

## University of Southampton Research Repository ePrints Soton

Copyright © and Moral Rights for this thesis are retained by the author and/or other copyright owners. A copy can be downloaded for personal non-commercial research or study, without prior permission or charge. This thesis cannot be reproduced or quoted extensively from without first obtaining permission in writing from the copyright holder/s. The content must not be changed in any way or sold commercially in any format or medium without the formal permission of the copyright holders.

When referring to this work, full bibliographic details including the author, title, awarding institution and date of the thesis must be given e.g.

AUTHOR (year of submission) "Full thesis title", University of Southampton, name of the University School or Department, PhD Thesis, pagination

## Appendix E

# FORTRAN 95 Codes

### VIX Head Files and Definitions

```
parameter(NQA = 300)
type node
  integer id
  real:: x,y,z
  real:: s,t,TAU
  real:: Up,U,Cp,Cpe
  real:: Dxt2,Dyt2,Dzt2
  real:: Dxs2,Dys2,Dzs2
  real:: Dps2,Dpt2
end type node

type section
  integer:: id,Nnd,Nlower
  integer:: istag,iblte(2),Nbl(2)
  integer,dimension(NQA,2):: IBL
  integer,dimension(2):: imatch,Nmatch,tr
  real:: sc,disp,leng,Vchord(3),aflow
  real,dimension(2):: xsep,xre,xtr
  real,dimension(NQA):: s,t,curv
  real,dimension(3,NQA):: P,Vn
!   BOUNDARY LAYER VARS
  real,dimension(NQA,2):: TAU,UINV
  real,dimension(NQA):: Up,Cp
  real,dimension(4,NQA,2):: Q
  real:: cli,clv,cd,Cdf,Cdp
end type section

!
!-----
!   NODE
!   id      -> pt identity
!   x,y,z   -> coordinates of node
!   Cp      -> pressure coefficient
!   Cpe     -> viscous pressure coefficient
!   U       -> absolute velocity
!   TAU     -> shear stress
!   s,t     -> surface coordinates
!   Dxt2,Dyt2,Dzt2 -> spline 2nd derivatives on parametric
!                   t direction for each coordinate
!   Dxs2,Dys2,Dzs2 -> spline 2nd derivatives on parametric
!                   s direction for each coordinate
!   Dps2,Dpt2  -> spline second derivative on parametric
!                   directions s and t for nodal Cp
!
!-----
!   SECTION
!   id      -> identity of section
!   istag   -> location of stagnation point
!   Nlower  -> index where leading edge is
!   Npt     -> No of points on section (max=100)
!   ite     -> index where trailing edge lives
!   p(..)   -> coordinates points
!   s(.)    -> curve coordinate
!   t(.)    -> spanwise curve coordinate
!   Vn(..)  -> normal vector on a section node
!   imatch(.) -> point of beginning of matching surface
```

```

!      Nmatch(.) -> point where ends matching surface
!      tr(.)      -> point where transition occurs
!      disp       -> displacement of section on x axis
!      sc         -> scale of section
!      leng       -> chord length
!      aflow      -> flow incidence on the particular section
!      Vchord     -> unitary vector of chord
!      xsep,ysep  -> point of separation  1 - upper
!                                           2 - lower
!      xtr,ytr    -> point of transition
!      BOUNDARY LAYER
!      Q          -> vector of boundary layer variables
!                  1 - nr or Ct
!                  2 - th
!                  3 - delts
!                  4 - Ue
!      Cli        -> inviscid lift coefficient
!      Clv        -> viscous lift coefficient
!      Cd         -> drag coefficient
!      Ue         -> edge velocity modulus (viscous)
!      Up         -> edge velocity modulus (potential)
!*****
!      Program created by Augusto Veiga (University of Southampton 2003) *
!*****

      type panel
         integer:: id,np,ibc
         integer,dimension(4)::ngb
         character(5):: tipo
         real:: s,t,area,fl,fd
         type (node),dimension(4):: nd,mid
         type (node):: co
         real:: Mx,My,Mz
         real:: cp,Up,rvp
         real:: u(3),vm(3)
      end type panel

!
!      PANEL
!      id         -> pt identity
!      ibc        -> panel type index:
!                  -1 trailing edge panel
!                  1  body panel
!                  2  wake attached to trailing edge
!                  3  free wake panels
!                  4  fixed wake
!      ngb        -> neighbouring panels      (they are 4)
!                  ngb(i)= -4 reflection plane
!                  ngb(i)= -2 discontinuous
!
!      x,y,z      -> coordinates of node
!      Cp         -> pressure coefficient
!      U          -> absolute velocity
!      co         -> collocation pt of panel
!      mid        -> mid edge between nodes (follows right hand rule)
!      vm         -> vector velocity in m/s (not that useful)
!      nd         -> panel nodes (they are 4)
!      area       -> area of panel
!      fl         -> lifting force of panel
!      fd         -> drag force of panel
!      rpv        -> viscous pressure resistance
!      Mx,My,Mz   -> momentum in relation to origin of root chord
!
!*****
!      Program created by Augusto Veiga (University of Southampton 2003) *
!*****

