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# Reply by Authors to J L Potter and J D Whitfield

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#### Nomenclature

vertical height of roughness above plate, ft k

heat flow rate along the centerline of the model

 $\stackrel{\scriptscriptstyle 1}{R}_0$ freestream unit Reynolds number per ft

Reynolds number based on fluid conditions at top of the  $R_k$ roughness elements and the height of the roughness

roughness Reynolds number for which a further increase  $R_k$ in roughness height causes no appreciable forward movement of the beginning of fully developed turbulent flow

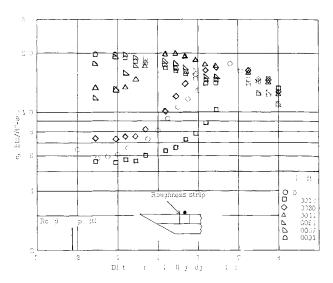
distance from the leading edge x

distance from the leading edge to the roughness position  $x_k$ 

THE salient points mentioned in the preceding comments were not discussed in Ref 1 because of a desire to make that manuscript as brief as possible However, these matters (including references of the preceding comment) are discussed in a more detailed report,2 which was under preparation at the time the original manuscript was submitted The authors certainly feel that any full length paper on transition should include a reference to the work of Potter and Whitfield

The discrepancy between the Reynolds numbers for natural transition in Figs 1 and 2 of Ref 1 is thought to result not only from small variations in the leading-edge thickness, but also from a small angle-of-attack variation between the two assemblies However, the angle of attack for each series of roughness tests was invariant The interchangeable leadingedge section method of varying the roughness elements has some inherent disadvantages since leading-edge thickness is a factor in determining the location of transition this method was chosen as a practical means of reducing the required time for a model change

Since this early work indicated that small roughness heights may delay transition, further research on this phenomenon is (This work has been further stimulated by being conducted the comments of Potter and Whitfield) Some new data now available have indicated that whereas some of the variation in the location of transition reported in Ref 1 was due to variation of the leading-edge thickness, under certain conditions transition is apparently slightly delayed when the surface roughness is less than the boundary layer thickness Figure 1 presents the heat flow rate distribution for the model with various height roughness elements at a unit Reynolds number of approximately  $8 \times 10^6/\mathrm{ft}$  and a Mach number of 6Test conditions are similar to those given in Ref 1 matic of the model shown in the figure illustrates the new



Heating rate distribution along flat plate with various size roughness,  $R_0 = 8.3 \times 10^6$ 

mounting technique with the same leading edge (thickness diameter of 0 0025 in ) being used for all tests and the small roughness strips being interchangeable This figure illustrates both the delay in transition obtained with the smallest roughness elements and the critical roughness height required for these freestream conditions (critical height is taken as 0 0091 in the figure) Figure 2 shows the variation of the critical roughness Reynolds number with freestream unit Reynolds number Included in this figure are the data from Fig 3 of Ref 1 and new data taken with one leading edge with roughness elements mounted on an interchangeable Although the new and old data do not coincide, which may be due in part to a different location  $(x_k)$  of the roughness elements, these data do not change the conclusions given A study of the effect of small roughness heights on transition is being continued

The definition and determination of the critical roughness Reynolds number varies considerably in the literature Difficulties arising from this were recognized, and as a result the roughness data were also compared in Ref 2 to other transition roughness data detected by a method sensitive to permanent changes in the boundary layer This comparison did not change the conclusions given in Ref 1

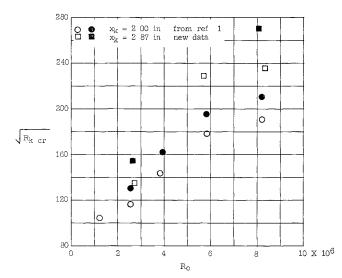


Fig 2 Variation of critical roughness Reynolds number with freestream Reynolds number Open symbols indicate that the roughness height is slightly less than the critical value, and solid symbols indicate that the roughness height is slightly greater than the critical value

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<sup>1</sup> Sterrett, J R and Holloway, P F, "Effects of controlled roughness on boundary-layer transition at a Mach number of 6.0." ATAA I 1 1951-1953 (1963)

 $6\,0$  ' AIAA J 1, 1951–1953 (1963)  $^2$  Holloway, P F and Sterrett, J R , "Effects of controlled surface roughness on boundary layer transition and heat transfer at Mach numbers of 4 8 and 6 0 " Proposed NASA TN D-2054

# Comment on "Interception of High-Speed Target by Beam Rider Missile"

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 $\blacksquare$  N Ref 1, it is possible to express the missile coordinates explicitly in terms of the normal elliptic integrals F and F

