLX,CYLY,4200,2.0)
673      C
674      C Transfer each layer in turn to the array holding
675      C the scaling parameters
676      C
677      DO 610 J = 2,LAYERS
678          DO 700 L = 1,600
679      C

```

LEVEL 1.1.1 (DEC 81). VS FORTRAN DATE: JUL 24. 1986 TIME: 13:15:29 NAME: DRAW PAGE: 14

ISN 57 680. PLOTX(L) = LAYRX(J,L)
ISN 58 681. PLOTY(L) = LAYRY(J,L)
ISN 59 682. 700 CONTINUE
ISN 683. C
ISN 684. C Draw accretion layer
ISN 685. C
ISN 60 686. CALL LINE(PLOTX,PLOTY,600,2,0)
ISN 61 687. 610 CONTINUE
ISN 688. C
ISN 689. C Close plotting library
ISN 690. C
ISN 62 691. CALL PLOT(1.0,1.0,999)
ISN 63 692. RETURN
ISN 64 693. END

SOURCE STATEMENTS = 64. PROGRAM SIZE = 12132 BYTES. PROGRAM NAME = DRAW PAGE: 12.

STATISTICS NO DIAGNOSTICS GENERATED.
***** OF COMPILED 2 *****

LEVEL 1.1.1 (DEC 81)

VS FORTRAN

DATE: JUL 24 1986 TIME: 12:15:00 NAME: MAIN PAGE: 15

OPTIONS IN EFFECT:

NOLIST NOMREF NOGOSTMT MODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(0) LANG VL(77) NOFIPS FLAG(I) NAME(MAIN) LINECOUNT(50)

```
694 C
695 C FUNCTION MBET - Calculate local estimate for droplet populations
696 C at and either above or below 20 microns
697 C using a cosine curve scaled to the deflection
698 C line c. g. and the maximum accretion angle
699 C
700 C
701 C Input Variables: as defined for main program:
702 C   ALPHI, BZERO, ALPHA, SMDROP
703 C
704 C
705 C
1 706 C DOUBLE PRECISION FUNCTION MBET(ALPHI, SMDROP, BZERO, ALPHI)
2 707 C
3 708 C DOUBLE PRECISION BZERO, ALPHI, ALPHA
4 709 C
5 710 C LOGICAL
6 711 C
7 712 C PARAMETER (HALFP1 = 3.1415926535897932/2.)
8 713 C
9 714 C IF (ALPHI .GE. ALPHI) THEN
10 715 C   MBET = 0.0
11 716 C ELSE IF (SMDROP) THEN
12 717 C   MBET = BZERO * (BZERO * (DSQRT(DCOS(ALPHI)) - 1.0)) /
13 718 C           (1.0 - DSQRT(DCOS(ALPHI)))
14 719 C ELSE
15 720 C   MBET = BZERO * (DCOS((ALPHA/ALPHI) * HALFP1))
16 721 C ENDIF
17 722 C RETURN
18 723 C
19 C
20 C
21 C
22 C
23 C
```

STATISTICS

SOURCE STATEMENTS = 13. PROGRAM SIZE = 586 BYTES. PROGRAM NAME = MBET PAGE: 15

STATISTICS NO DIAGNOSTICS GENERATED

***** END OF COMPILED 3 *****

LEVEL 1.1.1 (DEC 91)

VS FORTRAN

DATE: JUL 24, 1986 TIME: 13:15:30 NAME: MAIN PAGE: 16

OPTIONS IN EFFECT: NOLIST NOMAP NODECK SOURCE TERM PROJECT FIXED
OPTIMIZE(O) LANGVL(77) NOFTPS FLAG1) NAME(MAIN) LINECOUNT(60)

```

724
725      C   * Subroutine PARAN - Estimates Beta zero and alpha max
726      C   * free the input parameters according to
727      C   * an empirical fit to direct trajectories
728      C   * calculated with program VISION.
729      C
730      C   Input variables : as defined for Main program:
731      C   CYLRAD,DROPDI,VEL
732      C
733      C   Output variables : as defined for Main program:
734      C   ALPHM,BZERO
735      C
736      C   Other variables :
737      C   RE   - free stream Reynolds number
738      C   K    - Langmuir inertia parameter
739      C   PHI   - Langmuir parameter
740      C   TERM1, - terms used in the fitting function
741      C   TERM2,
742      C   TERM3,
743      C   TERM4
744      C   A(2,12) - array containing constants for
745      C   the fitting functions
746      C
747      C
748      C
749      C
750      C   SUBROUTINE PARAM (DROPDI,VEL,CYLRAD,BZERO,ALPHM)
751      C
752      C   DOUBLE PRECISION ALPHM,BZERO,CYLRAD,DROPDI,K,
753      C   PHI,RE,TERM1,TERM2,TERM3,VEL
754      C
755      C   DOUBLE PRECISION A(2,12)
756      C   DATA (A(1,1), I=1,12)/1.215,-7.420,-3,-0.561,-0.644,-0.164,
757      C   3.050,-3,0.430,2.22,-0.45,-0.767,-0.806,-0.068/
758      C   DATA (A(2,1), I=1,12)/2.327,-4.090,-3,-0.388,-0.593,-0.689,
759      C   8.50,-3,0.383,1.757,-0.298,-0.42,-0.96,-0.179/
760      C
761      C   Calculate free stream Reynolds number, phi and k assuming
762      C   water density is 1000, air density 1.323 and air's
763      C   viscosity 1.669 in MKS units. These correspond to
764      C   an air temperature of -10 C and pressure 100 kPa.
765      C
766      C   RE = DROPDI * VEL / 1.262E-5
767      C   K = (VEL * DROPDI * CYLRAD / 3.329E6) / CYLRAD
768      C   PHI = (RE * RE) / K
769      C
770      C   Calculate Beta zero and alpha max
771      C
772      C
773      C   DO 10 I = 1,2
774      C   TERM1 = (A(1,7)* (K ** A(1,2))*DEXP(A(I,3)*K**A(I,4)))
775      C   + A(1,6)*(PHI ** A(I,7))
776      C

```

LEVEL 1.1.1 (DEC 81)

DATE: JUL 24. 1986

VS FORTRAN

NAME: PARAM PAGE: 17

```
12   777.      TERM3 = (A(1,8) * (K ** A(1,9)) *DEXP(A(1,10)*K*A(I,11)))  
13   778.      + A(1,12)  
14   779.      IF (I .EQ. 1) THEN  
15   780.          BZERO = TERM1 - (TERM2 + TERM3)  
16   781.      ELSE  
17   782.          ALPHM = TERM1 - (TERM2 + TERM3)  
18   783.      ENDIF  
19   784.      10  CONTINUE  
20   785.      C Check for extreme values of K  
21   786.      C  
22   787.      C  
19   788.      IF (K .LT. 0.17) THEN  
19   789.          ALPHM = 0.0  
19   790.          BZERO = 0.0  
20   791.      ELSE IF (K .GT. 1000.) THEN  
21   792.          ALPHM = 1.56  
21   793.          BZERO = 0.99  
22   794.      ENDIF  
23   795.      RETURN  
24   796.  
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STATISTICS SOURCE STATEMENTS = 27. PROGRAM SIZE = 1362 BYTES. PROGRAM NAME = PARAM PAGE: 16.

STATISTICS NO DIAGNOSTICS GENERATED.

***** END OF COMPILED 4 *****

LEVEL 1.1.1 (DEC 81) DATE: JUL 24, 1986 TIME: 13:15:31 NAME: MAIN PAGE: 18
 OPTIONS IN EFFECT: NOLIST NOMAP NOXREF NOGOSTMT NODECK TERM OBJECT FIXED
 OPTIMIZE(0) LANGVL(77) NOFIPS FLAG(I) NAME(MAIN) LINECOUNT(60)

```

 1. . . . . 2. . . . . 3. . . . . 4. . . . . 5. . . . . 6. . . . . 7. . . . . 8. . . . . 9.
    C
    797. C
    798. C
    799. C     SUBROUTINE ROTAT1 - calculates position angles and
    800. C     slopes for each point on the
    801. C     current surface. Position angles
    802. C     are with respect to the current
    803. C     free stream direction.
    804. C
    805. C     Input Variables : as defined for Main program;
    806. C     J.RSPEED
    807. C
    808. C     Output Variables : ALPHA - array of position angles
    809. C     measures w.r.t. center of
    810. C     cylinder and free stream
    811. C     as defined for Main program;
    812. C     SLOPE
    813. C
    814. C     Common Variables : CYLX,CYLY,LAYRX,LAYRY
    815. C
    816. C
    817. C
    818. C
    819. C     SUBROUTINE ROTAT1(J,RSPEED,SLOPE,ALPHA)
    ISN 1     820. C
    ISN 2     821. C     DOUBLE PRECISION ALPHA(600),CYLX(1202),CYLY(1202),
    ISN 3     822. C     + LAYRX(30,600),LAYRY(30,600),SLOPE(600)
    ISN 4     823. C
    ISN 5     824. C     DOUBLE PRECISION RSPEED
    ISN 6     825. C     INTEGER     J
    ISN 7     826. C     PARAMETER     PI = 3.1415926535897932,HALFPI=PI/2.)
    ISN 8     827. C     COMMON CYLX,CYLY,LAYRX,LAYRY
    ISN 9     828. C     calculate slopes
    ISN 10    829. C
    ISN 11    830. C
    ISN 12    831. C
    ISN 13    832. C
    ISN 14    833. C
    ISN 15    834. C
    ISN 16    835. DO 590 I=2,599
    ISN 17    836. IF (LAYRY(J-1,I+1) EQ LAYRY(J-1,I-1)) THEN
    ISN 18    837.     SLOPE(I) = 1.E10
    ISN 19    838. ELSE
    ISN 20    839.     SLOPE(I) = (LAYRX(J-1,I+1)-LAYRX(J-1,I-1)) / DABS
    ISN 21    840.     +(LAYRY(J-1,I+1)-LAYRY(J-1,I-1))
    ISN 22    841. ENDIF
    ISN 23    842. C     Smooth over any discontinuities in slope caused
    ISN 24    843. C     by accreting over the previous layer's accretion
    ISN 25    844. C     limit.
    ISN 26    845. C
    ISN 27    846. IF (LAYRY(J-1,I-1) GT 0.0) THEN
    ISN 28    847.     IF (LAYRY(J-1,I+1) LT LAYRY(J-1,I-1)) THEN
    ISN 29    848.       IF (SLOPE(I) LT SLOPE(I-1)) SLOPE(I) = SLOPE(I-1)
    ISN 30    849.
  
```

LEVEL .1.1.1 (DEC - 81)

DATE: JUL 24.. 1986

TIME: 13:15:31

NAME: ROTAT1 PAGE: 19

VS FORTRAN

```
ISN 17    850
ISN 18    851      C
          852      C Calculate position angles.
          853      C
          854      IF (LAYRX(J-1,I) EQ. 1.0) THEN
          855          ALPHA(I) = DABS(HALFPI - ((J-2)*RSPEED))
          856          ELSE IF (LAYRX(J-1,I) GT. 1.0) THEN
          857              ALPHA(I) = HALFPI - ((J-2)*RSPEED) +
          858                  DATAN((LAYRX(J-1,I) - 1.) * LAYRY(J-1,I))
          859          ELSE
          860              IF (DABS(LAYRY(J-1,I)) LT. 1.00-10) LAYRY(J-1,I) = 0.0
          861              IF (DABS(LAYRX(J-1,I)) LT. 1.00-10) LAYRX(J-1,I) = 0.0
          862              ALPHA(I) = DABS(DATAN(LAYRY(J-1,I) /
          863                  (1.0 - LAYRX(J-1,I))) - ((J-2)*RSPEED))
          864          ENDIF
          865          590 CONTINUE
          866          RETURN
          867      C
          868      END
          869      END
```

STATISTICS

SOURCE STATEMENTS = 29. PROGRAM SIZE = 1998 BYTES. PROGRAM NAME = ROTAT1

PAGE: 18.

STATISTICS

NO DIAGNOSTICS GENERATED.

***** END OF COMPILEATION 5 *****

LEVEL 1.1.1 TOEC 81) VS FORTRAN

DATE: JUL 24, 1986 TIME: 13:15:31

NAME: MAIN PAGE: 20

OPTIONS IN EFFECT: MOLIST NOMAP NOXREF NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(O) LANGVL(77) NOFIPS FLAG(I) NAME(MAIN) LINECOUNT(60)



C *****
C * SUBROUTINE ROTAT2 - shifts array indices of current layer
C * by an amount corresponding to the
C * rotation increment. Points from the
C * cylinder array must be added to one
C * end.
C *
C * Input variables : as defined in main program:
C * J,ROTN
C *
C * Common variables : CYLX,CYLY,LAYRX,LAYRY
C *
C * Other variables : K,M - index limits for shifting
C *
C *
C * SUBROUTINE ROTAT2(J,ROTN)
C *
C * DOUBLE PRECISION CYLX(1202),CYLY(1202),LAYRX(30,600),
C * LAYRY(30,600)
C *
C * INTEGER J,ROTN
C *
C * COMMON CYLX,CYLY,LAYRX,LAYRY
C *
C * Renumber layer array to face new direction of air stream,
C * dropping increment number of points at the front, end.
C *
C * 895 C
C * 896 C
C * 897 C
C * K = 600-ROTN
C * M = 600-ROTN+1
C * DO 580 I = 1,K
C * LAYRX(J,I) = LAYRX(J,I+ROTN)
C * LAYRY(J,I) = LAYRY(J,I+ROTN)
C * 580 CONTINUE
C *
C * Add increment number of points from the bare cylinder to
C * the back, end.
C *
C * 907 C
C * DO 585 I = M,600
C * LAYRX(J,I) = CYLX(299 + (J-1)*ROTN + 1)
C * LAYRY(J,I) = CYLY(299 + (J-1)*ROTN + 1)
C * 585 CONTINUE
C * RETURN
C * END
C *
ISN 11 908
ISN 12 909
ISN 13 910
ISN 14 911
ISN 15 912
ISN 16 913

SOURCE STATEMENTS = 16, PROGRAM SIZE = 780 BYTES, PROGRAM NAME = ROTAT2 PAGE: 20
STATISTICS
STATISTICS
NO DIAGNOSTICS GENERATED.
***** END OF COMPILEATION 6 *****

LEVEL 1.1.1 (DEC 81)

VS FORTRAN

DATE: JUL 24, 1986 TIME: 13:15:31

NAME: MAIN PAGE: 21

OPTIONS IN EFFECT: NOLIST NOMAP NOXREF NOGOSTAT NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(O) LANGVL(77) NOFIPS FLAG(P) NAME(MAIN) LINECOUNT(60)

```
914. C
915. C      Function SPCBET Estimates Beta(alpha) for a
916. C      simulated droplet size distribution. For an
917. C      myd ette smaller or larger than 20 microns.
918. C      The polynomial function is scaled to (alpha
919. C      max + tail length), where the tail length
920. C      is a function of the myd.
921. C
922. C      Input Variables : as defined for Main Program:
923. C      ALPHA,ALPHM,BZERO,SMDROP,TAILEN
924. C
925. C      Other Variables : X - scaled angle = Alpha /
926. C                  (Alpha + Tailen)
927. C
928. C
929. C
930. C
931. C      DOUBLE PRECISION FUNCTION SPCBET(ALPHA,SMDROP,BZERO,
932. C                                         ALPHM,TAILEN)
933. C
934. C      DOUBLE PRECISION ALPHA,ALPHM,BZERO,TAILEN,X
935. C      LOGICAL     SMDROP
936. C
937. C      IF (ALPHA .GE. (ALPHM+TAILEN)) THEN
938. C          SPCBET = 0.0
939. C
940. C      ELSE
941. C          X = ALPHA / (ALPHM+TAILEN)
942. C          IF (SMDROP) THEN
943. C              SPCBET = BZERO * (1.0 + 0.25 * X - 3.084 *
944. C                         (X ** 2) + 1.866 * (X ** 3))
945. C          ELSE
946. C              SPCBET = BZERO * (1.0 + 0.0287 * X - 1.536 *
947. C                         (X ** 2) + 2.484 * (X ** 3) - 4.112 *
948. C                         (X ** 4) + 2.539 * (X ** 5))
949. C
950. C      ENDIF
951. C      IF (SPCBET .LT. 0.0) SPCBET = 0.0
952. C
953. C
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997. C
998. C
999. C
```