```

## Two-Dimensional Panel Method

```

      subroutine panel (X,Y,IFLAG,Nlower
& ,Nupper,Nodtot,Nmax,V,alpha,

```

```

& naca,tau)

integer Nmax
real,dimension(Nmax)::x,y,xmid,ymid,costhe,sinthe
real,dimension(Nmax)::V,CP

real:: alpha,Tau,gamma,CD,CL1,CL2
integer:: Nodtot,Nupper,Nlower,IFLAG

real PI, PI2INV, XNU,cosalf,sinalf,thick,camber,beta
real, dimension(Nodtot,Nodtot+1) :: A
integer N

C..... Begin
C
    PI = 4. * atan(1.)
    PI2INV = 1. / (2. * PI)
    V = 0
    IF (ALPHA.gt.PI/2.) GOTO 400
!    if (ALPHA.gt.90.) goto 400
!    close (5)
C
C
C..... Initializing data
C
    print *, '*** Inicializacao dos dados - aguarde ... '
    call SETUP(x,y,xmid,ymid,costhe,sinthe,cosalf,sinalf,PI,
& PI2INV,Nodtot,Nmax,Nupper,Nlower,alpha,tau,
& NACA,IFLAG,XNU)
C
    COSALF = cos(ALPHA)    !*PI/180.)
    SINALF = sin(ALPHA)    !*PI/180.)

    A=0
C
C
C..... Influence coefficient matrix assembly
C
    print *, '*** Montagem da matriz de coeficientes - aguarde ... '
    call COFISH(x,y,xmid,ymid,costhe,sinthe,Nodtot,pi,pi2inv,
& alpha,cosalf,sinalf,Nmax,A)
C
C..... Gauss Elimination solution system
C
    print *, '*** Solucao do sistema de equacoes - aguarde ... '
    call GAUSS2(A,Nodtot-1,1,Nodtot)
C
C
C..... Velocity and pressure coefficient
C
    print *, '*** Calculo das velocidades e pressoes - aguarde ... '
    call VELDIS(A,x,y,Nlower,Nupper,Nodtot,xmid,ymid,V,CP,gamma,
& Nmax,XNU)
C
C..... Calculo dos coeficientes do aerofolio
C
    print *, '*** Calculo dos coefs. adimensionais - aguarde ... '
    call FANDM(A,x,y,Nlower,Nupper,Nodtot,xmid,ymid,V,CP,gamma,
& Nmax,sinalf,cosalf,CD,CL1,CL2)

    400 write (*,9999)
        !stop
    9999 format (//, ' End of panel method - Univ. Southampton/COPPE(C)
& 2001')

C
C----- Fim do programa principal
C
    end subroutine

```





[illegible]

```

        F1 = FI(K-1)
        S2 = 0.
        F2 = FI(K)
        S3 = sqrt((XK3 - XK2)**2 + (YK3 - YK2)**2)
        F3 = FI(K+1)
    endif
    DELTA = (S3 - S1) * (S2 - S1) * (S2 - S3)
    DELTB = (S2**2 - S1**2) * (F3 - F1)
    *      - (S3**2 - S1**2) * (F2 - F1)
    V(K) = DELTB / DELTA      !- DELTB / DELTA
    CP(K) = 1. - V(K) * V(K)
    !      CF(K) = 0.075/(log(V(K)/XNU) -2.0)**2
100 continue
c
c
c..... upper surface
c
do 110 K=Nlower,NUPPER-1
    L = k      !      NLOWER
    if (K.eq.Nlower.or.K.eq.NUPPER-1) then
        if (K.eq.Nlower) then
            XK1 = .5 * (X(L) + X(L+1))
            YK1 = .5 * (Y(L) + Y(L+1))
            XK2 = .5 * (X(L+1) + X(L+2))
            YK2 = .5 * (Y(L+1) + Y(L+2))
            XK3 = .5 * (X(L+2) + X(L+3)) ! x e y com o n° de nós
            YK3 = .5 * (Y(L+2) + Y(L+3))
            S1 = 0.
            F1 = FI(L)      !Fi varia com o n° de painéis
            S2 = sqrt((XK2 - XK1)**2 + (YK2 - YK1)**2)
            F2 = FI(L+1)
            S3 = S2 + sqrt((XK3 - XK2)**2 + (YK3 - YK2)**2)
            F3 = FI(L+2)
        endif
        if (K.eq.NUPPER-1) then
            XK1 = .5 * (X(L-2) + X(L-1))
            YK1 = .5 * (Y(L-2) + Y(L-1))
            XK2 = .5 * (X(L-1) + X(L))
            YK2 = .5 * (Y(L-1) + Y(L))
            XK3 = .5 * (X(L) + X(L+1))
            YK3 = .5 * (Y(L) + Y(L+1))
            S3 = 0.
            F3 = FI(L)
            S2 = - sqrt((XK3 - XK2)**2 + (YK3 - YK2)**2)
            F2 = FI(L-1)
            S1 = S2 - sqrt((XK2 - XK1)**2 + (YK2 - YK1)**2)
            F1 = FI(L-2)
        endif
    else
        XK1 = .5 * (X(L-1) + X(L))
        YK1 = .5 * (Y(L-1) + Y(L))
        XK2 = .5 * (X(L) + X(L+1))
        YK2 = .5 * (Y(L) + Y(L+1))
        XK3 = .5 * (X(L+1) + X(L+2))
        YK3 = .5 * (Y(L+1) + Y(L+2))
        S1 = - sqrt((XK2 - XK1)**2 + (YK2 - YK1)**2)
        F1 = FI(L-1)
        S2 = 0.
        F2 = FI(L)
        S3 = sqrt((XK3 - XK2)**2 + (YK3 - YK2)**2)
        F3 = FI(L+1)
    endif
    DELTA = (S3 - S1) * (S2 - S1) * (S2 - S3)
    DELTB = (S2**2 - S1**2) * (F3 - F1)
    *      - (S3**2 - S1**2) * (F2 - F1)
    V(L) = DELTB / DELTA
    CP(L) = 1. - V(L) * V(L)
    !      CF(L) = 0.075/(log(V(L)/XNU) -2.0)**2
110 continue
c
    return
c