The linear differential equation considered by the authors was

$$(d/d\gamma)(\cot\theta) + \frac{1}{2}\cot\gamma\cot\theta + \frac{1}{2} = 0 \tag{1}$$

Let  $ds_m$  be an element of arc of the missile trajectory and

$$ds_m = V_m dt = (V_m/V_t)V_t dt = \tau R d(\cot \theta)$$

With  $\tau R = c$  and cot  $\theta = x_m/y_m$ , we have,

$$ds_m = cd\left(\frac{x_m}{y_m}\right) = c\left(\frac{dx_m}{y_m} - \frac{x_m dy_m}{{y_m}^2}\right)$$
 (2)

Since

$$dx_m = \cos \gamma ds_m \qquad dy_m = \sin \gamma \ ds_m$$

Eq (2) becomes

$$ds_m = c \left( \frac{\cos \gamma}{y_m} - \frac{x_m \sin \gamma}{y_m^2} \right) ds_m \tag{3}$$

Therefore,

$$y_m^2 = c(y_m \cos \gamma - x_m \sin \gamma) \tag{4}$$

This equation clearly shows that the velocity of the missile is tangent to the circle centered at the origin and of radius  $y_m^2/c$ , a result mentioned in Ref  $\, 2$  using different arguments Therefore,

$$\cot\theta = \frac{x_m}{y_m} = \cot\gamma - \frac{y_m}{c\sin\gamma} \tag{5}$$

Using Eq. (5) in (1), one obtains the differential equation

$$2\sin\gamma(dy/d\gamma) - \cos\gamma y + 1 = 0 \tag{6}$$

which is also linear with  $y = y_m/c$ 

Integrating Eq. (6) yields

$$y = \cos\phi \left( C + \frac{F(\phi, k) - 2E(\phi, k)}{2^{1/2}} \right) + \sin\phi (1 + \cos^2\phi)^{1/2}$$
 (7)

and using (4) we obtain the x coordinate as

$$x = -\left(C + \frac{F - 2E}{2^{1/2}}\right) \times \left[C + \frac{F - 2E}{2^{1/2}} + \tan\phi(1 + \cos^2\phi)^{1/2}\right]$$
(8)

where F and E are normal elliptic integrals of the first and

second kind with moduli  $k=1/2^{1/2}$  and arguments  $\phi=\arccos(\sin\gamma)^{1/2}$ 

The constant C is determined by the initial conditions We also have

$$\cot\theta = \frac{x}{y} = -\frac{1}{\cos\phi} \left( C + \frac{F - 2E}{2^{1/2}} \right) \tag{9}$$

which is Eq (9) in Ref 1

Using the initial conditions, when x = y = 0,  $\theta = \theta_0 = \gamma_0$ , we have for the constant C,

$$C = \frac{2E(\phi_0) - F(\phi_0)}{2^{1/2}} - \tan\phi_0 (1 + \cos^2\phi_0)^{1/2}$$
 (10)

By these expressions it can easily be seen that:

1) y=1 is an asymptote since x becomes infinite for  $\gamma=0$  except when

$$C\,-\,\frac{2E(\pi/2)\,-\,F(\pi/2)}{2^{1/2}}\,=\,0$$

or C = 0.599 This gives  $\theta_0 = 37^{\circ}$ 

2) The points where the velocity is directed vertically are such that  $\gamma = \pi/2$ ,  $\phi = 0$  Therefore  $x = -C^2$  and y = C

Hence, they are situated on the parabola  $y^2 = -x$ , a result also mentioned in Ref. 2 using different arguments

#### References

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## "Equilibrium" Gas Composition Computation with Constraints

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IT has been suggested in a recent note<sup>1</sup> that one may "freeze" a particular species in a chemical equilibrium computation by introducing a fictitious second set of species and manipulation of the real and imaginary sets of species, within the usual procedures of the computation routine. The purpose of this note is to point out that such a constraint may be imposed on the equilibrium calculation in another manner when one is using the popular minimization of free energy technique <sup>2</sup> In this approach the Gibbs free energy is minimized subject to the constraints of the mass balances

$$\sum_{i=1}^{n} a_{ji} x_i = b_j \tag{1}$$

where  $a_{ji}$  is the number of atoms of element j in species i,  $x_i$  is the moles of species i per unit mass, and  $b_j$  is the number of gram-atoms of element j per unit mass

It is possible, however, to impose a priori relations (constraints) among the  $x_i$  composition variables by considering the constraints to be pseudo elements, as long as the relations are of the form of Eq. (1), i.e., linear. Thus the formula matrix  $a_{ji}$  would have a column for each species and m rows for each of the m real chemical elements, as usual, with an additional row for each constraint. Possible constraints

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