SOURCE STATEMENTS = 16. PROGRAM SIZE = 708 BYTES. PROGRAM NAME = SPCBET PAGE:

STATISTICS NO DIAGNOSTICS GENERATED

END OF COMPIRATION 7

LEVEL 1.1.1 (DEC 81) VS FORTRAN

OPTIONS IN EFFECT: NOLIST NODECK NOGOSTMT NOFILE OPTIMIZE(O) LANGLVL(77) NOFIPS FLAG(I) NAME(MAIN) LINECOUNT(60)

DATE: JUL 24, 1986 TIME: 13:15:32 NAME: MAIN PAGE: 22

```
954 C
955 C *****
956 C * SUBROUTINE WTBT * Calculates the average local
957 C * collision efficiency according to
958 C * the input volume weighted droplet
959 C * size spectrum.
960 C
961 C * Input Variables : as defined for Main program:
962 C * ALPHA,AM,BZ,SMDROP,WGHT
963 C
964 C * Output Variables : as defined for Main program:
965 C * BETA
966 C
967 C
968 C
969 C
970 C
971 C
972 C
973 C
974 C
975 C
976 C
977 C
978 C
979 C
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998 C
999 C
```

SUBROUTINE WTBT(ALPHA,SMDROP,BZ,AM,WGHT,BETA)

DOUBLE PRECISION AM(18),BZ(18),WGHT(18)

DOUBLE PRECISION ALPHA,BETA

LOGICAL SMDROP

BETA = 0.0

DO 22 L=1,18

IF (WGHT(L) .NE. 0.0) THEN

BETA = BETA + WGHT(L) * MNBET(

ALPHA(1),BZ(L),AM(L),SMDROP)

ENDIF

CONTINUE

22 RETURN

END

***** END OF COMPILE B *****

STATISTICS

STATISTICS

NO DIAGNOSTICS GENERATED.

SOURCE STATEMENTS = 12, PROGRAM SIZE = 654 BYTES, PROGRAM NAME = WTBT , PAGE: 22.

LEVEL 1.1.1 (DEC- 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:15:32 NAME: MAIN PAGE: 23

SUMMARY OF MESSAGES AND STATISTICS FOR ALL COMPILENTIONS

STATISTICS SOURCE STATEMENTS = 161, PROGRAM SIZE = 2744 BYTES, PROGRAM NAME = MAIN PAGE: 1.
STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPILATION 1 *****

STATISTICS SOURCE STATEMENTS = 64, PROGRAM SIZE = 12132 BYTES, PROGRAM NAME = DRAW PAGE: 12.
STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPILATION 2 *****

STATISTICS SOURCE STATEMENTS = 13, PROGRAM SIZE = 586 BYTES, PROGRAM NAME = MONBET PAGE: 15.
STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPILATION 3 *****

STATISTICS SOURCE STATEMENTS = 27, PROGRAM SIZE = 1362 BYTES, PROGRAM NAME = PARAM PAGE: 16.
STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPILATION 4 *****

STATISTICS SOURCE STATEMENTS = 29, PROGRAM SIZE = 1998 BYTES, PROGRAM NAME = ROTAT1 PAGE: 18.
STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPILATION 5 *****

STATISTICS SOURCE STATEMENTS = 16, PROGRAM SIZE = 780 BYTES, PROGRAM NAME = ROTAT2 PAGE: 20.
STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPILATION 6 *****

STATISTICS SOURCE STATEMENTS = 16, PROGRAM SIZE = 708 BYTES, PROGRAM NAME = SPCBET PAGE: 21.
STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPILATION 7 *****

STATISTICS SOURCE STATEMENTS = 12, PROGRAM SIZE = 654 BYTES, PROGRAM NAME = WTBET PAGE: 22.
STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPILATION 8 *****

***** SUMMARY STATISTICS ***** O DIAGNOSTICS GENERATED HIGHEST SEVERITY CODE IS 0

Appendix III

Approximation of the mvd from Stagnation Line Thickness

The measurement of water droplet sizes and size distributions is one of the major experimental problems in icing studies. The very common measurement technique, known as the oil slide method, has been shown by Makkonen and Stallabrass (1984) to overestimate the median volume diameters of observed droplet spectra in comparison to measurements made with an optical sizing probe, the Forward Scattering Spectrometer Probe, or FSSP, manufactured by Particle Measuring Systems and described by Knollenberg (1981). Makkonen and Stallabrass provide an empirical correction formula for oil slide measured mvds in the range 10 to 40 μm , as follows:

$$\text{mvd}_{\text{FSSP}} = 8.8 + 0.27 \text{ mvd}_{\text{oil slide}} \quad \text{A.1}$$

All of the wind tunnel experiments performed in the PROST tunnel and described in this thesis have had droplet size distributions determined by the oil slide method. The mvds derived from this data appear to be too large, not only in light of Makkonen and Stallabrass' results, but also because the theoretical collection efficiencies calculated

for these mvd's imply a minimum stagnation line thickness much larger than that observed.

In order to supply the most accurate input data possible to the model, these oil slide measurement should be corrected. However, some of the circular cylinder cases lie outside of the range given for Equation A.1. For these cases an approximate mvd has been derived from the observed stagnation line thicknesses of accreted ice, using the method outlined below, similar to the rotating cylinder methods which have been used for the same purpose.

For cases which lie inside the range of validity, these approximate mvd's agree well (within about 10 per cent) with the results of Equation A.1. The correlation of Makkonen and Stallabrass has therefore been used for all cases within the appropriate range..

Outside of that range, however, it is important to remember that the approximation given here is derived independently of the actual size spectrum, and represents a droplet size corresponding to the average value of β_0 for the time period of the accretion layer. It has been used here only because of the absence of a reliable calibration for oil slide measurements of droplets larger than 40 μm .

The method is as follows:

1. The stagnation line thickness (Th) of the chosen layer is measured from a photograph of the accretion in cross-section.
2. A first estimate of β_0 is calculated from applying

the dry growth equation to the stagnation line:

$$\beta_0 = (\rho_0 \text{ Th}) / (\text{U lwc } \tau) \quad \text{A.2}$$

A best estimate of ρ_0 may be made as in Chapter IV. For lwc, the first guess is made using the oil slide mvd, and the measured rate of mass collected on a small rotating cylinder, and calculated by the algorithm given by Stallabrass (1978).

3. Calculate the Stokes parameter

$$\phi = 9 \rho_a^2 D_c U / (\mu \rho_d); \quad \text{A.3}$$

then read from Figure II.3, a value of K corresponding to ϕ and β_0 .

4. Derive a new value of mvd from the definition of K:

$$K = (\rho_d / 18 \mu) \left(\text{mvd}^2 U / D_c \right) \quad \text{A.4}$$

5. Recalculate lwc using the new mvd.

6. Recalculate β_0 from new lwc. Repeat steps 2 through 6 until the values of β_0 , mvd and lwc appear to have converged.

For the examples given below, the rotating cylinder diameter was 2.46 mm.

Example 1:oil slide mvd = 24.2 μm

Th = 1.49 mm

U = 10.7 m/sec

 τ = 600 sec ρ_0 = 750 kg/m³

rotating cylinder mass = 0.029 g/min

revised mvd = 13.8 μm mvd from Equation A.1 = 15.3 μm **Example 2:**oil slide mvd = 136 μm

Th = 13.5 mm

U = 10.24 m/sec

 τ = 1800 sec ρ_0 = 890 kg/m³

rotating cylinder mass = 0.122 g/min

revised mvd = 62 μm

References

- Knollenberg, R.G. 1981: Techniques for Probing Cloud Microstructure. Clouds, Their Formation, Optical Properties, and Effects, ed.s P.V. Hobbs and A. Deepak. Academic Press, 15-92.
- Makkonen, L. and Stallabrass, J.R. 1984: Ice Accretion on Cylinders and Wires. National Research Council, Division of Mechanical Engineering, Technical Report TR-LT-005, 50 pp.
- Stallabrass, J.R. 1978: National Research Council, Division of Mechanical Engineering, Technical Report TR-LT-92, 26 pp., 7 fig.s, + 2 App.s.

Appendix IV

A Geometrical Derivation of Equation V.3

Figure A.1 illustrates two droplet trajectories impinging on an airfoil, at zero angle of attack. If the initial separation of the trajectories is Δy , then the distance separating them (in the direction normal to the lower trajectory) will have increased at the point of impact to, say, $(\Delta y + a)$. Referring to Figure A.1, if the angle γ is the departure from horizontal of the tangent to the upper trajectory at the point X, it follows that for small $\Delta \ell$:

$$\cos(\alpha + \gamma) = (\Delta y + a) / \Delta \ell = \beta + (a / \Delta \ell). \quad A.5$$

using the definition of β given in Chapter II.

For airfoils, γ is always $\ll \alpha$, so:

$$\cos(\alpha + \gamma) \approx \cos \alpha \quad A.6$$

Now if it is assumed that $(a / \Delta \ell)$ is approximately a constant, and applying the boundary condition at $\alpha = 0$; this gives:

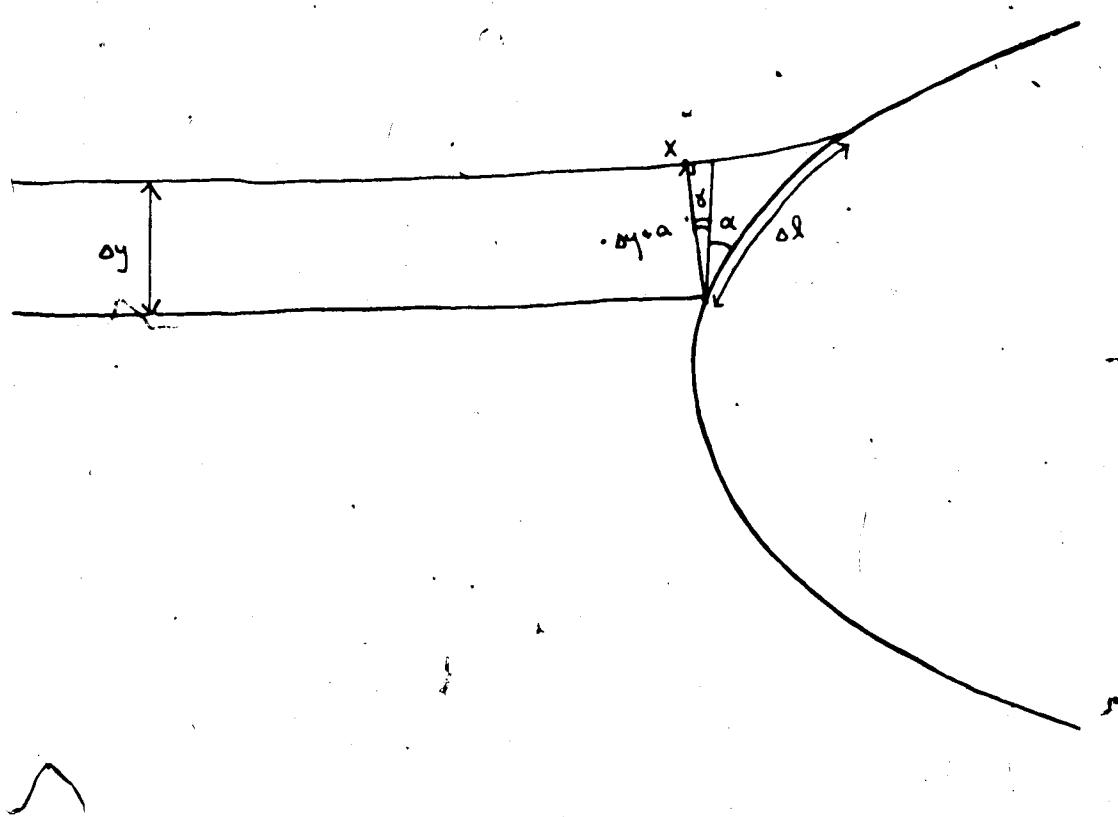


Figure A.1 - Droplet trajectories impinging on an airfoil.

$$\beta \approx \cos \alpha - (1 - \beta_0)^f,$$

A.7

which is Equation V.3. The above approximation should then hold when the impinging trajectories have small curvature.

Appendix V

Program Listing: OMNIFOIL

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OPTIONS IN EFFECT: MOLIST NOMAP NOXREF NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(O) LANG VL(77) NDFIPS FLAG(1) NAME(MAIN) LINECOUNT(EQ)

PAGE:

SEQ

..... 1 2 3 4 5 6 7 8 9

***** Airfoil Rime Icing Model *****

Karen J. Flinstad June, 1986

1. C
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51. C

This program will calculate and plot rime ice accretion layers on an airfoil, as well as calculate the mass and mean density of each layer. Local collision efficiencies for each point on the surface are calculated as a cosine function of the current local surface angle and scaled to the value of stagnation line collision efficiency (Beta zero), which is determined from droplet trajectory calculations.

Beta zero is updated on every fourth layer to account for the changing flow around the accretion. Layer time steps should be chosen so that Beta zero is updated at reasonable intervals, say after a maximum thickness of no more than 1/10 of the chord length has been accreted.

The maximum accretion angle, alpha max, is determined from Beta zero.

When the angle of attack is zero, a "seedulated spectrum" version of local collision efficiency distribution may be used to approximate the effects of droplet size spectra with median volume diameters less than about 20 microns. Larger drops are well represented by the monodisperse version of collision efficiency parameterisation at the MHD.

For other angles of attack, an approximate "skewed" distribution is used, which should apply to all droplet sizes.

Local density is approximated by a polynomial function of the local surface slope.

The output consists of a plot description file written into -pdf attached to unit 8, and a mass and density table written into -mass attached to unit 9. (A negative sign identifies temporary files on MTS.)

The plot description file is produced using the system subroutine library *PLOTLIB. All of the plotting is done in a single program subroutine, so that other available plotting routines may be substituted if desired.

An additional external subroutine, from the *IMSLPLIB library, is used in the solution of the potential flow. Other matrix inversion routines may be substituted.

Please note: In order to run *PLOTLIB subroutines (or any subroutines which have not been compiled in Fortran77 and which contain character strings or variables in their

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Argument list). In conjunction with this program, (or any Fortran77 program) the command EPROCESS (list of subroutines) MUST appear in ROW 1, COLUMN 1 if the subroutines are called directly from MN. If they are called from another subroutine, the EPROCESS command must be placed IN COLUMN 1 OF THE LINE IMMEDIATELY FOLLOWING THE END STATEMENT OF THE PREVIOUS ROUTINE.

To compile (on MTS), first empty -mass and -pdf. Then
\$RUN *FORTRAN5 SCARDS=OMNIFOIL SPUNCH=FOILRUN

To run, enter \$RUN FOILRUN+IMSLDPLIB+PLOTLIB.
You \$111 be prompted for the following input parameters:

median volume droplet DIAMETER in microns

y or n to simulated spectrum option

airfoil chord length in metres

NACA airfoil height as fraction of chord

free stream air speed in metres per second

liquid water content in kg per cubic metre

Up step per layer in seconds

total number of layers

angle of attack of the free stream in degrees

OR, input data may be entered in a file. You must then specify "5 = filename" in the \$RUN command.