```





```

150 CAMBER = EPSMAX * (1. - Z)
    DCAMDX = - EPSMAX
    goto 120
c
c
c----- End NACA45
c
    end

```

## VIX 3D Main Code

```

Program VII3d
!use AVDef
!use DFLib

!      This program reads the streamlines given by PALISUPAN
!      as object str where str has the properties
!      id          -> identity of streamline
!      Npt         -> No of points on streamline (max=100)
!      p(..)       -> coordinates of points
!      u(..)       -> stream velocity on each point
!      V(..)       -> vector of velocities
!      imatch(..)  -> point of beginning of matching surface
!      Nmatch(..)  -> point where ends matching surface
!      tr(..)      -> point where transition occurs
!      Nsec        -> number of sections (z cte)

!      flags       -> false if Xsep=0
!                  true, otherwise
!
!      With data on streamlines, the program calculates:
!      - influence matrix for each streamline
!      - boundary layer var distribution
!      - make a Newton-Raphson solver for Lag entrainment method
!      - make viscous corrections for potential stream velocity
!      - print out viscous flow characteristics for each velocity
!*****
!*      Program created by Augusto Elisio Lessa Veiga      *
!*      FSIG - University of Southampton/2003              *
!*      Sugestions are welcome                             *
!******
    include 'section.inc'
    include 'panel.inc'
    include 'xfoil.inc'

    integer Nstr,Npan,Npb,Nmax,Nte,Nsec
    real:: Rey,visc,alpha,Dlimit,pi,EPS1
    real,dimension(3):: tol
    real,dimension(2,150):: xsep,Xre,Xtr
    type (node),allocatable,dimension(:):: nd,bnd,wnd
    type (section),allocatable,dimension(:):: sec,wsec
    type (panel),allocatable,dimension(:):: pan,bpan,wpan
    character*70 :: arqname,arqname2
    logical:: fl,inviscid,fsharp
    BIJ = 0
    CIJ = 0
    DIJ = 0
    ! pflag -> flag that indicates if it's to apply
    !       wall pressure correction
    pi = 4. * atan(1.)
    call read_set(arqname,arqname2,Rey,visc,alpha,ACRIT,
    &              Nit,Nsec,Nmax,
    &              xtr,TFORCE,VACCEL,EPS1,fl,inviscid,fsharp)
    alpha = alpha*pi/180
    ALFA = alpha
    open(1,file ='BL_log.txt')
    write(1,*) 'New problem'
    close(1)

    if (inviscid) then
        call read_prev(Nsec,Nmax,Nw)

```

```

        allocate (sec(Nsec),wsec(Nsec))
        call read_section(sec,wsec,Nsec,Nmax,Nw)
        write(*,*) '*****'
        write(*,*) '*                V I X                *'
        write(*,*) '*          Copyright, Augusto E. L. Veiga          *'
        write(*,*) '*          University of Southampton, 2004          *'
        write(*,*) '*                Version 1.0                *'
        write(*,*) '*****'
        write(*,*) 'previously interpolated sections...'
        N = Nmax
    else
        call read_N(arqname,Nt,Npan) !reads the number of sections
                                   !and nodes
        allocate(nd(Nt))           !allocate vector str
        allocate(pan(Npan),sec(Nsec),wsec(Nsec))
        !allocate(sec(Nsec))
        !initialize structure sec
        call read_nodes(arqname,nd,pan,Nt,Npan) !read body nodes
        if (mod(nmax,2)==0) then
            Nmax=Nmax+1
        Else
            Nmax=Nmax
        endif
        call read_uns(arqname2,pan,Npan,nd,nt,Nte)
        call find_wk(pan,Npan,Npw,Iwake)
        allocate(wpan(Npw))
        ! sections are interpolated on this routine
        write(*,*) '*****'
        write(*,*) '*                V I X                *'
        write(*,*) '*          Copyright, Augusto E. L. Veiga          *'
        write(*,*) '*          University of Southampton, 2004          *'
        write(*,*) '*                Version 1.0                *'
        write(*,*) '*****'
        write(*,*) 'Interpolating sections...'
        call wk_surface(pan,wpan,Npan,Npw,Iwake,wsec,Nsec,fl)
        call surface(pan,Npan,sec,Nsec,Nmax,Nte,fl)
        ! there is no need anymore for such heavy structure
        deallocate(nd,pan,wpan)
        N = sec(1).Nnd
        Nw = wsec(1).Nnd
    endif
    write(*,*)
    write(*,*) '...calculating viscous flow'
    !i=int(Nsec/2)
    do i=1,Nsec
        !calculate geometric curvature for each node
        call G_curv(sec(i),Nmax,pflag)
        !find very first point on leading edge
        call find_Nlower(sec(i),Nmax)
        REINF = Rey
        QINF = 1.0

        XSTRIP(1) = xtr(1,i) !localizes transition upper part
        XSTRIP(2) = xtr(2,i) !lower part
        XTE = sec(i).p(1,1)
        XLE = sec(i).p(1,sec(i).nlower)
        YTE = sec(i).p(2,1)
        YLE = sec(i).p(2,sec(i).nlower)
        SLE = sec(i).s(sec(i).nlower)
        !fill up vars for xlib library
        call secset(sec(i),wsec(i),N,Nw,QINV,x,y,s)
        !ALFA = -sec(i).aflow
        open(2,file = 'BL_log.txt',position = 'APPEND')
        write(2,10) i
    close(2)
        ! This part solves the viscous flow for each section
        call VIX(sec(i),wsec(i),Nmax,EPs1,NQX,Nit,Nsec,i,fsharp)
        sec(i).tr(1) = ITRAN(1)
        sec(i).tr(2) = ITRAN(2)
        sec(i).Nbl(1) = NBL(1)
    enddo