Subprograms

..... Subprograms

INTERNAL:

91 SUBROUTINE DRAW - Plots airfoil profile and all
92 accreted ice layers with appropriate legend

93 SUBROUTINE DROPS - Integrates droplet trajectories
94 and calculates stagnation line collision efficiency.
95 Subroutines called only from within DROPS (CALLID,
96 FLOWXY, KCALC, POFLO, TRAJEC) are described in the program
97 summary to DROPS.

98 DOUBLE PRECISION FUNCTION ONEBET - Estimates local
99 collision efficiency as a function of surface slope for
100 a monodispersed droplet population at the end.

101 DOUBLE PRECISION FUNCTION PHIBET - calculates local
102 collision efficiency as a function of slope w.r.t. to a
103 free stream direction, and distance from the maximum

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108. C local value of collision efficiency (set to 0.95 times
 109. C the value for zero angle of attack). The location of
 110. C the maximum is estimated from a correlation derived
 111. C from previous numerical results

112. C

113. C

114. C

115. C

116. C

117. C

118. C

119. C

120. C

121. C

122. C

123. C

124. C

125. C

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162. C

163. C

DOUBLE PRECISION FUNCTION SPCBET - Estimates local

114. C collision efficiency as a function of surface slope
 115. C for an average distribution shape of droplet sizes,
 116. C scaled to the chord. The function assumes that small
 117. C droplets dominate in the spectrum, with chord less than
 118. C about .20 microns.

EXTERNAL:

123. C *IMSLDPLIB - The subroutine LEQT1F is used to solve
 124. C the matrix of potential vortices for the potential
 125. C flow calculation in subroutine DROPS.

126. C *PLOTLIB - The subroutines PLOTS, METRIC, PLOT, AXIS2,
 127. C LINE, SYMBOL and NUMBER (called from SUBROUTINE DRAW) are
 128. C found in the public system subroutine library *PLOTLIB.
 129. C (See note above regarding external subroutines in Fortran77).

130. C ***** Variable Dictionary *****
 131. C
 132. C ALPHA - array of angles between surface normal
 133. C and the free stream direction

134. C ALPHAM - Alpha max, the maximum impingement surface
 135. C angle

136. C APHI - parameter for determining the distribution
 137. C of collision efficiency and of density
 138. C about the position of maximum c.e. when
 139. C phi is non-zero.

140. C AREA - cross sectional area between
 141. C adjacent points in the current layer
 142. C (dimensioned in metres)

143. C AZERO - the surface angle of maximum collision
 144. C efficiency when phi is non-zero,
 145. C determined from an approximating
 expression.

146. C BETA - local collision efficiency.

147. C BZERO - stagnation line value of the local collision
 148. C efficiency

149. C CHORD - length of airfoil chord plus accreted ice
 150. C CHORD1 - length of bare airfoil chord in metres

164 C				
165 C	DIAG	- cross diagonal of the quadrilateral of AREA		
166 C	DROPDI	- droplet diameter in microns (INPUT)		
167 C	H	- airfoil thickness as a fraction of the chord length (INPUT)		
168 C	HALFPI	- pi/2 (Fortran77 constant PARAMETER)		
169 C	LAREA	- cross sectional area of the current layer in sq. metres		
170 C	LAYERS	- integer number of layers		
171 C	LAYRX, LAYRY	- array containing non-dimensional x and y coordinates of the accretion surfaces for up to 30 accretion layers		
172 C	Ldens	- mean density of the current layer in kg/cubic metre		
173 C	LMASS	- mass of the current layer in kg per unit length		
174 C	LWC	- liquid water content in kg per cubic metre (INPUT)		
175 C	L1, L2, L3	- dimensioned sides of the quadrilateral area defined by two adjacent points on the current surface and two on the surface below		
176 C	L4	- array of masses per unit length for areas defined by adjacent points. In kg		
177 C	NPOINT	- number of points defining the initial airfoil surface (note not the same as for potential flow calculation)		
178 C	PFOILX,	- arrays of non-dimensional x,y points of the front of the initial surface. for plotting		
179 C	PFOILY			
180 C	PHI	- angle of attack of the free stream (INPUT)		
181 C	PI	- pi (Fortran77 constant PARAMETER)		
182 C	RHO	- array of local densities in kg/cubic metre		
183 C	SIMSPC	- logical variable. true if simulated spectrum option is chosen		
184 C	SIMOPT	- character variable for reading in response to simulated spectrum option query (INPUT)		
185 C				
186 C				
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1 2 3 4 5 6 7 8 9

220 C SLOPE - array of local surface slopes calculated at each surface point

221 C S1,S2 - quantities used to determine the quadrilaterals for the layer area calculation

222 C TAILEN - filter length for optional smoothing filter applied to C. curve

223 C TAU - accretion time step in seconds (INPUT)

224 C TH - accumulated stagnation line thickness

225 C THICK - array of ice thicknesses normal to surface

226 C TMASS - total mass of the accretion

227 C VEL - free stream air speed in m per sec (INPUT)

228 C XFOIL, YFOIL - arrays containing dimensional x and y coordinates of the airfoil surface

229 C *****

230 C 231 C 232 C 233 C 234 C 235 C 236 C 237 C 238 C 239 C 240 C 241 C 242 C 243 C 244 C 245 C 246 C 247 C 248 C 249 C 250 C 251 C 252 C 253 C 254 C 255 C 256 C 257 C 258 C 259 C 260 C 261 C 262 C 263 C 264 C 265 C 266 C 267 C 268 C 269 C 270 C 271 C 272 C 273 C 274 C 275 C

ISN 1 255.0 256.0 257.0 258.0 259.0 260.0 261.0 262.0 263.0 264.0 265.0 266.0 267.0 268.0 269.0 270.0 271.0 272.0 273.0 274.0 275.0

ISN 2 255.0 256.0 257.0 258.0 259.0 260.0 261.0 262.0 263.0 264.0 265.0 266.0 267.0 268.0 269.0 270.0 271.0 272.0 273.0 274.0 275.0

ISN 3 255.0 256.0 257.0 258.0 259.0 260.0 261.0 262.0 263.0 264.0 265.0 266.0 267.0 268.0 269.0 270.0 271.0 272.0 273.0 274.0 275.0

ISN 4 255.0 256.0 257.0 258.0 259.0 260.0 261.0 262.0 263.0 264.0 265.0 266.0 267.0 268.0 269.0 270.0 271.0 272.0 273.0 274.0 275.0

ISN 5 255.0 256.0 257.0 258.0 259.0 260.0 261.0 262.0 263.0 264.0 265.0 266.0 267.0 268.0 269.0 270.0 271.0 272.0 273.0 274.0 275.0

ISN 6 255.0 256.0 257.0 258.0 259.0 260.0 261.0 262.0 263.0 264.0 265.0 266.0 267.0 268.0 269.0 270.0 271.0 272.0 273.0 274.0 275.0

ISN 7 255.0 256.0 257.0 258.0 259.0 260.0 261.0 262.0 263.0 264.0 265.0 266.0 267.0 268.0 269.0 270.0 271.0 272.0 273.0 274.0 275.0

SLOPE - array of local surface slopes calculated at each surface point

S1,S2 - quantities used to determine the quadrilaterals for the layer area calculation

TAILEN - filter length for optional smoothing filter applied to C. curve

TAU - accretion time step in seconds (INPUT)

TH - accumulated stagnation line thickness

THICK - array of ice thicknesses normal to surface

TMASS - total mass of the accretion

VEL - free stream air speed in m per sec (INPUT)

XFOIL, YFOIL - arrays containing dimensional x and y coordinates of the airfoil surface

DOUBLE PRECISION AZERO,ALPHM,APHI,AREA,BZERO,
CHORD,CHORD1,DIAG,DROPDI,H,LAREA,LDENS,
LAYR(20,300),PFOILX(302),PFOILY(302),
RHO(300),SLOPE(300),THICK(300),
XFOLI(600),YFOIL(600)

DOUBLE PRECISION ALPHM,APHI,AREA,BZERO,
CHORD,CHORD1,DIAG,DROPDI,H,LAREA,LDENS,
LMASS,LWC,L1,L2,L3,L4,MASS,PHI,SI,
S2,TAILEN,TAU,TH,TMASS,VEL

INTEGER LAYER,INPOINT

CHARACTER*1 SIMOPT

LOGICAL SIMSPC

PARAMETER (PI = 3.1415926535897932,HALFPI = PI/2)

COMMON /MN/XFOIL,YFOIL,GAMMA/DR/LAYRX,LAYRY,PFOILX,PFOILY

Prepare mass table and plot description files.

OPEN (UNIT = 8, FILE = '-MASS', STATUS = 'UNKNOWN')
OPEN (UNIT = 9, FILE = '-PDF', STATUS = 'UNKNOWN')

Set up headings for mass and density table in -MASS.

WRITE (8,110)

110 FORMAT (//,.15X,'Mass of Ice Accreted',//,.5X,'Layer no.',
7X,'Mass',7X,'Mean Density',//,
20X,(kg/m)',9X,(kg/cu.m)',//,

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```

276.      +
277.      C Write program title and intro to screen.
278.      C
279.      C
12   280.      WRITE (6,120)          * * * * *
     120 FORMAT (//,*,*,*,*,*,*,*,*,*,*) AIRFOIL RIME ICING MODEL
13   281.      * * * * * //,17X, Version Mar. 12, 1986 //.
282.      * * * * * //,17X, K. J. Flinstad //,17X,
283.      * * * * * //,17X, Meteorology Division //,12X, The University of Alberta
284.      * * * * * //,17X, //, Please enter parameters to begin accretion:
285.      * * * * *
286.      C Prompt user for input parameters.
287.      C
288.      C      WRITE (6,*) 'Median Volume Droplet Diameter in microns? '
     289.      C      READ (5,*) DROPDF
290.      C      Convert to metres
291.      C
292.      C
293.      C      DROPDF = DROPDF*1.0E-6
294.      C
295.      C      Set flags for requested treatment of droplet spectrum.
296.      C
297.      C      298.      WRITE (6,*) 'Do you wish simulated spectrum weighting? '
     299.      C      WRITE (6,*) 'READ (5,*) SIMOPT
300.      C      301.      8 FORMAT (1A1)
302.      C      303.      C For simulated spectrum calculate 'tail' length from
     304.      C      median volume diameter
305.      C
306.      C      306.      IF ((SIMOPT EQ 'Y') OR ((SIMOPT EQ 'Y')) THEN
     307.      C      SIMSPC = TRUE
     308.      C      TAILEN = 0.13 * DEXP(-DROPDF / 7.0E-5)
     309.      C      ELSE
     310.      C      SIMSPC = FALSE
     311.      C      TAILEN = 0.0
     312.      C      ENDIF
313.      C
314.      C      WRITE (6,*) 'Airfoil chord length in metres? '
     315.      C      WRITE (6,*) 'READ (5,*) CHORD1
316.      C      317.      C Set current chord length equal to bare airfoil
     318.      C      319.      C
     320.      C      CHORD = CHORD1
321.      C
322.      C      WRITE (6,*) 'Airfoil height as fraction of chord? '
     323.      C      WRITE (6,*) 'READ (5,*) H
324.      C
325.      C      WRITE (6,*) 'Free stream speed in m / sec? '
     326.      C      WRITE (6,*) 'READ (5,*) VEL
327.      C      328.      C
     329.      C      WRITE (6,*) 'Liquid water content in kg / cubic metre? '
     330.      C      WRITE (6,*) 'READ (5,*) Q
331.      C

```

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      READ (5,*) LWC
      WRITE (6,*) 'Time step per layer in seconds? '
      WRITE (6,*) TAU
      READ (5,*) TAU

      WRITE (6,*) 'Total number of layers? '
      READ (6,*) LAYERS
      READ (5,*) LAYERS

      C Bare surface will be layer 1.. so add 1 to LAYERS
      LAYERS = LAYERS + 1

      WRITE (6,*) 'Angle of attack in degrees? '
      READ (6,*) PHI
      READ (5,*) PHI

      C Convert angle of attack to radians, and calculate
      C surface angle of the max collision efficiency point
      C for non-zero attack angles (AZERO).
      C
      AZERO = DLOG10(PHI + 1) / (0.001 * DEXP(300.5 * HI))
      PHI = (PHI + 2.0 * PI) / 360.0
      AZERO = (AZERO + 2.0 * PI) / 360.0

      C Fill array of points defining the initial surface of the
      C airfoil in dimensional coordinates. The complete
      C surface contains NPOINT points.
      NPOINT = 600

      DO 100 I = 1, NPOINT
        XFOIL(I) = 0.5 * (.1,0 - DCOS(2.0 * PI * DBLE(I-1)/
        DBLE(NPOINT)))
        YFOIL(I) = 5.0 * H * (0.2969 * DSQRT(XFOIL(I)) -
        0.126 * XFOIL(I) - 0.3516 * XFOIL(I) ** 2 -
        0.2843 * XFOIL(I) ** 3 -
        0.1015 * XFOIL(I) ** 4) * CHORD1
        IF (I .GT. 30) YFOIL(I) = -YFOIL(I)
      100 CONTINUE

      C Fill arrays of non-dimensional coordinates for later plotting
      DO 101 I=1, 1450
        PFOILX(150+I) = XFOIL(I) / CHORD1
        PFOILY(150+I) = YFOIL(I) / CHORD1
        PFOILX(I) = XFOIL(450+I) / CHORD1
        PFOILY(I) = YFOIL(450+I) / CHORD1
      101 CONTINUE

      C Copy front part of surface to array which will contain
      C all of the layers.
      DO 200 I = 1, NPOINT / 4

```

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```

      LAYRX(1,1) = XFOIL((3 * NPOINT / 4) + 1)
      LAYRY(1,1) = YFOIL((3 * NPOINT / 4) + 1)
      CONTINUE
      DO 210 I = (NPOINT / 4) + 1, NPOINT / 2
      LAYRX(1,I) = XFOIL(I - (NPOINT / 4))
      LAYRY(1,I) = YFOIL(I - (NPOINT / 4))
      210 CONTINUE
      C
      C Call Subroutine DROPS to calculate Beta zero. Alpha max 18
      C then determined from Beta zero.
      CALL DROPS(DROPDS, VEL, CHORD1, NPOINT, BZERO)
      ALPHM = DACOS(1.0 - BZERO)

      ISN    76      400
      ISN    77      401
      C
      Initialize total mass and begin loop to accrete ice layers
      402      C
      403      C
      404      C      TMASS = 0.0
      405      C      DO 300 J = 2,LAYERS
      406      C      Check for recalculation of BZERO every fifth layer
      407      C
      408      C      IF ((J .EQ. 6) .OR. (J .EQ. 11) .OR. (J .EQ. 16)) THEN
      409      C
      410      C      copy new shape to arrays for potential flow calculation in DROPS
      411      C
      412      C      DO 220 I=1,NPOINT / 4
      413      C      XFOIL(I + (3*NPOINT/4)) = LAYRX(J-1,1)
      414      C      YFOIL(I + (3*NPOINT/4)) = LAYRY(J-1,1)
      415      C      CONTINUE
      416      C      DO 230 I = (NPOINT / 4) + 1, NPOINT / 2
      417      C      XFOIL(I - NPOINT / 4) = LAYRX(J-1,1)
      418      C      YFOIL(I - NPOINT / 4) = LAYRY(J-1,1)
      419      C      CONTINUE
      420      C
      421      C      DO 240 I = 1, NPOINT / 4
      422      C      Update length of airfoil chord plus accreted ice
      423      C
      424      C      CHORD = CHORD + (((J-2) * BZERO + VEL * LWC * TAU)/880.0)
      425      C      Update BZERO, and ALPHM if necessary
      426      C
      427      C      CALL DROPS(DROPDS, VEL, CHORD, NPOINT, BZERO)
      428      C      IF (DACOS(1.0 - BZERO) LT. ALPHM) ALPHM =
      429      C      DACOS(1.0 - BZERO)
      430      C      ENDIF
      431      C
      432      C      Calculate slope and surface angle from finite differences for
      433      C      each point on current surface
      434      C
      435      C      DO 400 J = 2,299
      436      C      SLOPE(I) = (LAYRX(J-1,1+1) - LAYRX(J-1,1-1)) /
      437      C      DABS(LAYRY(J-1,1+1) - LAYRY(J-1,1-1))
      438      C      IF ((LAYRY(J-1,1+) .EQ. LAYRY(J-1,1-1))
      439      C      SLOPE(I) = 1.0E10
      440      C      IF ((LAYRY(J-1,1+1) LT. LAYRY(J-1,1-1)) THEN
      441      C      ALPHAI = PI - DATAN(SLOPE(I))
      442      C
      443      C

```

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100      444      ELSE
101      445      ALPHA(I) = DATAN(SLOPE(I))
102      446      ENDIF
103      447      400 continue
104      448      C Calculate for each point local density, collision efficiency and
105      449      C ice thickness. The method of calculation for Beta and rho
106      450      C depends on the values of SIMSPC and PHI.
107      451      C
108      452      DO 300 I = 2,289
109      453      IF (PHI .EQ. 0.0) THEN
110      454      RHO(I) = 880.
111      455      (1.0 - 0.163 * DABS(ALPHA(I)/(ALPHB+TAILEN))) *
112      456      * (1.0 - 0.163 * DABS(ALPHA(I)/(ALPHB+TAILEN))) *
113      457      * (0.235 * (DABS(ALPHA(I)/(ALPHB+TAILEN))) ** 2) -
114      458      * (2.357 * (DABS(ALPHA(I)/(ALPHB+TAILEN))) ** 3) +
115      459      * (1.374 * (DABS(ALPHA(I)/(ALPHB+TAILEN))) ** 4) +
116      460      * (0.0188 * (DABS(ALPHA(I)/(ALPHB+TAILEN))) ** 5)) +
117      461      ELSE
118      462      ALPHA(I) = ((-ALPHA(I)) - AZERO) / (HALFPI - AZERO)
119      463      RHO(I) = 880.0 *
120      464      (1.0 - (0.163 * DABS(ALPHA(I)/HALFPI)))
121      465      * (0.235 * DABS(ALPHA(I)/HALFPI) ** 2) -
122      466      * (2.357 * DABS(ALPHA(I)/HALFPI) ** 3) +
123      467      * (1.374 * DABS(ALPHA(I)/HALFPI) ** 4) +
124      468      * (0.0188 * DABS(ALPHA(I)/HALFPI) ** 5)) +
125      469      ENDIF
126      470      IF (ALPHA(I) .GT. HALFPI) RHO(I) = 170.
127      471      C Make sure density does not become too small
128      472      C
129      473      C
130      474      IF (RHO(I) .LT. 170.) RHO(I) = 170.
131      475      C
132      476      C Calculate local collision efficiency
133      477      C
134      478      C
135      479      IF (PHI .NE. 0.0) THEN
136      480      BETA = 0.95 * SPCBET(ALPHA,BZERO,HALFPI,O,O)
137      481      ELSE IF (SIMSPC) THEN
138      482      BETA = SPCBET(ALPHA(I),BZERO,ALPHA,TAILEN)
139      483      ELSE
140      484      BETA = ONEBET(ALPHA(I),BZERO,ALPHA)
141      485      ENDIF
142      486      C
143      487      C Calculate local ice thickness perpendicular to surface
144      488      C
145      489      THICK(I) = BETA * VEL * TAU * LMC / RHO(I)
146      490      C Define new layer surface coordinates.
147      491      C
148      492      LAYRX(J,I) = LAYRX(J-1,I) - (THICK(I) *
149      493      * DCOS(DATAN(SLOPE(I)))) +
150      494      * LAYRY(J,I) = LAYRY(J-1,I) + (THICK(I) *
151      495      * DSIN(DATAN(SLOPE(I)))) +
152      496      500 continue
153      497      LAYRX(J,1) = LAYRX(J-1,1)
154      498      LAYRX(J,300) = LAYRX(J-1,300)
155      499

```

LEVEL 1.1.1 (DEC) 811

VS FORTRAN DATE: JUN 24, 1986 TIME: 03:15:48 NAME: MAIN PAGE: 10

```

      126      500
      127      501
      128      502
      129      503
      130      504
      131      505
      132      506
      133      507
      134      508
      135      509
      136      510
      137      511
      138      512
      139      513
      140      514
      141      515
      142      516
      143      517
      144      518
      145      519
      146      520
      147      521
      148      522
      149      523
      150      524
      151      525
      152      526
      153      527
      154      528
      155      529
      156      530
      157      531
      158      532
      159      533
      160      534
      161      535
      162      536
      163      537
      164      538
      165      539
      166      540
      167      541
      168      542
      169      543
      170      544
      171      545
      172      546
      173      547
      174      548
      175      549
      176      550
      177      551
      178      552
      179      553
      180      554
      181      555

```

```

      LAYER(J,1) = LAYER(J-1,1)
      LAYER(J,300) = LAYER(J-1,300)

      C Calculate mass and mean density of the layer from
      C the sum of areas between surface points on successive
      C layers. First, initialize layer mass and area.
      C

      LMASS = 0.0
      LAREA = 0.0
      DO 560 I = 2, 286
      IF ((THICK(I), GT, 0.0) AND (THICK(I+1), GT, 0.0)) THEN
        L4 = THICK(I)
        L2 = THICK(I+1)
        L1 = DSORT((LAYERX(J-1,I+1) - LAYERX(J-1,I))**2 + (LAYER
          /J-1,I+1) - LAYER(J-1,I))**2)
        L3 = DSORT((LAYERX(J,I+1) - LAYER(J,I))**2 + (LAYER
          (J,I+1) - LAYER(J,I))**2)
        DIAG = DSORT((LAYERX(J,I+1) - LAYERX(J-1,I))**2 + (LAYER
          (J,I+1) - LAYER(J,I))**2)
        S1 = 0.5 * (DIAG+L1+L2)
        S2 = 0.5 * (DIAG+L3+L4)
        AREA = DSORT(S1 * (S1-DIAG) * (S1-L2) +
          DSORT(S2 * (S2-DIAG) * (S2-L3) * (S2-L4)))
        MASS = ((RHO(I) * RHO(I+1)) / 2.0) * AREA
        LMASS = MASS + MASS
        LAREA = LAREA + AREA
      ENDIF
      CONTINUE
      560  C Calculate mean density.
      529  C
      530  C
      531  C
      532  C
      533  C
      534  C
      535  C
      536  C
      537  C
      538  C
      539  C
      540  C
      541  C
      542  C
      543  C
      544  C
      545  C
      546  C
      547  C
      548  C
      549  C
      550  C
      551  C
      552  C
      553  C
      554  C
      555  C

```

```

      WRITE (8,570) J-1, LMASS, LDENS
      570  FORMAT (/,.8X,12.7X, F9.9, T9, F9.9)
      530  C Add mass of current layer to total mass.
      539  C
      TMASS = TMASS + LMASS
      540  C
      541  C Shift surface points closer to nose by a decreasing
      C interval between points
      542  C
      543  C
      544  C
      545  C
      546  C
      547  C
      548  C
      549  C
      550  C
      551  C
      552  C
      553  C
      554  C
      555  C

```

```

      IF ((J, GT, 2) AND (PM1, EQ, 0.0)) THEN
        DO 560 I=1,149
        IF ((THICK(I+1)-1), GT, 0.0) THEN
          LAYER(J,151-1) = LAYER(J,151-1) -
            ((0.4 * (20 / DBLE(I+1))) *
             DABSLAYER(J,151-1) - LAYER(J,152-1))
          LAYER(J,151-1) = LAYER(J,151-1) +
            ((0.4 * (20 / DBLE(I+1))) *
             DABS(LAYER(J,151-1) - LAYER(J,152-1)))
        ENDIF
        IF (THICK(I+1), GT, 0.0) THEN

```

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```

      1.....2.....3.....4.....5.....6.....7.....8.....9.

  ISN 156   556.    LAYRX(J,I+151) = LAYRX(J,I+151)
  ISN     557.    +   ((0.4 * (20./DBLE(I+19))) *
  ISN     558.    +   DABS(LAYRX(J,I+150)-LAYRX(J,I+151))
  ISN 157   559.    LAYRY(J,I+151) = LAYRY(J,I+151)
  ISN     560.    +   ((0.4 * (20./DBLE(I+19))) *
  ISN     561.    +   DABS(LAYRY(J,I+150)-LAYRY(J,I+151)))
  ISN 158   562.    ENDIF
  ISN 159   563.    CONTINUE
  ISN 160   564.    ENDIF
  ISN 161   565.    C 900 CONTINUE
  ISN     566.    C Write total mass accreted to table
  ISN     567.    C 568. C Write (8.590) TMASS
  ISN     569.    C 570. WRITE (8.590) TMASS
  ISN 162   571.    C 590 FORMAT (8X,/, Total mass accreted per metre: ,F9.5)
  ISN 163   572.    C 573. C Call subroutine to plot airfoil and accreted layers
  ISN     574.    C 575. CALL DRAW(ALPHM,BZERO,CHORD1,DROPDI,LAYERS,LWC,PHI,
  ISN     576.    C 577. SIMSPC,TAILEN,TAU,VEL)
  ISN 164   578.    C 579. 999 STOP
  ISN 165   579.    END
  ISN 166   579.    END

*STATISTICS* SOURCE STATEMENTS = 162, PROGRAM SIZE = 21362 BYTES, PROGRAM NAME = MAIN PAGE: 1
*STATISTICS* NO DIAGNOSTICS GENERATED.
***** END OF COMPIRATION 1 *****

```

```

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:15:50 NAME: MAIN PAGE: 12
REQUESTED OPTIONS (PROCESS): SC(Axis2.Symbol)

OPTIONS IN EFFECT: NOLIST NOMAP NOXREF NOGOSTMT NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(0) LANGLVL(77) NOFIPS FLAG(1) NAME(MAIN) LINECOUNT(60)

* * * * * 1.....2.....3.....4.....5.....6.....7.....8.....9

581 C
582 * SUBROUTINE DRAW - to plot airfoil and all accretion
583 * layers, with appropriate information
584 * in the legend. This subroutine uses
585 * the plotting library *PLOTLIB
586 *
587 *
588 * Input Variables : as defined for Main program :
589 * ALPHM,BZERO,CHORD1,DROPDI,LAYERS,LWC,
590 * PHI,SIMSPC,TAILEN,TAU,VEL
591 *
592 * Common Variables : PLOTX,PLOTY,PFOILX,PFOILY
593 *
594 * Other Variables : PLOTX,PLOTY - arrays containing x,y
595 * plotting data and scaling information
596 * for accretion layer currently being
597 * plotted
598 *
599 *

600 C SUBROUTINE DRAW(ALPHM,BZERO,CHORD1,DROPDI,LAYERS,
601 * LWC,PHI,SIMSPC,TAILEN,TAU,VEL)
602 +
603 C
604 C DOUBLE PRECISION PFOILX(302),PFOILY(302),LAYRX(20,300),
605 * LAYRY(20,300),PL0TX(302),PL0TY(302),
606 C
607 C DOUBLE PRECISION ALPHM,BZERO,CHORD,DROPDI,LWC,PHI,
608 * TAILEN,TAU,VEL
609 C
610 C INTEGER LAYERS
611 C LOGICAL SIMSPC
612 C
613 F COMMON /DR/LAYRX,LAYRY,PFOILX,PFOILY
614 C
615 C To plot airfoil surface.
616 C place scaling data at the end of coordinate arrays.
617 C
618 C
619 C PFOILX(301) = -0.4
620 C PFOILX(302) = 0.1
621 C PFOILY(301) = -0.25
622 C PFOILY(302) = 0.1
623 C PL0TX(301) = -0.4
624 C PL0TX(302) = 0.1
625 C PL0TY(301) = -0.25
626 C PL0TY(302) = 0.1
627 C Open plotting library
628 C
629 C CALL PLOTS
630 C CALL METRIC(O)
631 C

```

LEVEL 1.1.1 (DEC 81)

DATE: JUL 24, 1986 TIME: 13:15:50

NAME: DRAW PAGE: 13

VS FORTRAN

1.....1.....2.....3.....4.....5.....6.....7.....8.....9

```

632 C Draw 4 by 6 .inch box
633 C
634 C - CALL PLOT(2.0,2.0,-3)
635 C - CALL PLOT(0.0,2.0,3)
636 C - CALL PLOT(0.0,6.0,3)
637 C - CALL PLOT(9.0,6.0,2)
638 C - CALL PLOT(9.0,0.0,2)
639 C
640 C Label axes and draw tics
641 C CALL AXIS2(0.0,0.0,'X/CHORD',-7.9,0.0,0,-0.4,0,1,0)
642 C CALL AXIS2(0.0,0.0,'Y/CHORD',7.6,0.9,0,-0.25,0,1,0.5)
643 C
644 C Write legends tq top left corner
645 C
646 C
647 C
648 C
649 C
650 C
651 C
652 C
653 C
654 C
655 C
656 C
657 C
658 C
659 C
660 C
661 C
662 C
663 C
664 C
665 C
666 C
667 C
668 C
669 C
670 C
671 C
672 C
673 C
674 C
675 C
676 C Draw initial airfoil surface
677 C
678 C
679 C
680 C Transfer each layer in turn to the array holding
681 C the scaling parameters
682 C
683 C
684 C
685 C
686 C
687 C
52 688 C
53 689 C
54 690 C
55 691 C
56 692 C
191

```

191

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:15:50 NAME: DRAW PAGE: 14
.....1.....2.....3.....4.....5.....6.....7.....8.....9

ISN 57 688. 700 CONTINUE
689. C
690. C Draw accretion layer
691. CALL LINE(PLOTX,PLOTY,300,2,0)
ISN 58 692. 610 CONTINUE
ISN 59 693. C
694. C Close Plotting library
695. C
696. C CALL PLOT(10,10,999)
ISN 60 697. RETURN
ISN 61 698.
ISN 62 699. END

SOURCE STATEMENTS = 62. PROGRAM SIZE = 7298 BYTES. PROGRAM NAME = DRAW PAGE: 12.

STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPILED 2 *****

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:15:50 NAME: MAIN PAGE: 15

OPTIONS IN EFFECT: NOLIST NOMAP NOXREF NOGOSTMT NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(O) LANG(VL(77)) NOTIPS FLAG(1) NAME(MAIN) LINECOUNT(60)

193