```

```

sec(i).Nbl(2) = NBL(2)
do is = 1,2
  do ip =2,iblte(is)
    sec(i).ibl(ip,is) = IPAN(ip,is)
  enddo
  sec(i).iblte(is) = iblte(is)
  do in = 1,sec(i).Nbl(is)
    sec(i).TAU(in,is) = TAU(in,is)
  enddo
enddo
enddo
!perform lift, drag and viscous pressure resistance calculations
!call trefftz(wsec,Nsec,20,Dtrefftz) !calculates inviscid induced drag
call panmk(wsec,sec,Nsec,Nw,Nmax,CLvis,CLinv,cdf,cd,
&          cdi,cdiv,area,Cmx,Cmy,Cmz,zcp)
!call vii_graph(sec,Nsec,Nmax)
write(*,*)
write(*,*) '...printing results'
call print_result(sec,Nsec,Nw,CLvis,CLinv,cdf,cd,
&              cdi,cdiv,area,Cmx,Cmy,Cmz,zcp)
call print_blvar(sec,Nsec,Nmax)
call post_process(sec,Nsec,Nmax) !organize sectional plots
!reorganise panels using the sections again and print
! files to be used by PANVISE
!call makepan
deallocate(sec,wsec)
10  format('Section =',i4)
END

subroutine VIX(sec,wsec,Nmax,EPS1,NQX,NIT,Nsec,isec,fsharp)
! This subroutine receives the following variables:
! Geometry:      x,y,z and s of each section
! Inviscid flow:  QINV for each section
!                QINV for wake section
! Data           N -> number of section points
!                Nw-> number of wake section points
!                Nlower -> leading edge point that
!                        divides upper and lower parts
!                tr(.) -> transition point on lower and upper
!                        parts
!                Nsep -> separation points on lower and upper parts
!                Nre -> reattachment points on lower and upper parts
!                isec -> index of section
!                        if isec=1 or isec=Nsec, then viscous flow is not cal
!                        culated
! ...And spills out the following:
! Viscous flow:
! QVIS           -> viscous velocity
! Dstr           -> displacement thickness
! Thet          -> momentum thickness
! Ctau          -> sqrt (max shear coefficient
! H             -> shape parameter
! Cf            -> friction coefficient
! Dis           -> dissipation coefficient
!
! *****
!           A T E N T I O N ! ! !
!
! Points are input on counterclockwise order and this continues like that.
! Do not use the inverse order or you may experience problems
! *****
! This program was modified by Augusto Elisio Lessa Veiga and
! uses parts of the GNU software XFOIL
! *****
!* Author: Augusto Elisio Lessa Veiga *
!* University of Southampton, 2004 *
!* (Made in Brasil) *
! *****
include 'section.inc'
include 'xfoil.inc'
include 'xbl.inc'

```

```

type(section) :: sec,wsec
real,dimension(N+Nw):: Up,xbd,sbd
real,dimension(Nw):: Upw
real,dimension(3,N):: p
real,dimension(3,Nw):: pw
real,dimension(Izx+Iwx):: z
real:: clsec,cdsec,EPS1
real,dimension(ix,2):: pcor
integer N,Nw,Nlower,Nsec,isec
logical:: fsharp

do i = 1,N
  Up(i) = sec.Up(i)
  xbd(i) = X(i)
  sbd(i) = S(i)
enddo
j = 0
do i = N+1,N+Nw
  j = j+1
  Up(i) = wsec.up(j)
enddo
!main settings
pcor = 0 !22/05/2005
PI = 4.0*ATAN(1.0)
HOPI = 0.50/PI
QOPI = 0.25/PI
DTOR = PI/180.0
C---- default Cp/Cv (air)
GAMMA = 1.4
GAMM1 = GAMMA - 1.0
C---- initialize freestream Mach number to zero
MATYP = 1
MINF1 = 0.
MINF = 0.

CL = 0.
CM = 0.
CD = 0.

SIGTE = 0.0
GAMTE = 0.0
SIGTE_A = 0.
GAMTE_A = 0.

SIG = 0.

SHARP = .true. !if trailing edge is sharp
LIMAGE = .FALSE. !if image airfoil is present
LGAMU = .TRUE. !if GAMU arrays exist for current airfoil geometry
LQINU = .TRUE. !if QINVU arrays exist for current airfoil geometry
LVISC = .TRUE. !if viscous option is invoked
LALFA = .TRUE. !if alpha is specified, .FALSE. if CL is specified
LWAKE = .TRUE. !if wake geometry has been calculated
C LPACC = .TRUE. if each point calculated is to be saved
  LBLINI = .FALSE. !if BL has been initialized
  LIPAN = .TRUE. !if BL->panel pointers IPAN have been calculated
C LQAIJ = .TRUE. if dPsi/dGam matrix has been computed and factored
  LADIJ = .FALSE. !if dQ/dSig matrix for the airfoil has been computed
  LWDIJ = .FALSE. !if dQ/dSig matrix for the wake has been computed
C LQVDES = .TRUE. if viscous Ue is to be plotted in QDES routines
C LQSPEC = .TRUE. if Qspec has been initialized
C LQREFL = .TRUE. if reflected Qspec is to be plotted in QDES routines
  LVCONV = .FALSE. !if converged BL solution exists
C LCPREF = .TRUE. if reference data is to be plotted on Cp vs x/c plots
C LLOCK = .TRUE. if source airfoil coordinates are clockwise
C LPFILE = .TRUE. if polar file is ready to be appended to
C LPFILX = .TRUE. if polar dump file is ready to be appended to
C LPPSHO = .TRUE. if CL-CD polar is plotted during point sequence
C LBFLAP = .TRUE. if buffer airfoil flap parameters are defined
C LFLAP = .TRUE. if current airfoil flap parameters are defined
C LEIW = .TRUE. if unit circle complex number array is initialized
C LSCINI = .TRUE. if old-airfoil circle-plane arc length s(w) exists

```

```

C   LFOREF      .TRUE. if CL,CD... data is to be plotted on Cp vs x/c plots
C   LNORM       .TRUE. if input buffer airfoil is to be normalized
C   LGSAME      .TRUE. if current and buffer airfoils are identical
C
C   LPLCAM      .TRUE. if thickness and camber are to be plotted
C   LQSYM       .TRUE. if symmetric Qspec will be enforced
C   LGSYM       .TRUE. if symmetric geometry will be enforced
C   LQGRID      .TRUE. if grid is to overlaid on Qspec(s) plot
C   LGGRID      .TRUE. if grid is to overlaid on buffer airfoil geometry plot
C   LGTICK      .TRUE. if node tick marks are to be plotted on buffer airfoil
C   LQSLOP      .TRUE. if modified Qspec(s) segment is to match slopes
C   LGSLOP      .TRUE. if modified geometry segment is to match slopes
C   LCSLOP      .TRUE. if modified camber line segment is to match slopes
C   LQSPPL      .TRUE. if current Qspec(s) in in plot
C   LGEOP      .TRUE. if current geometry in in plot
C   LCPGRD      .TRUE. if grid is to be plotted on Cp plots
C   LBLGRD      .TRUE. if grid is to be plotted on BL variable plots
C   LBLSYM      .TRUE. if symbols are to be plotted on BL variable plots
C   LCMINP      .TRUE. if min Cp is to be written to polar file for cavitation
C   LHMOMP      .TRUE. if hinge moment is to be written to polar file
C
C   LPGRID      .TRUE. if polar grid overlay is enabled
C   LPCDW       .TRUE. if polar CDwave is plotted
C   LPLIST      .TRUE. if polar listing lines (at top of plot) are enabled
C   LPLEGN      .TRUE. if polar legend is enabled
C
C   LPLOT       .TRUE. if plot page is open
C   LSYM        .TRUE. if symbols are to be plotted in QDES routines
C   LIQSET      .TRUE. if inverse target segment is marked off in QDES
C   LCLIP       .TRUE. if line-plot clipping is to be performed
C   LVLAB       .TRUE. if label is to be plotted on viscous-variable plots
C   LCURS       .TRUE. if cursor input is to be used for blowups, etc.
C   LLAND       .TRUE. if Landscape orientation for PostScript is used

      call stagpoint (Up,sbd,GAM,xbd,N,Nw,IST,SST,SST_GO,SST_GP,fsharp)
      sec.istag = IST

      call SIC(sec,wsec,N,Nw)
!      DATA EPS1 / 1.0E-4 /
C
      NITER = 10
      QINF = 1.0
C
C
C---- set velocities on wake from airfoil vorticity for alpha=0, 90
C      CALL QWCALC
C
C---- set velocities on airfoil and wake for initial alpha
C      CALL QISET
C
C
C---- locate stagnation point arc length position and panel index
!      CALL STFIND
C
C---- set BL position -> panel position pointers
      CALL IBLPAN
      sec.iblte(1) = iblte(1)
      sec.iblte(2) = iblte(2)
C
C---- calculate surface arc length array for current stagnation point location
      CALL XICALC
C
C---- set BL position -> system line pointers
      CALL IBLSYS
C
C
C---- set inviscid BL edge velocity UINV from QINV
      CALL UICALC
C

      IF (.NOT.LBLINI) THEN