```
      C *****  
      C * FUNCTION ONEBET - Calculates local collision efficiency  
      C * estimate for a monodisperse droplet populations  
      C *  
      C * Input Variables : as defined for main program:  
      C *          ALPHA,BZERO,ALPHM,SHDROP  
      C *  
      C *  
      C *          DOUBLE PRECISION FUNCTION ONEBET(ALPHA,BZERO,ALPHM)  
      C *          DOUBLE PRECISION BZERO,ALPHM,ALPHA  
      C *  
      C *          IF (DABS(ALPHA) .GT. ALPHM) THEN  
      C *              ONEBET = 0.0  
      C *          ELSE  
      C *              ONEBET = DCOS(ALPHA) - (1.0 - BZERO)  
      C *          ENDIF  
      C *          RETURN  
      C *          END  
      C *  
      ISN   1    710  
      ISN   1    711  
      ISN   2    712  
      ISN   2    713  
      ISN   3    714  
      ISN   3    715  
      ISN   4    716  
      ISN   5    717  
      ISN   6    718  
      ISN   7    719  
      ISN   8    720  
      ISN   9    721  
      C *  
      C * SOURCE STATEMENTS = 9. PROGRAM SIZE = 870 BYTES. PROGRAM NAME = ONEBET PAGE: 15.  
      C * STATISTICS* NO DIAGNOSTICS GENERATED.  
      C * STATISTICS* ***** END OF COMPILATION 3 *****
```

LEVEL 1.1.1 (DEC 81) VS FORTRAN { DATE: JUL 24, 1986 TIME: 13:15:50 NAME: MAIN PAGE: 16

OPTIONS IN EFFECT: NOLIST NOMAP NOXREF NODECK SOURCE TERM OBJECT FIXED
 OPTIMIZE(O) LANGLVL(77) NOFIPS FLAG(1) NAME(MAIN) LINECOUNT(60)

```

*.... *....1.....2.....3.....4.....5.....6.....7.....8.....9.....
```

```

722. C *****
723. C * Function SPCBET - Calculates the local c.e. for a
724. C * simulated droplet size distribution
725. C * scaled to the median volume droplet,
726. C * and to (alpha max + tail). The tail
727. C * length was determined in the main program.
728. C *
729. C * Input Variables : as defined for Main program:
730. C *          ALPHA,BZERO,SMDROP,TAILEN
731. C *
732. C * Other Variables : X - scaled angle = Alpha /
733. C *          (Alpha + Taillen)
734. C *
735. C *
736. C *
737. C DOUBLE PRECISION FUNCTION SPCBET(ALPHA,BZERO,  

     .          ALPHM,TAILEN)
    ISN   1   738. +
    ISN   2   739. +
    ISN   3   740. C
    ISN   4   741. C DOUBLE PRECISION ALPHA,ALPHM,BZERO,TAILEN,X
    ISN   5   742. C
    ISN   6   743. C ALPHA = DABS(ALPHA)
    ISN   7   744. C IF (ALPHA .GE. (ALPHM+TAILEN)) THEN
    ISN   8   745. C      SPCBET = 0.0
    ISN   9   746. C ELSE
    ISN  10   747. C      X = ALPHA / (ALPHM+TAILEN)
    ISN  11   748. C      SPCBET = BZERO * (1.000 + 0.039 * X - 1.842 *
    ISN  12   749. C                  * (X ** 2) - 0.543 * (X ** 3) +
    ISN  13   750. C                  * (X ** 4) - 0.444 * (X ** 5))
    ISN  14   751. C      IF (SPCBET .LT. 0.0) SPCBET = 0.0
    ISN  15   752. C      IF (SPCBET .GT. BZERO) SPCBET = BZERO
    ISN  16   753. C ENDIF
    ISN  17   754. C RETURN
    ISN  18   755. C END

  •STATISTICS• SOURCE STATEMENTS = 13. PROGRAM SIZE = 608 BYTES. PROGRAM NAME = SPCBET PAGE: 16
  •STATISTICS• NO DIAGNOSTICS GENERATED.
  ••••• END OF COMPILEATION 4 •••••
```

LEVEL 1.1.1 (DEC 81)

VS FORTRAN

DATE: JUL 24, 1986 TIME: 13:15:51

NAME: MAIN PAGE: 17

OPTIONS IN EFFECT: NOLIST NOMAP NOXREF NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(O) LAMESVL(77) NDFIPS FLAG(I)- NAME(MAIN) LINECOUNT-(60)

..... 1 2 3 4 5 6 7 8 9

756 C
757 C SUBROUTINE DROPS

758 C
759 C
760 C This routine calculates a trajectory pair for water
C droplets in a potential flow field around a NACA airfoil or
761 C similar accretion shape. The stagnation line collision
762 C efficiency is then determined.

763 C
764 C The equation of motion of the droplet is integrated
765 C by the Haun method (Meslinger and Arakawa, 1976; Numerical
766 C Methods used in Atmospheric Models, Volume 1, W.M.O. GARP
767 C Publications Series No. 17). The time step is set to
768 C 1/400 of the time to travel one chord length at the free
769 C stream speed. A maximum of 3000 steps are allowed.
770 C by which time the droplet is presumed to have collided
771 C with the airfoil or to have passed it.

772 C
773 C The droplet's initial x-coordinate is set to an
774 C upstream distance of 5.0 chord lengths. At this
775 C distance the x-speed is assumed to be equal to the air
776 C speed in the x-direction, and the droplet y-speed is
777 C set to zero.

778 C
779 C A first trajectory is integrated beginning at y = zero,
780 C which is then incremented by a small amount (DELTAY)
781 C for the second trajectory. If both trajectories
782 C intercept the airfoil, the local collision efficiency, and
783 C the angle between the focal surface normal and free stream
784 C direction are calculated.

785 C
786 C The formulation of Beard and Pruppacher (J.A.S., Vol. 26,
787 C p. 1066, 1969) is used for the drag coefficient. Potential
788 C flow calculations are made according to the method of Kennedy
789 C and Marsden (Can. Aero. and Space J., Vol. 22, p. 243, 1976).

790 C
791 C One external subroutine is required, from the IMSL library:
792 C the subroutine LEQTIF, to solve a system of equations
793 C in the potential flow calculation.

794 C
795 C The following internal subprograms are called from DROPS
796 C
797 C
798 C
799 C
800 C
801 C SUBROUTINE COLLID - calculates final impact position
802 C if the droplet collides with the surface.

803 C
804 C SUBROUTINE FLOWXY- calculates the air speeds at a
805 C given x,y location using the results of POFLO,
806 C and droplet accelerations (from the equations of
807 C motion) for the current iteration and time step.
808 C

LEVEL 1.1.1 (DEC) 81)

VS FORTRAN DATE: JUL 24, 1986 TIME: 13:15:51

NAME: MAIN PAGE: 18

809. C SUBROUTINE KCALC - calculates the matrix elements .
 810. C K used in POFLO and FLOWXY .
 811. C
 812. C SUBROUTINE POFLO - calculates the potential vorticities
 at the surface of the airfoil according to the
 method of Kennedy and Marsden .
 813. C
 814. C
 815. C SUBROUTINE TRAJEC - calculates droplet trajectories
 by integrating the droplet equations of motion by
 the trapezoid rule with one iteration step .
 816. C
 817. C
 818. C
 819. C
 820. C
 821. C Input Variables:
 822. C CHORD - airfoil chord length in metres
 823. C DROPDI - droplet diameter in microns
 824. C
 825. C
 826. C
 827. C
 828. C NPOINT - number of points in the complete airfoil
 surface (must be divisible by 15)
 829. C VEL - free stream air speed in metres per second
 830. C
 831. C
 832. C Output Variables:
 833. C BZERO - stagnation line collision efficiency
 834. C
 835. C Common Variables: (with Main)
 836. C
 837. C
 838. C GAMMA - vector of potential vorticities for
 potential flow calculation
 839. C
 840. C
 841. C XFOIL - vectors containing points which define
 YFOIL - the complete dimensioned airfoil surface .
 842. C
 843. C
 844. C Other Variables:
 845. C
 846. C ALPHA - angle between the local surface normal
 and the free stream direction at
 the impact point
 847. C
 848. C
 849. C
 850. C
 851. C ARC - arc length along the airfoil surface
 between the impact points of the two
 trajectories (estimated by a straight
 line segment)
 852. C
 853. C DELTAT - time step in seconds
 854. C
 855. C
 856. C DELTAY - difference in initial y position of
 the two trajectories
 857. C
 858. C
 859. C FINX1, FINX2 - x-positions of the droplets at impact
 for each trajectory
 860. C
 861. C
 862. C
 863. C FINY1, FINY2 - y-positions of the droplets at impact
 for each trajectory
 864. C

LEVEL 1.1.1 (DEC 81)

VS FORTRAN DATE: JUL 24, 1986 TIME # 13:15:51

NAME: MAIN PAGE: 19

..... 2

..... 3

VS FORTRAN

..... 4

..... 5

..... 6

..... 7

..... 8

..... 9

C HALFPI - PI / 2 (FORTRAN77 constant parameter)
 866 C
 867 C HIT1, - logical flags for each trajectory
 868 C HIT2 which indicate whether a collision
 869 C with the airfoil has occurred
 870 C
 871 C NSTEPS - maximum number of time steps to be
 872 C integrated
 873 C
 874 C PI - PI (FORTRAN77 constant parameter)
 875 C
 876 C SLOPE1, - slopes of the airfoil or accretion
 877 C SLOPE2 surface at the impact points
 878 C
 879 C XINIT, - initial droplet x,y position in metres
 880 C YINIT
 881 C
 882 C
 883 C
 884 C
 885 C SUBROUTINE DROPS(DROPDI, VEL, CHORD, NPOINT, BZERO)
 ISN 1 886 C
 ISN 2 887 C DOUBLE PRECISION XFOIL(600), YFOIL(600), GAMMA(41)
 888 C
 ISN 3 889 C DOUBLE PRECISION ARC,BZERO,CHORD,DELTAT,DELTAY,
 890 C DROPO1,FINX1,FINX2,FINY1,FINY2,
 891 C SLOPE1,SLOPE2,VEL,XINIT,YINIT
 892 C
 893 C INTEGERS NSTEPS,NPOINT
 ISN 4 894 C LOGICAL HIT1,HIT2
 ISN 5 895 C
 ISN 6 896 C
 ISN 7 897 C
 ISN 8 898 C
 ISN 9 899 C
 ISN 10 900 C
 ISN 11 901 C
 902 C Set maximum number of time steps, initial x-coordinate, and
 903 C the difference in initial y-coordinates
 904 C
 905 C
 ISN 8 906 C
 ISN 9 907 C DELTAY = 1.0E-8
 ISN 10 908 C XINIT = 5.0 * CHORD
 ISN 11 909 C YINIT = 0.0
 910 C Calculate time step
 911 C
 912 C DELTAT = CHORD / (VEL * 400)
 ISN 12 913 C
 914 C Calculate potential vorticity matrix for air speed calculations
 915 C
 916 C CALL POFL0(NPOINT,CHORD)
 ISN 13 917 C
 918 C Integrate first trajectory
 919 C
 920 C

```

LEVEL 1.1.1 (DEC 81)          VS FORTRAN          DATE: JUL 24, 1986   TIME: 13:15:51      NAME: DROPS PAGE: 20
                                * 1.....1.....2.....3.....4.....5.....6.....7.....8.....9.....*
ISN    14   921.           CALL TRAJEC (XINIT,YINIT,DROPDI,VEL,CHORD,DELTAT,NSTEPS,
ISN    14   922.           + NPOINT,HIT1,FINX1,FINY1,SLOPE1)
ISN    14   923.           C Correct initialy position and integrate second trajectory
ISN    14   924.           C
ISN    15   925.           C YINIT = YINIT + DELTAY
ISN    15   926.           CALL TRAJEC (XINIT,YINIT,DROPDI,VEL,CHORD,DELTAT,NSTEPS,
ISN    16   927.           + NPOINT,HIT2,FINX2,FINY2,SLOPE2)
ISN    16   928.           C
ISN    16   929.           C Test for result of trajectory (miss, graze or collision)
ISN    17   930.           C
ISN    17   931.           C IF ((HIT1) AND. (HIT2)) THEN
ISN    17   932.           C
ISN    17   933.           C Calculate impact parameters and local collision efficiency
ISN    18   934.           C
ISN    18   935.           C ARC = DSORT(((FINX2 - FINX1) ** 2) + ((FINY1 - FINY2)
ISN    18   936.           C ** 2))
ISN    19   937.           C BZERO = DELTAY / ARC
ISN    19   938.           C
ISN    19   939.           C If droplet didn't hit the foil, write message to screen
ISN    19   940.           C
ISN    19   941.           C ELSE IF ((NOT HIT1) OR. (NOT HIT2)) THEN
ISN    20   942.           C     WRITE (6,40)
ISN    21   943.           C     FORMAT (//,'Droplet did not reach the airfoil.')
ISN    22   944.           C
ISN    23   945.           C
ISN    23   946.           C     RETURN
ISN    24   947.           C
ISN    25   948.           C     END
ISN    25   949.           C
*STATISTICS* SOURCE STATEMENTS = 25. PROGRAM SIZE = 926 BYTES. PROGRAM NAME = DROPS PAGE: 17
*STATISTICS* NO DIAGNOSTICS GENERATED
***** END OF COMPILED 5 *****

```

LEVEL 1.1.1 (DEC 81)

VS FORTRAN DATE: JUL 24, 1986 TIME: 13:15:51

NAME: MAIN PAGE: 21

OPTIONS IN EFFECT: NDLIST NOXREF NOGOSTMT NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(0) LANG(VL(77)) NOFIPS FLAG(1) NAME(MAIN) LINECOUNT(60)

* 1 2 3 4 5 6 7 8 9

949 C
950 C
951 C

SUBROUTINE TRAJEC

Integrates trajectory of droplet whose initial position is specified in the parameter list. Calls POFL0 to set up matrix for potential flow calculations. FLOWXY for iterative calculations of air speed and droplet acceleration, and COLLID to determine impact points.

INPUT VARIABLES:

CHORD	- airfoil chord length
DELTAT	- time step
DROPOL	- droplet diameter
NPOINT	- number of surface points
NSTEPS	- maximum number of time steps
VEL	- free stream air speed
XINIT	- initial x position
YINIT	- initial y position

OUTPUT VARIABLES:

FINALX	- droplet x velocity on impact
FINALY	- droplet y velocity on impact
HIT	- logical flag set to TRUE if a collision occurs
SLOPE	- slope of the surface at impact

COMMON VARIABLES:

CO	- vector of cosines of angles of the control elements
DELTA	- vector of half-lengths of the control elements
GAMMA	- output array of POFL0, passed to FLOWXY
SI	- vector of sines of angles of the control elements
XCON	- x,y coordinates of the control points
YCON	- vectors of initial surface coordinates
XFOIL	-
YFOIL	-

OTHER VARIABLES:

DVX	- arrays containing x,y components of droplet acceleration
DVY	- counter for searching surface array
L	- initial value of counter L
VM,VY	- arrays containing x,y components of droplet speed for each step
X,V	- arrays containing x,y positions for each time step
C	-