```

```

C
C----- set initial Ue from inviscid Ue
DO IBL=1, NBL(1)
    UEDG(IBL,1) = UINV(IBL,1)
ENDDO

C
DO IBL=1, NBL(2)
    UEDG(IBL,2) = UINV(IBL,2)
ENDDO

C
ENDIF
!initial lift calculation (inviscid)
Nref = iqx
call clcalc2(N,Nref,sec,gam_a,alfa,minf,qinf,pcor,
&             XCMREF,YCMREF,CL,CM,CDP,CL_ALF,CL_MSQ)
!    CALL CLCALC(N,X,Y,GAM,GAM_A,ALFA,MINF,QINF, XCMREF,YCMREF,
!    &             CL,CM,CDP,CL_ALF,CL_MSQ)
sec.cli = CL

C
C
C----- Newton iteration for entire BL solution
NITER = Nit
if (isec>1 .and. isec<Nsec) then

    WRITE(*,*) 'Solving BL system ...'
    open(1,file = 'BL_log.txt',position = 'APPEND')
    DO 1000 ITER=1, NITER

C
C----- fill Newton system for BL variables
CALL SETBL

C
C----- solve Newton system with custom solver
CALL BLSOLV

C
C----- update BL variables
CALL UPDATE(sec,pcor) !output pcor (22/05/2005)

C
C    IF(LALFA) THEN
C----- set new freestream Mach, Re from new CL
C    CALL MRCL(CL,MINF_CL,REINF_CL)
C    CALL COMSET
C    ELSE
C----- set new inviscid speeds QINV and UINV for new alpha
C    CALL QISET
C    CALL UICALC
C    ENDIF
C
C----- calculate edge velocities QVIS(.) from UEDG(..)
CALL QVFUE

C
C----- set GAM distribution from QVIS
CALL GAMQV

C
C----- relocate stagnation point
!    CALL STMOVE
C
C----- set updated CL,CD
!Nref = iqx
call clcalc2(N,Nref,sec,gam_a,alfa,minf,qinf,pcor,
&             XCMREF,YCMREF,CL,CM,CDP,CL_ALF,CL_MSQ)
sec.clv = CL
!    CALL CLCALC(N,X,Y,GAM,GAM_A,ALFA,MINF,QINF, XCMREF,YCMREF,
!    &             CL,CM,CDP,CL_ALF,CL_MSQ)
CALL CDCALC
sec.cdf = CDF !sectional frict. Cd
sec.cd = CD
if (iter==1) then
!    sec.cli = CL
!endif
C
C----- display changes and test for convergence
IF(RLX.LT.1.0)

```

```

& WRITE(1,2000) ITER, RMSBL, RMXBL, VMXBL, IMXBL, ISMXBL, RLX
IF(RLX.EQ.1.0)
& WRITE(1,2010) ITER, RMSBL, RMXBL, VMXBL, IMXBL, ISMXBL
  CDP = CD - CDF
  WRITE(1,2020) ALFA/DTOR, CL, CM, CD, CDF, CDP
C
  IF(RMSBL .LT. EPS1) THEN
    LVCONV = .TRUE.
    AVISC = ALFA
    MVIS = MINF
    GO TO 90
  ENDIF
C
1000 CONTINUE
  WRITE(1,*) 'VISCAL: Convergence failed'
C
  90 CONTINUE
  close(1)

endif !avoiding tip sections
!filling up vectors
sec.clv = CL
DO IS=1, 2
  DO IBL=2, NBL(IS)
    I = ibl !IPAN(IBL,IS)
    sec.Q(1,i,is) = Ctau(i,is) !shear stress or critical amp
    sec.Q(2,i,is) = thet(i,is) !momentum thick
    sec.Q(3,i,is) = dstr(i,is) !recording displacement thick
    sec.Q(4,i,is) = uedg(i,is) !recording viscous velocity
    sec.UINV(i,is) = UINV(i,is) !recording inviscid velocity
  enddo
  ! wake variables
  if (is==2) then
    iwk=0
    do ibl=iblte(is)+1,iblte(is)+wsec.Nnd
      iwk = iwk+1
      i = ibl
      wsec.Q(1,iwk,is) = Ctau(i,is) !shear stress or critical amp
      wsec.Q(2,iwk,is) = thet(i,is) !momentum thick
      wsec.Q(3,iwk,is) = dstr(i,is) !recording displacement thick
      wsec.Q(4,iwk,is) = uedg(i,is) !recording viscous velocity
    enddo
  endif
enddo
if (isec>1 .and. isec<Nsec) then
  clsec = cl
  cdsec = cd
  sec.CD = CD
  sec.Cdf = Cdf
  sec.Cdp = Cdp
  sec.iter = iter
else
  clsec = sec.cli
  cdsec = 0
  sec.CD = 0
  sec.Cdf = 0
  sec.Cdp = 0
  sec.iter = 0
endif
RETURN
C.....
2000 FORMAT
& (/1X,I3,' rms: ',E10.4,' max: ',E10.4,3X,A1,' at ',I4,I3,
& ' RLX:',F6.3)
2010 FORMAT
& (/1X,I3,' rms: ',E10.4,' max: ',E10.4,3X,A1,' at ',I4,I3)
2020 FORMAT
& ( 1X,3X,' a =', F7.3,' CL =',F8.4 /
& 1X,3X,' Cm =', F8.4,' CD =',F9.5,
& ' => CDF =',F9.5,' CDP =',F9.5)
END subroutine ! VIX