```

LEVEL 1.1.1 (DEC 81)   VS FORTRAN    DATE: JUL 24, 1986    TIME: 13:15:51    NAME: MAIN    PAGE: 22

      C*****1.....2.....3.....4.....5.....6.....7.....8.....9.....0.....
1002. C
1003. C      SUBROUTINE TRAJEC(XINIT,YINIT,DROP01,VEL,CHORD,DELTAT,
1004. C
1005. C      +
1006. C      DOUBLE PRECISION X(2500),Y(2500),GAMMA(41),XFOIL(600),
1007. C      YFOIL(600),VX(2500),VY(2500),DVX(2500),DVY(2500),
1008. C      DVY(2500),DELTA(40),CO(40),SI(40),
1009. C
1010. C      +
1011. C      DOUBLE PRECISION XINIT,YINIT,DROP01,CHORD,VEL,DELTAT,
1012. C      FINALX,FINALY,SLOPE
1013. C
1014. C      INTEGER NSTEPS,NPOINT,LIMIT
1015. C      LOGICAL HIT
1016. C      COMMON /MN/ XFOIL,YFOIL,GAMMA/DRP/DELTA,CO,SI,XCON,YCON
1017. C
1018. C
1019. C
1020. C
1021. C      Initialize values for integration
1022. C
1023. ISN 7      X(1) = XINIT
1024. ISN 8      Y(1) = YINIT
1025. ISN 9      VV(1) = 0.0
1026. ISN 10     VX(1) = VEL
1027. ISN 11     Y(1) = 0
1028. ISN 12     HIT = FALSE
1029. C
1030. C      Begin integration
1031. C
1032. ISN 13     I = I + 1
1033. ISN 14     IF (I .GE. NSTEPS) GOTO 250
1034. C
1035. C      Calculate current step air speeds and accelerations
1036. C
1037. ISN 15     CALL FLOWXY(X(I)),Y(I),VX(I),VY(I),VEL,CHORD,DROP01,
1038. C              NPOINT,DVX(I),DVY(I))
1039. C
1040. C      Calculate next step droplet speeds and positions
1041. C
1042. ISN 16     VX(I+1) = VX(I) - (DELTAT * DVX(I))
1043. ISN 17     VV(I+1) = VV(I) + (DELTAT * DVY(I))
1044. ISN 18     X(I+1) = X(I) + (VX(I) * DELTAT) - (O.S * DVX(I))
1045. C              * (DELTAT * DELTAT)
1046. ISN 19     Y(I+1) = Y(I) + (VY(I) * DELTAT) + (O.S * DVY(I))
1047. C              * (DELTAT * DELTAT)
1048. C
1049. C      Calculate next step air speeds and accelerations
1050. C
1051. ISN 20     CALL FLOWXY(X(I+1)),Y(I+1),VX(I+1),VY(I+1),VEL,
1052. C              CHORD,DROP01,NPOINT,DVX(I+1),DVY(I+1))
1053. C
1054. C      Iterate calculations of next step droplet speed and
1055. C      position using averaged values of accelerations
1056. C
1057. ISN 21     VX(I+1) = VX(I) - ((DELTAT / 2.0) * (DVX(I)))

```

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUN 24, 1986 TIME: 13:18:51 NAME: TRAJEC PAGE: 23

```

1058      DVX(I+1)) )
1059      VY(I+1) = VV(I) + ((DELTAT / 2.0) * DVY(I) +
1060                  DVY(I+1)))
1061      X(I+1) = X(I) + (VX(I) + DELTAT) * (0.25 *
1062      (DVX(I) + DVX(I+1)) * (DELTAT * DELTAT))
1063      Y(I+1) = Y(I) + (VY(I) + DVY(I) * DELTAT) * (0.25 *
1064      (DVY(I) + DVY(I+1)) * (DELTAT * DELTAT))
1065
1066      C If trajectory has reached the nose
1067      C test whether trajectory has intercepted the airfoil surface
1068      C Set LIMIT to begin search on upper or lower surface
1069      C If the droplet has reached to within one current chord
1070      C length of the nose
1071
1072      LIMIT = 0
1073      L = LIMIT
1074      L = LIMIT
1075      CONTINUE
1076      L = L + 1
1077      IF (L .GE. (LIMIT + 125)) GOTO 100
1078
1079      C Search for adjacent pair of points which bracket the
1080      C y-coordinate of current trajectory position (if pair
1081      C cannot be found, continue with next integration step)
1082
1083      IF ((YFOIL(L+1) .GT. Y(I+1)) .AND. (YFOIL(L) .LT.
1084                      Y(I+1))) THEN
1085
1086      C Check if current position has passed the airfoil surface
1087      C between the two points
1088
1089      IF ((XFOIL(L+1) .LT. X(I+1)) .OR. (XFOIL(L) .LT.
1090                  X(I+1))) THEN
1091          HIT = TRUE
1092
1093      C Determine the impact point on the surface, and local
1094      C surface slope at that point - then exit from subroutine
1095
1096      CALL COLLID(X(I), X(I+1), Y(I), Y(I+1), XFOIL(L),
1097                  YFOIL(L), XFOIL(L+1), YFOIL(L+1),
1098                  FINALX_FINALY)
1099      SLOPE = (YFOIL(L+1) - YFOIL(L)) / DASS(XFOIL
1100                  (L+1) - XFOIL(L))
1101
1102
1103      C If not inside surface, continue integration
1104
1105      C
1106      GOTO 100
1107      ENDIF
1108
1109      C Else continue search for bracketing pair
1110
1111      C
1112      GOTO 200
1113      ENDIF

```

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:15:51 NAME: TRAJEC PAGE: 24
* * * * * 2.....3.....4.....5.....6.....7.....8.....9.....

ISN 42 1114 250 CONTINUE
ISN 43 1115 C RETURN
ISN 44 1116 END
STATISTICS SOURCE STATEMENTS = 44. PROGRAM SIZE = 122570 BYTES. PROGRAM NAME = TRAJEC PAGE: 21.
STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPIRATION 6 *****

LEVEL 1.1.1 (DEC 81)

VS FORTRAN

DATE: JUL 24. 1986 TIME: 13:15:52

NAME: MAIN PAGE: 25

OPTIONS IN EFFECT: NOLIST NOXREF NOGOSTMT NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(0) LANGVL(77) NOFIPS FLAG(1) NAME(MAIN) LINECOUNT(60).

1.....1.....2.....3.....4.....5.....6.....7.....8.....9.....

1118 C *
1119 C *
1120 C * SUBROUTINE POFL0 - Uses the method of Kennedy and
1121 Marsden (ref) to solve for the matrix of potential
1122 vorticities about the airfoil surface. A reduced
1123 number of points is used to define the surface for
1124 this subroutine than for the accretion section.
1125 Subroutine KCALC is called to calculate the matrix
1126 of K values.
1127 C *
1128 C * INPUT VARIABLES : CHORD - airfoil chord length
1129 C * NPOINT - number of points defining
1130 C * airfoil surface
1131 C *
1132 C * COMMON VARIABLES : CO,SI - vectors of cosines and sines
1133 C * of the control element angles
1134 C * DELTA - vector containing control
1135 C * element half lengths
1136 C * GAMMA - vector containing the
1137 C * potential vorticities on
1138 C * the airfoil surface
1139 C * XCON.. - vectors containing control
1140 C * point coordinates
1141 C * XFOIL - vectors of surface
1142 C * YFOIL - point coordinates
1143 C *
1144 C * OTHER VARIABLES : IDGT, routine
1145 C * IER
1146 C * K - vector input to matrix
1147 C * solver subroutine LEQT1F
1148 C * N - NPOINT/15, the no. of surface
1149 C * points used to calculate
1150 C * the stream function
1151 C * WKAREA, - input required for LEQT1F
1152 C * XF,YF - reduced vectors of surface
1153 C * points (length NPOINT/15)
1154 C *
1155 C *
1156 C *
1157 C *
ISN 1 1158 C SUBROUTINE POFL0(NPOINT,CHORD)
1159 C
ISN 2 1160 C DOUBLE PRECISION GAMMA(41),XF0IL(600),YFOIL(600).
1161 C + XCON(41),YCON(41),XF(41),YF(41).
1162 C + K(41,48),WKAREA(41),DELTA(40).
1163 C + CO(40),\$1(40),CHORD
1164 C
ISN 3 1165 C INTEGER NPOINT,N, IDGT,IER
1166 C
ISN 4 1167 C COMMON /MN/ XFOIL,YFOIL,GAMMA/DRP/DELTA,CO,SI,XCON,YCON
1168 C Construct new surface from every 15th point of old surface
1169 C non-dimensionalised to the current chord length
1170 C

```

1171      C      N = NPOINT / 15
1172      C      XF(1) = XFOIL(1) / CHORD
1173      C      YF(1) = YFOIL(1) / CHORD
1174      C      XF(N+1) = XF(1)
1175      C      YF(N+1) = YF(1)
1176      C      DO 10 I=2, N
1177      C      XF(I) = XFOIL((15 * (I-1)) + 1) / CHORD
1178      C      YF(I) = YFOIL((15 * (I-1)) + 1) / CHORD
1179      C      10 CONTINUE
1180      C
1181      C      Form arrays of control points
1182      C
1183      C      DO 20 I = 1,N
1184      C      XCON(I) = ((XF(I) + XF(I+1)) / 2.0
1185      C      YCON(I) = (YF(I) + YF(I+1)) / 2.0
1186      C      20 CONTINUE
1187      C
1188      C      Define the trailing point
1189      C
1190      C      XCON(N+1) = ((XF((N/2)+1) - XCON(N/2)) * 0.01) + XF((N/2)+1)
1191      C      YCON(N+1) = 0.0
1192      C
1193      C      Calculate R (as defined in Kennedy and Marsden) for each
1194      C      control point, store in GAMMA
1195      C
1196      C      DO 30 J = 1,N+1
1197      C      GAMMA(J) = /YCON(J))
1198      C
1199      C      Set last column of K(I,J) to 1.0
1200      C
1201      C      IF (J .EQ. N+1) THEN
1202      C          DO 35 I=1,N+1
1203      C              K(I,N+1) = 1.0
1204      C          35 CONTINUE
1205      C
1206      C      For all other J, calculate the control element half lengths
1207      C      and functions of the angle to the x-axis
1208      C
1209      C
1210      C      ELSE
1211      C          DELTA(J) = DSQRT(((XF(J+1) - XF(J)) ** 2) +
1212      C                         ((YF(J+1) - YF(J)) ** 2)) / 2
1213      C          CO(J) = (XF(J+1) - XF(J)) / (2.0 * DELTA(J))
1214      C          SI(J) = (YF(J+1) - YF(J)) / (2.0 * DELTA(J))
1215      C
1216      C      Calculate K for all remaining I,J
1217      C
1218      C      DO 40 I=1,N
1219      C          CALL KCALC(XCON(J), YCON(J), XCON(I), YCON(I),
1220      C
1221      C          40 CONTINUE
1222      C
1223      C
1224      C      Call IMSLPLIB routine to solve matrix equation;
1225      C      output matrix of potential vorticites is in GAMMA
1226      C

```

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:15:52 NAME: POFLO PAGE: 27

1227 C
1228 IDGT = B
1229 CALL EOT1F(K,1,N+1,GAMMA, IDGT,WKAREA,IERR)
1230 WRITE (6,*) GAMMA
1231 C
1232 RETURN
1233 END

STATISTICS SOURCE STATEMENTS = 39. PROGRAM SIZE = 16358 BYTES. PROGRAM NAME = POFLO PAGE: 25.

STATISTICS NO DIAGNOSTICS GENERATED.

***** END OF COMPILED 7 *****

LEVEL 1.1.1 (DEC 81) VS FORTRAN
 OPTIONS IN EFFECT: NOLIST NOMAP NOXREF NOGOSTMT NODECK SOURCE TERM OBJECT FIXED
 OPTIMIZE(O) LANG.VL(77) NOFIPS FLAG(I) NAME(MAIN) LINECOUNT(60)

		NAME: MAIN	PAGE:	26
		NOTEST	SEQ	
		2	3	4
		5	6	7
		8	9	
1234	C			
1235	C			
1236	C	ROUTINE FLCWXY - calculates for the current step:		
1237	C	air speed components from matrix equation for		
1238	C	potential flow streamline using the output		
1239	C	matrix GAMMA from POFLO, and KCALC to generate the		
1240	C	matrix elements K. Droplet accelerations are then		
1241	C	calculated according to the equations of motion		
1242	C	from the air speed, droplet Reynolds number, and		
1243	C	drag coefficient according to the formulation of		
1244	C	Beard and Pruppacher.		
1245	C			
1246	C	INPUT VARIABLES : CHORD - chord length		
1247	C	DROPDI - droplet diameter		
1248	C	NPOINT - no of airfoil surface points		
1249	C	VEL - free stream air speed		
1250	C	VX,VY - current droplet speeds		
1251	C	X,Y - current droplet positions		
1252	C			
1253	C			
1254	C	OUTPUT VARIABLES : DVX,DVY - current droplet accelerations		
1255	C			
1256	C			
1257	C	COMMON VARIABLES : CO,SI,GAMMA,DELTA,XCON,YCON, Y11 ,YFOIL		
1258	C			
1259	C			
1260	C	OTHER VARIABLES : COEFF - combined coefficients of the		
1261	C	equations of motion		
1262	C	DRAGCO - drag coefficient		
1263	C	DROPRE - droplet Reynolds number		
1264	C	GRIDX, GRIDY - gridpoints used in the finite		
1265	C	difference calculation of air		
1266	C	speed from the stream function		
1267	C	N - NPOINT/15, the no. of surface		
1268	C	points used to calculate the		
1269	C	stream function		
1270	C	K, R - variables for the calculation		
1271	C	of stream function		
1272	C	PSI - stream function of the air flow		
1273	C	field		
1274	C	SUMK - array containing the sum of K		
1275	C	times GAMMA(J) for 1 gridpoints		
1276	C	UX,UY - current air speed components		
1277	C			
1278	C			
1279	C			
1280	C	ROUTINE FLOWXY(X,Y,VX,VY,VEL,CHORD,DROPDI,NPOINT,		
1281	C			
1282	C			
1283	C	DOUBLE PRECISION CO(40),DELTA(40),GAMMA(41),GRIDX(4),		
1284	C	GRIDY(4),PSI(4),SI(40),SUMK(4),		
1285	C	XCON(41),YCON(41),XFoil(600).		
1286	C			
ISN	1			
ISN	2			

LEVEL 1.1.1 (DEC) 81 VS FORTRAN DATE: JUL 24. 1986 TIME: 13:15:53 NAME: FLOWXY PAGE: 29

```

1287.      +
1288.      C DOUBLE PRECISION COEFF,CHORD,DRAGCO,DROPDI,DRPRE,DWX,
1289.          + DVY,K,R,UX,UV,VEL,VX,VY,X,Y
1290.      C
1291.      C INTEGER NPOINT
1292.      C COMMON /MN/XFOIL,YFOIL,GAMMA/DRP/DELTA,CO,SI,XCON,YCON
1293.      C
1294.      C
1295.      C Set up non-dimensional grid for finite difference calculation
1296.      C of the streamfunction derivatives
1297.      C
1298.      C GRIDX(1) = X / CHORD
1299.      C GRIDX(2) = X / CHORD
1300.      C GRIDX(3) = (X - (DROPDI / 2.0)) / CHORD
1301.      C GRIDX(4) = (X + (DROPDI / 2.0)) / CHORD
1302.      C GRIDY(1) = (Y + (DROPDI / 2.0)) / CHORD
1303.      C GRIDY(2) = (Y - (DROPDI / 2.0)) / CHORD
1304.      C GRIDY(3) = Y / CHORD
1305.      C GRIDY(4) = Y / CHORD
1306.      C
1307.      C Define N as in POFLO
1308.      C
1309.      C Calculate for each grid point, R and the sum of products
1310.      C N = NPOINT / 15
1311.      C
1312.      C K times GAMMA(J), for J = 1, N
1313.      C
1314.      C
1315.      C DO 10 I = 1,4
1316.      C     R = (GRIDY(I))
1317.      C     SUMK(I) = 0.0
1318.      C     DO 20 J = 1,N
1319.      C       CALL KCALF(XCON(J),YCON(J),GRIDX(I),GRIDY(I),
1320.      C                      DELTA(J),CO(J),SI(J),K)
1321.      C       SUMK(I) = SUMK(I) + (GAMMA(J) * K)
1322.      C     CONTINUE
1323.      C
1324.      C Calculate stream functions
1325.      C
1326.      C     PSI(I) = -SUMK(I) + R - GAMMA(N+1)
1327.      C
1328.      C
1329.      C Calculate dimensional air speeds at unit x,y from finite
1330.      C difference derivatives of the stream functions
1331.      C
1332.      C     UX = VEL * CHORD * (PSI(1) - PSI(2)) / DRPDI
1333.      C     UV = VEL * CHORD * (PSI(3) - PSI(4)) / DRPDI
1334.      C
1335.      C Calculate Reynolds number and drag coefficient
1336.      C
1337.      C     DRPRE = DRPDI * DSQRT(((VX - UX) ** 2) + ((VY - UV)
1338.          + ** 2)) / 1.262E-5
1339.      C     IF (DRPRE .LT. 2.0) THEN
1340.          - DRAGCO = (24.0 / DRPRE) + (2.45 / (DRPRE ** 0.045))
1341.      C ELSE IF (DRPRE .LT. 21.0) THEN
1342.          - DRAGCO = (24.0 / DRPRE) + (2.76 / (DRPRE ** 0.198))

```

```

LEVEL 1.1.1 (DEC 81)          VS FORTRAN          DATE: JUL 24, 1986   TIME: 13:15:53   NAME: FLOWXY PAGE: 30
*.....*.....1.....2.....3.....4.....5.....6.....7.....8.....9

ISN 31    1343    ELSE IF (DROPRE .GT. 200.0) THEN
ISN 32    1344    DRAGCO = (24.0 / DROPRE) + (4.73 / (DROPRE ** 0.37))
ISN 33    1345    + (6.24E-3 * (DROPRE ** 0.38))
ISN 34    1346    ELSE
ISN 35    1347    DRAGCO = (24.0 / DROPRE) + (4.536 / (DROPRE ** 0.368))
ISN      1348    ENDIF

ISN      1349    C Calculate accelerations according to the equations of motion
ISN      1350    C
ISN      1351    COEFF = 9.9225E-4 * (DRAGCD / DROPDF)
ISN      1352    DVX = COEFF * DSQRT((VX - UX) ** 2) + ((VY - UV) ** 2)
ISN      1353    + (VX - UX)
ISN      1354    DVY = -COEFF * DSQRT((VX - UX) ** 2) + ((VY - UV) ** 2)
ISN      1355    + (VY - UV)
ISN      1356    C
ISN      1357    RETURN
ISN      1358.
ISN      1359. END

*STATISTICS* SOURCE STATEMENTS = 40, PROGRAM SIZE = 2150 BYTES, PROGRAM NAME = FLOWXY PAGE: 28.
*STATISTICS* NO DIAGNOSTICS GENERATED.
***** END OF COMPILATION 8 *****
```

LEVEL 1.1.1 (DEC 81)

VS FORTRAN

DATE: JUL 24. 1986 TIME: 13:15:53 NAME: MAIN PAGE: 31

OPTIONS IN EFFECT: MOLIST NOMREF NOGOSTMT NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(O) LANGVL(77) NOFIPS FLAG(I) NAME(MAIN) LINECOUNT(GO)

```

1360. C ****
1361. C * SUBROUTINE KCALC - calculates a single element of the
1362. C * matrix K for the control points I
1363. C *
1364. C *
1365. C *
1366. C * INPUT DATA : CO,SI - direction angles of the segment
1367. C * DELTA - half length of the control
1368. C * element J
1369. C *
1370. C * XCI,XCJ - x-coordinates of control points
1371. C * I and J, or of grid point I
1372. C *
1373. C * YCI,YCJ - y-coordinates of above
1374. C *
1375. C * OUTPUT VARIABLES :
1376. C * K - matrix element
1377. C *
1378. C * OTHER VARIABLES : A,B - geometrical quantities defined
1379. C * in Figure 2. Kennedy and Marsden
1380. C * squares of R1 and R2, also
1381. C * R1SQ,
1382. C * R2SQ - defined in Figure 2 of K&M
1383. C * TERM1, - individual terms of the
1384. C * expression for K
1385. C * TERM2,
1386. C * TERM3,
1387. C *
1388. C *
1389. C *
1390. C *
1391. C *
1392. C *
1393. C *
1394. C *
1395. C *
1396. C *
1397. C *
1398. C *
1399. C *
1400. C *
1401. C *
1402. C *
1403. C *
1404. C *
1405. C *
1406. C *
1407. C *
1408. C *
1409. C *
1410. C *
1411. C *
1412. C *

ISN 1
ISN 2
ISN 3
ISN 4
ISN 5
ISN 6
ISN 7
ISN 8
ISN 9
ISN 10
ISN 11
ISN 12
ISN 13
ISN 14
ISN 15
ISN 16

      B = ((XCI-XCJ) * CO) + ((YCI-YCJ) * SI)
      A = ((YCI-YCJ) * CO) - ((XCI-XCJ) * SI)
      R1SQ = (A*A) + ((B+DELTA)*(B+DELTA))
      R2SQ = (A*A) + ((B-DELTA)*(B-DELTA))
      TERM1 = (B + DELTA) * DLOG(R1SQ)
      TERM2 = (B - DELTA) * DLOG(R2SQ)
      DENOM = (A*A) + (B*B) - (DELTA*DELTA)
      IF (DENOM .LE. 1.00-30) THEN
        IF (DABS(A) .LT. 1.00-30) THEN
          TERM3 = 2.0 * A * PI
        ELSE
          TERM3 = 2.0 * A * (DATAN((B + DELTA) / A) -
          DATAN((B - DELTA) / A))
        ENDIF
      ENDIF
    
```

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:15:53 NAME: KCALC - PAGE: 32
..... 1 2 3 4 5 6 7 8 9

```
1413. ELSE
1414. TERM3 = 2.0 * A * DATAN(2.0 * A * DELTA * DENOM)
1415. ENDIF
1416. C   K = (TERM1 - TERM2 + TERM3 - (4.0 * DELTA)) / (4.0 * PI)
1417. C
1418. C   RETURN
1419. C   END
1420.
1421.
1422. SOURCE STATEMENTS = 22. PROGRAM SIZE = 1060 BYTES. PROGRAM NAME = KCALC PAGE: 31.
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STATISTICS NO DIAGNOSTICS GENERATED.
***** END OF COMPILATION 9 *****

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 00:15:54 NAME: MAIN PAGE: 33

OPTIONS IN EFFECT: NOLIST NOMAP NOXREF NOGOSTMT NODECK SOURCE TERM OBJECT FIXED
OPTIMIZE(0) LANG(LV(77)) NOFIPS FLAG(1) NAME(MAIN) LINECOUNT(60)

```

1421. C
1422. C
1423. C * SUBROUTINE COLLID - Calculates the x,y position
1424. C * of impact on the airfoil surface
1425. C
1426. C
1427. C * INPUT VARIABLES: XOUT, - last droplet position outside
1428. C * YOUT, - the airfoil
1429. C * XIN, - first droplet position inside
1430. C * YIN, - the airfoil
1431. C * XF1,XF2 - x and y coordinates of the end-
1432. C * points of the surface segment
1433. C
1434. C * OUTPUT VARIABLES: FINALX, - estimated coordinates of the
1435. C * FINALY, - impact point on the surface
1436. C
1437. C
1438. C * OTHER VARIABLES: SLOPTR - slope of the trajectory segment
1439. C * which crosses the airfoil surface
1440. C * SLOPSF - slope of the surface segment
1441. C * Intercepted by the trajectory
1442. C * INT1, - intercepts for the linear
1443. C * INT2, - equations of the two segments
1444. C
1445. C
1446. C
1447. C
1448. C
1449. C * SUBROUTINE COLLID (XOUT,XIN,YOUT,YIN,XF1,YF1,XF2,YF2,
1450. C * FINALX,FINALY)
1451. C
1452. C * DOUBLE PRECISION XOUT,XIN,YOUT,YIN,XF1,XF2,YF1,YF2,
1453. C * SLOPTR,SLOPSF,INT1,INT2,FINALX,FINALY
1454. C
1455. C Calculate slope of trajectory at impact point
1456. C
1457. C * SLOPTR = (YIN - YOUT) / (XIN - XOUT)
1458. C
1459. C Calculate slope of surface segment at impact point
1460. C
1461. C * SLOPSF = (YF2 - YF1) / (XF2 - XF1)
1462. C
1463. C Find intercepts of the two linear equations
1464. C
1465. C * INT1 = YIN - (SLOPTR * XIN)
1466. C * INT2 = YF2 - (SLOPSF * XF2)
1467. C
1468. C Find the impact point as the intersection point
1469. C of the two segments
1470. C
1471. C * FINALX = (INT2 - INT1) / (SLOPTR - SLOPSF)
1472. C * FINALY = (SLOPTR * FINALX) + INT1
1473. C

```

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24. 1986 TIME: 13:15:54 NAME: C0LLID PAGE: 34
ISN 9 1474 RETURN
ISN 10 1475 END
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***** END OF COMPIRATION 10 *****

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:15:54 NAME: MAIN PAGE: 35

SUMMARY OF MESSAGES AND STATISTICS FOR ALL COMPILENTIONS

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***** END OF COMPILATION 2 *****

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***** END OF COMPILATION 3 *****

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***** END OF COMPILATION 4 *****

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***** END OF COMPILATION 5 *****

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***** END OF COMPILATION 6 *****

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***** END OF COMPILATION 7 *****

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***** END OF COMPILATION 8 *****

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***** END OF COMPILATION 9 *****

STATISTICS SOURCE STATEMENTS = 10. PROGRAM SIZE = 512 BYTES. PROGRAM NAME = C0LLID PAGE: 33

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NAME : PAGE :

NAME : MAIN

LEVEL 1.1.1 (DEC 81) VS FORTRAN DATE: JUL 24, 1986 TIME: 13:15:54

STATISTICS NO DIAGNOSTICS GENERATED.

***** END OF COMPIRATION 10 *****

***** SUMMARY STATISTICS ***** O DIAGNOSTICS GENERATED. HIGHEST SEVERITY CODE IS 0.

Appendix VI

The MVD Approximation for Collision Efficiency

The task of computing volume-weighted averages of collision efficiencies for a droplet size spectrum is often simplified by substituting a monodisperse size spectrum at the median volume diameter of the distribution. As discussed in Chapter III, this can lead, in some cases, to a significant underestimation of the maximum accretion angle. However, it has often been assumed that the weighted overall and stagnation line collision efficiencies are well approximated by the mvd values. In this Appendix the validity of that assumption is examined.

Table A.1 lists a variety of atmospheric and marine droplet size spectra, as relative numbers of droplets in 10 or 3 μm wide size bins. These spectra were either measured in the field, or in icing wind tunnels, or estimated from empirical parameterized distributions, and cover a range of types and mvds.

They are derived from the following sources: Bain and Gayet (1984), Battan and Reitan (1957), Best (1951), Khrgian-Mazin, as reported by Pruppacher and Klett (1980), McComber and Touzot (1981), Preobrazhenskii (1973), and Squires (1958). Those marked FROST were measured by the oil

mid-band droplet diameter	relative number of drops per unit volume					requires Trade-wind q	
	Squires 1	Squires 2	Squires 3	Squires 4	Squires dark stratus		
5 um	0.724	0.069	0.014	0.002	0.25	6.5	9
15	2.397	0.340	0.082	0.012	0.65	6.0	16
25	2.449	0.518	0.148	0.024	1.0	5.4	29
35	1.766	0.557	0.189	0.035	0.8	5.1	21
45	1.074	0.505	0.203	0.043	0.8	2.5	3.5
55	0.590	0.414	0.198	0.047	0.7	0.5	
65	0.303	0.317	0.180	0.049	0.35	0.2	
75	0.149	0.232	0.156	0.048	0.15	0.1	
85	0.070	0.164	0.131	0.046	0.10	0.05	
95	0.032	0.112	0.106	0.042	0.06		
105	0.014	0.075	0.085	0.038	0.05		
115	0.006	0.049	0.066	0.034	0.04		
125	0.003	0.032	0.051	0.030	0.03		
135	0.001	0.021	0.039	0.026	0.04		
145		0.013	0.029	0.022	0.01		
155		0.008	0.022	0.019			
165		0.005	0.016	0.016			
175		0.003	0.012	0.013			
185		0.002	0.009	0.011			
195		0.001	0.006	0.009			
205			0.004	0.007			
215			0.003	0.006			
225				0.005			
235				0.004			
245				0.003			

Table A.1 - Droplet size spectra from many different sources
(see text for references). -

relative number of drops per unit volume

mid-band droplet diameter	Preobrazhenskii						
	FROST 5	FROST 4	FROST 3	FROST 2	FROST 1	Battan and Reitan stratus	Battan and Reitan fairweather cumulus
5 μ m	600	100	127	94	97	3	17
15	70	15	327	167	213	37	68
25	13	1.1	104	33	86	63	64
35	0.26	0.2	13	6	33	58	31
45				4	14	51	12
55					1	35	11
65						26	9
75						17	3
85						14	2
95						16	4
105						12	9
115						5	2
125						4	3
135						8	2
145						7	1
155						2	2
165						1	2
175						2	1
185							2

Table A.1 - continued.

relative number of drops per unit volume

mid-band droplet diameter	McComber and Touzot	mid-band droplet diameter	Bain and Gayet 3	Bain and Gayet 2	Bain and Gayet 1
5.1 μm	945	1.5	20	7	13.0
15.3	1093	4.0	41	12	15.5
25.5	481	7.5	30	23	15.5
35.7	211	10.5	7.5	33	18.0
45.9	74	13.5	2.5	18	15.5
56.1	38	16.5		5	12.5
66.3	30	19.5		2	6.0
76.5	13	22.5			2
86.7	19	25.5			2
96.9	11				
107.1	10				
117.3	3				
127.5	1				
137.7	1				

percent volume

mid-band droplet diameter	McComber and Touzot	Best 1	Best 2	Best 3	Best 4	Best 5
5 μm	63.2	22	4	3	36	
15	36.8	41	24	31	47	
25		26	39	53	15	
35		9	26	13	2	
45		2	6	1		
55						

Table A.1 - continued.

slide technique in the FROST icing wind-tunnel facility at The University of Alberta.

For each spectrum, the mvd was calculated following the method derived by Lozowski (1978). Assuming a uniform distribution of droplets within each bin, the partial volume of each is given by:

$$v_i = (n_i/24) \pi w^3 (i^4 - (i-1)^4), \quad A.8$$

where n_i is the number of drops in the i th bin, and w is the bin width, say 10 μm . Then, if V is the total volume of all droplets, let

$$u_i = (\Sigma v_i) / V. \quad A.9$$

Now find u_k such that $u_{k-1} < 0.5$, and $u_k > 0.5$. Finally,

$$\begin{aligned} \text{mvd} = w & ((0.5 - u_{k-1}) / (u_k - u_{k-1})) (k^4 - (k-1)^4) \\ & + (k-1)^4)^{0.25}, \end{aligned} \quad A.10$$

Overall and stagnation line collision efficiencies have been calculated for the mvd of each spectrum. For the complete spectrum, weighted average collision efficiencies are calculated using the fractional volume of each bin as the weighting factor for the collision efficiency of the mvd of that bin.

All of the collision efficiencies are calculated according to Table II.6, and for the same conditions, i.e. $U = 10 \text{ m/sec}$ and $D_c = 0.0254 \text{ m}$. The results are shown in Table A.2 and Figures A.2 and A.3. For a different type of comparison, the conditions have been varied for a single spectrum (labelled as Bain and Gayet 2 in Table A.1) to give a complete range of collision efficiencies. These results are compared in Table A.3 and in Figures A.4 and A.5.

Also calculated are collision efficiencies for the droplet at the 42nd volume percentile (d_{42}), on a suggestion made by Lozowski and Makkonen (private communication, 1986), based on a theoretical analysis using Gaussian quadrature to estimate the overall collision efficiency for a droplet spectrum. These data are included in [redacted] figures and tables listed above.