```



```

!*****
! This subroutine was taken from XFOIL code
!*****

      SUBROUTINE CDCALC
      INCLUDE 'XFOIL.INC'

C
      SA = SIN(ALFA)
      CA = COS(ALFA)

C
      IF(LVISC .AND. LBLINI) THEN
C
C----- set variables at the end of the wake
      THWAKE = THET(NBL(2),2)
      URAT = UEDG(NBL(2),2)/QINF
      UEWAKE = UEDG(NBL(2),2) * (1.0-TKLAM) / (1.0 - TKLAM*URAT**2)
      SHWAKE = DSTR(NBL(2),2)/THET(NBL(2),2)

C
C----- extrapolate wake to downstream infinity using Squire-Young relation
C      (reduces errors of the wake not being long enough)
      CD = 2.0*THWAKE * (UEWAKE/QINF)**(0.5*(5.0+SHWAKE))

C
      ELSE

C
      CD = 0.0

C
      ENDIF

C
C----- calculate friction drag coefficient
      CDF = 0.0
      DO 20 IS=1, 2
        DO 205 IBL=3, IBLTE(IS)
          I = IPAN(IBL,IS)
          IM = IPAN(IBL-1,IS)
          DX = (X(I) - X(IM))*CA + (Y(I) - Y(IM))*SA
          CDF = CDF + 0.5*(TAU(IBL,IS)+TAU(IBL-1,IS))*DX * 2.0/QINF**2
        205 CONTINUE
      20 CONTINUE

C
      RETURN
      END ! CDCALC
!*****
! This subroutine include on CL the three-dimensional effects
!*****

      SUBROUTINE CLCALC2(N,Nref,sec,gam_a,ALFA,MINF,QINF,pcor,
&                      XREF,YREF,
&                      CL,CM,CDP, CL_ALF,CL_MSQ)
      include 'section.inc'

C-----
C      Integrates surface pressures to get CL and CM.
C      Integrates skin friction to get CDF.
C      Calculates dCL/dAlpha for prescribed-CL routines.
C      Modified by Augusto Veiga
C-----

      real,dimension(Nref,2):: pcor
      type(section):: sec
      REAL:: MINF,v
      real:: dui(N),gam_a(N),x(N),y(N)

C
C----- moment-reference coordinates
ccc      XREF = 0.25
ccc      YREF = 0.

C
C      transforming Vpot into pressure coefficient
      do i = 1,N
        sum = 0
        x(i) = sec.p(1,i)
        y(i) = sec.p(2,i)
        do k = 1,3
          v = sec.vpot(k,i)
          sum = sum + v**2
        enddo
      enddo

```

```

        sec.cp(i) = 1.0-sum
        enddo
!transforming pcor into dui
        j = sec.istag
        do i = 1,sec.iblte(1)
            dui(j) = pcor(i,1)
            j = j-1
        enddo
        j = sec.istag
        do i = 2,sec.iblte(2)
            j = j+1
            dui(j) = pcor(i,2)
        enddo

        SA = SIN(ALFA)
        CA = COS(ALFA)

C
        BETA = SQRT(1.0 - MINF**2)
        BETA_MSQ = -0.5/BETA

C
        BFAC      = 0.5*MINF**2 / (1.0 + BETA)
        BFAC_MSQ = 0.5          / (1.0 + BETA)
&      - BFAC      / (1.0 + BETA) * BETA_MSQ

C
        CL = 0.0
        CM = 0.0

        CDP = 0.0

C
        CL_ALF = 0.
        CL_MSQ = 0.

C
        I = 1
        CGINC = sec.cp(i) + dui(i)**2
        CPG1   = CGINC/(BETA + BFAC*CGINC)
        CPG1_MSQ = -CPG1/(BETA + BFAC*CGINC)*(BETA_MSQ + BFAC_MSQ*CGINC)

C
        CPI_GAM = -2.0*cginc
        CPC_CPI = (1.0 - BFAC*CPG1)/ (BETA + BFAC*CGINC)
        CPG1_ALF = CPC_CPI*CPI_GAM*GAM_A(I)

C
        DO 10 I=1, N
            IP = I+1
            IF(I.EQ.N) IP = 1

C
            CGINC = sec.cp(i) + dui(i)**2
            CPG2   = CGINC/(BETA + BFAC*CGINC)
            CPG2_MSQ = -CPG2/(BETA + BFAC*CGINC)*(BETA_MSQ + BFAC_MSQ*CGINC)

C
            CPI_GAM = -2.0*cginc
            CPC_CPI = (1.0 - BFAC*CPG2)/ (BETA + BFAC*CGINC)
            CPG2_ALF = CPC_CPI*CPI_GAM*GAM_A(IP)

C
            DX = (X(IP) - X(I))*CA + (Y(IP) - Y(I))*SA
            DY = (Y(IP) - Y(I))*CA - (X(IP) - X(I))*SA
            DG = CPG2 - CPG1

C
            AX = (0.5*(X(IP)+X(I))-XREF)*CA + (0.5*(Y(IP)+Y(I))-YREF)*SA
            AY = (0.5*(Y(IP)+Y(I))-YREF)*CA - (0.5*(X(IP)+X(I))-XREF)*SA
            AG = 0.5*(CPG2 + CPG1)

C
            DX_ALF = -(X(IP) - X(I))*SA + (Y(IP) - Y(I))*CA
            AG_ALF = 0.5*(CPG2_ALF + CPG1_ALF)
            AG_MSQ = 0.5*(CPG2_MSQ + CPG1_MSQ)

C
            CL      = CL      + DX* AG
            CDP     = CDP     - DY* AG
            CM      = CM      - DX*(AG*AX + DG*DX/12.0)
&              - DY*(AG*AY + DG*DY/12.0)

C
            CL_ALF = CL_ALF + DX*AG_ALF + AG*DX_ALF

```

```

        CL_MSQ = CL_MSQ + DX*AG_MSQ
C
        CPG1 = CPG2
        CPG1_ALF = CPG2_ALF
        CPG1_MSQ = CPG2_MSQ
10 CONTINUE
C
        RETURN
        END ! CLCALC2