Comparison of Figure A.4 with Figure 1 of Makkonen (1984), which shows the same type of calculations for the same spectrum used here as well as two other spectra given by Bain and Gayet, 1982), reveals a major disagreement in the sign of $E_{\text{spec.}} - E_{\text{mvd}}$. The disagreement is due to a misinterpretation of the original droplet spectrum diagram in Bain and Gayet (1982) on the part of Makkonen (private communication, 1986). Figure A.4 reflects the correct interpretation.

Spectrum	β_{spec}	E_{spec}	mvd	β_{mvd}	E_{mvd}	d42	β_{d42}	E_{d42}
Khrigian-Mazin 1	0.824	0.710	56.9 μm	0.852	0.745	52.1 μm	0.835	0.718
Khrigian-Mazin 2	0.906	0.836	93.6	0.924	0.862	86.1	0.914	0.846
Khrigian-Mazin 3	0.936	0.886	126.2	0.949	0.906	116.5	0.943	0.895
Khrigian-Mazin 4	0.955	0.919	165.9	0.964	0.933	154.5	0.960	0.927
Squires Hawaiian	0.886	0.804	81.4	0.908	0.835	70.9	0.889	0.804
Squires dark stratus	0.767	0.626	43.0	0.790	0.650	39.7	0.769	0.619
Squires Trade-wind Cu	0.702	0.532	34.9	0.730	0.565	33.2	0.714	0.543
Battan and Reitan Cu	0.393	0.224	18.0	0.459	0.249	16.6	0.419	0.211
Battan and Reitan stratus	0.364	0.199	17.2	0.434	0.225	15.9	0.396	0.191
FROST 1	0.555	0.362	24.7	0.603	0.402	22.6	0.565	0.357
FROST 2	0.583	0.402	26.2	0.628	0.432	23.4	0.580	0.375
FROST 3	0.680	0.516	35.0	0.731	0.566	31.2	0.694	0.515
FROST 4	0.928	0.873	127.7	0.949	0.907	111.3	0.939	0.889
FROST 5	0.919	0.859	112.3	0.940	0.890	107.1	0.936	0.883
Preobrazhenskii	0.838	0.745	78.5	0.903	0.827	67.9	0.883	0.793
McComber and Touzot	0.836	0.736	75.9	0.899	0.820	65.8	0.878	0.786
Bain and Gayet 1	0.108	0.036	8.8	0.088	0.010	8.4	0.065	0.003
Bain and Gayet 2	0.284	0.123	13.3	0.301	0.118	12.5	0.267	0.096
Bain and Gayet 3	0.421	0.232	17.5	0.444	0.233	16.6	0.418	0.210
Best 1	0.415	0.247	18.3	0.467	0.255	17.0	0.429	0.220
Best 2	0.582	0.401	26.9	0.639	0.445	25.0	0.610	0.409
Best 3	0.542	0.350	24.4	0.599	0.397	22.5	0.564	0.357
Best 4	0.313	0.160	15.3	0.375	0.173	13.1	0.291	0.111
Best 5	0.186	0.070	9.4	0.120	0.021	9.0	0.098	0.014
standard error			0.031	0.029	0.027	0.026		

Table A.2 | Overall and stagnation line collision efficiencies as plotted in Figures A.2 and A.3.

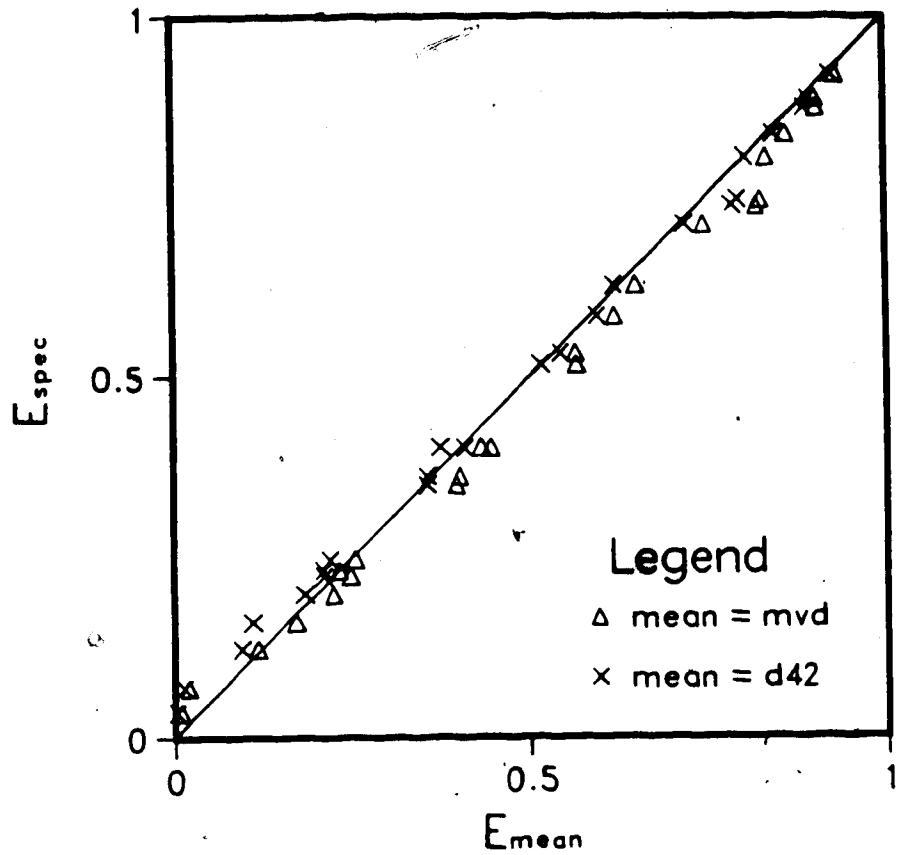


Figure A.2 - Overall collision efficiencies calculated for many different drop size spectra vs. E's calculated for the mvd and d42 of each spectrum. In each case U=10 m/sec, D_c=0.0254 m.

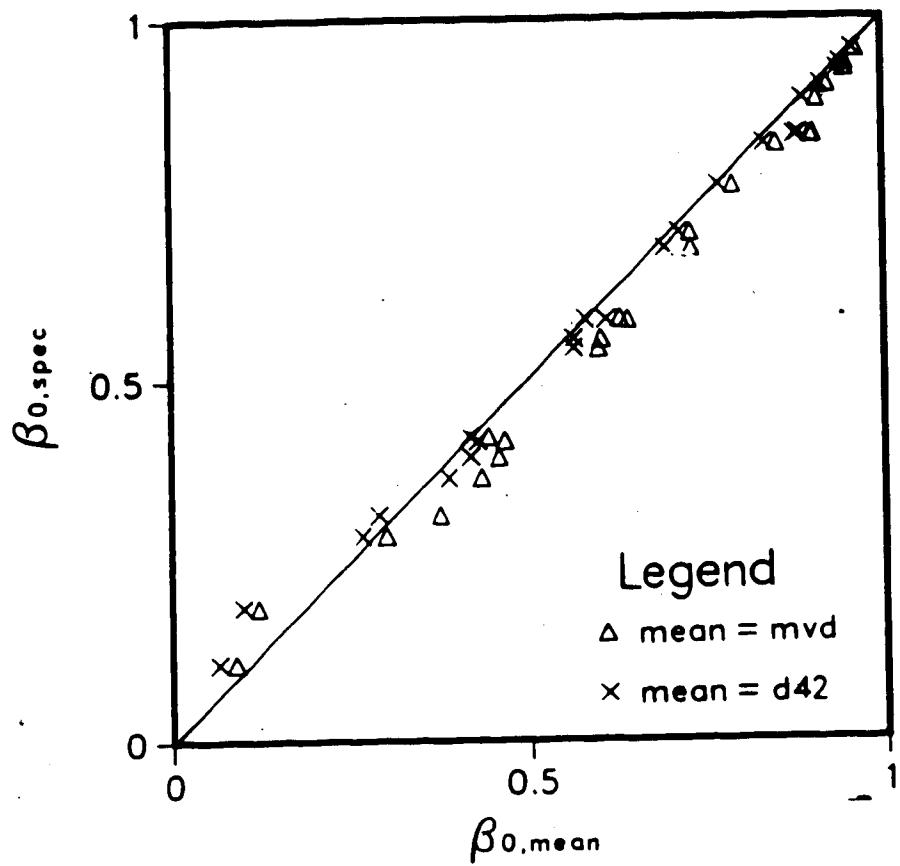


Figure A.3 - Stagnation line collision efficiencies calculated for many different drop size spectra vs. β_0 's calculated for the mvd and d42 of each spectrum. In each case $U=10$ m/sec, $D_c=0.0254$ m.

U	$\beta_{o\text{ spec}}$	E_{spec}	$\beta_{o\text{mvd}}$	E_{mvd}	$\beta_{o\text{d42}}$	E_{d42}
5m/sec	0.121	0.037	0.121	0.022	0.089	0.010
7	0.202	0.072	0.208	0.062	0.175	0.044
10	0.284	0.123	0.301	0.118	0.267	0.096
15	0.383	0.194	0.403	0.196	0.370	0.169
20	0.448	0.250	0.470	0.257	0.439	0.228
25	0.496	0.295	0.518	0.305	0.489	0.275
30	0.532	0.331	0.555	0.344	0.527	0.314
40	0.586	0.389	0.608	0.405	0.582	0.374
50	0.624	0.432	0.646	0.450	0.622	0.420
80	0.693	0.516	0.713	0.536	0.692	0.508
200	0.790	0.648	0.806	0.668	0.791	0.646
350	0.830	0.708	0.844	0.726	0.832	0.708
700	0.865	0.765	0.877	0.781	0.867	0.766
1500	0.892	0.811	0.901	0.825	0.894	0.812
standard error			0.012	0.013	0.011	0.011

Table A.3 - Overall and stagnation line collision efficiencies as plotted in Figures A.4 and A.5.

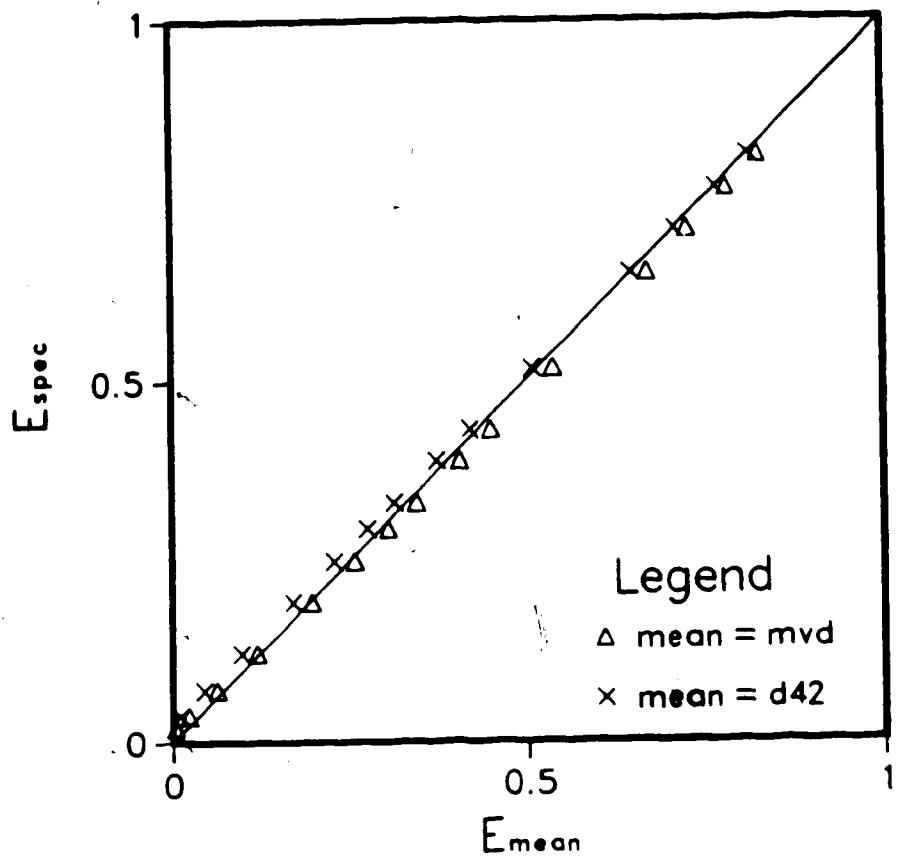


Figure A.4 - Overall collision efficiencies calculated for a single drop size spectrum (Bain and Gayet 2) at many different wind speeds, vs. E's calculated for the mvd and d42 of the same spectrum for the same wind speeds. $D_c = 0.0254$ m.

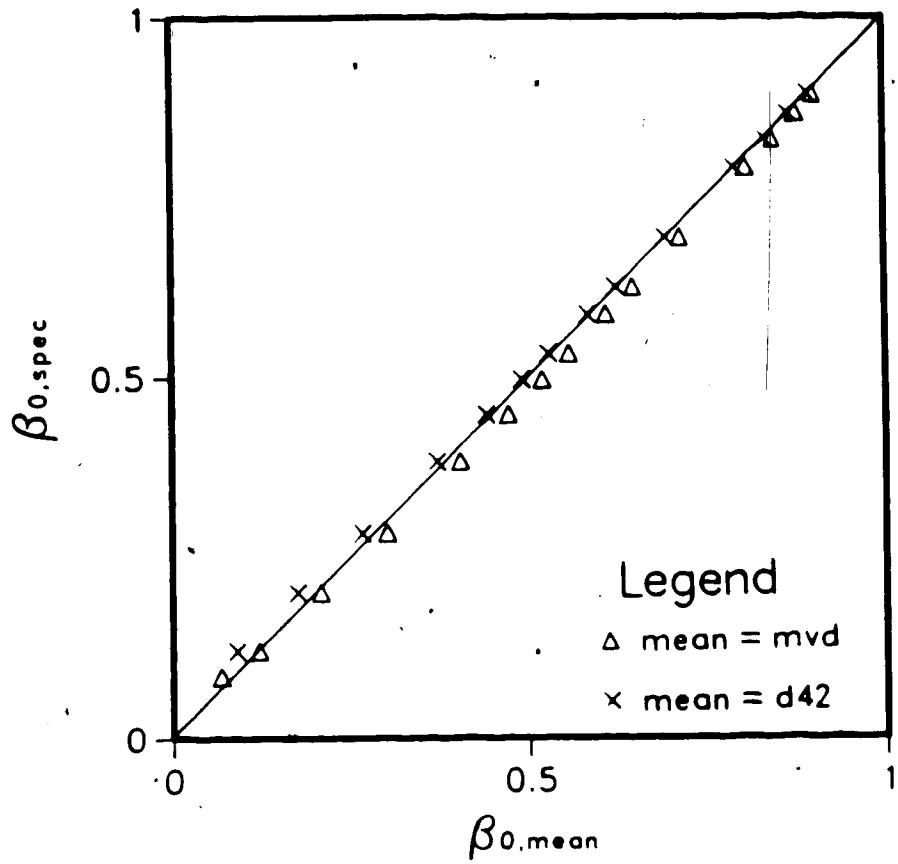


Figure A.5 - Stagnation line collision efficiencies calculated for a single drop size spectrum (Bain and Gayet 2) at many different wind speeds, vs. β_0 's calculated for the mvd and d42 of the same spectrum for the same wind speeds. $D_c = 0.0254$ m.

The conclusion which may be made from this exercise is that, although the standard errors are not significantly different over the entire range of collision efficiencies (see Tables A.2 and A.3), in the range of E or β_0 less than about 0.2 the mvd approximation is preferable. Above that range, for most types of spectra, the d42 approximation gives better agreement with the spectrum-weighted collision efficiencies. However, for certain types of spectra which depart significantly from a Gaussian distribution (e.g. the double peaked spectrum from McComber and Touzot) even the d42 value may be in error by more than 10 percent.

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