```

## VIX Surface Interpolation

```

        subroutine surface(pan,Npan,sec,Nsec,Nmax,Nte,fl)
        include 'section.inc'
        include 'panel.inc'
!       This subroutine calculates collocation points for each panel
!       calculates mean edges
!       calculates surface coordinate for each collocation point on s and t
!       gets points and interpolate cp using a spline distribution
!*****
!*       (C) Augusto Veiga, University of Southampton 2003      *
!*****
        integer Npan,Nsec,Nv,Nh,Nb,Nmax
        type(panel),dimension(Npan):: pan,panb
        type(section),dimension(Nsec):: sec
        real:: S(3),Dmax
        logical:: fl
!Calculating collocation points
Do i=1,Npan
        S=0
        do j=1,4
                S(1)=S(1)+pan(i).nd(j).x
                S(2)=S(2)+pan(i).nd(j).y
                S(3)=S(3)+pan(i).nd(j).z
        enddo
        pan(i).co.x=S(1)/4.
        pan(i).co.y=S(2)/4.
        pan(i).co.z=S(3)/4.
enddo
!call smooth_cp(pan,Npan)
!Calculating surface coordinates s and t
call org_pan(pan,panb,Npan,Nb,Nv,Nh,Nte)
call surf_coord(panb,Nb,Nv,Nh)
call surf_spl(panb,Nb,Nv,Nh,fl) !Makes a spline surface
do i = 1,3
        call press_int(panb,Nb,Nv,Nh,i,.false.)!Makes a V(x,y,z) surface
! i -> 1 = x
!       2 = y
!       3 = z
enddo
!dividing surface into sections with 100 points equally
! spaced each and following a plane which normal is the
! slope at yz plane
do i=1,Nsec
        sec(i).Nnd=Nmax
enddo
call interpol_surf(panb,Nb,Nv,Nsec,sec,Nmax)
call sec_normal(sec,Nsec,Nmax) !calculate normals on each point of section
!call calc_Up(sec,Nsec,Nmax) !calculate modular potential velocity
call vsec_plot(Sec,Nsec,Nmax) !plots spanwise velocity

return
end subroutine

        subroutine interpol_surf(pan,Npan,Nv,Nsec,sec,Nmax)
        include 'section.inc'
        include 'panel.inc'

        integer :: Npan,Nv,Nsec,Nmax,iv,ih
        type(panel),dimension(Npan) :: pan
        type (section),dimension(Nsec):: sec

```

```

real:: tc,sc,t

! divide body onto spanwise sections with tc spacement
tc= (pan(Nv).nd(4).t-pan(1).nd(1).t)/(Nsec-1)
t=pan(1).nd(1).t
k=1
id=0
ih=1
iv=1
i=1
do while (i<=Nsec)
  !This loop makes geometry interpolation just
  !finding s last points
  sc=1.0
  sec(i).id=id
  id=id+1
  sec(i).s(1)=0
  sec(i).s(Nmax)=sc
  sec(i).t(1)=t
  sc=(sec(i).s(Nmax)-sec(i).s(1))/(Nmax-1)
  if ( t<=pan(iv).nd(4).t) then
    ih=iv
    call interpol_sec(sec(i),1,pan,Npan,ih,iv,Nv,1) !x
    call interpol_sec(sec(i),1,pan,Npan,ih,iv,Nv,2) !y
    call interpol_sec(sec(i),1,pan,Npan,ih,iv,Nv,3) !z
    !call interpol_Vs(sec(i),1,pan,Npan,ih,iv,Nv) !Vm
    do j=2,Nmax
      if (j==Nmax) then
        sec(i).s(j)=pan(Npan-(Nv-iv)).nd(2).s
        sec(i).t(j)=t
      else
        sec(i).s(j)=sc+sec(i).s(j-1)
        sec(i).t(j)=t
      endif
    enddo
    call interpol_sec(sec(i),j,pan,Npan,ih,iv,Nv,1) !x
    call interpol_sec(sec(i),j,pan,Npan,ih,iv,Nv,2) !y
    call interpol_sec(sec(i),j,pan,Npan,ih,iv,Nv,3) !z
    !call interpol_Vs(sec(i),j,pan,Npan,ih,iv,Nv) !Vm
    t=t+tc
    i=i+1
  else if (i==Nsec) then
    ih=Nv
    !Last section
    call interpol_sec(sec(i),1,pan,Npan,ih,iv,Nv,1) !x
    call interpol_sec(sec(i),1,pan,Npan,ih,iv,Nv,2) !y
    call interpol_sec(sec(i),1,pan,Npan,ih,iv,Nv,3) !z
    !call interpol_Vs(sec(i),1,pan,Npan,ih,iv,Nv) !Vm
    do j=2,Nmax
      if (j==Nmax) then
        sec(i).s(j)=pan(Npan).nd(2).s
        sec(i).t(j)=pan(Npan).nd(3).t
      else
        sec(i).s(j)=sc+sec(i).s(j-1)
        sec(i).t(j)=t
      endif
    enddo
    call interpol_sec(sec(i),j,pan,Npan,ih,iv,Nv,1) !x
    call interpol_sec(sec(i),j,pan,Npan,ih,iv,Nv,2) !y
    call interpol_sec(sec(i),j,pan,Npan,ih,iv,Nv,3) !z
    !call interpol_Vs(sec(i),j,pan,Npan,ih,iv,Nv) !Vm
    enddo
    i=i+1
  else if (iv<Nv) then
    iv=iv+1
  endif
enddo
!Now we calculate the cossine of sectional segments
do isec = 1,Nsec
  do j = 1,Nmax
    if (j==Nmax) then
      do k = 1,3
        sec(isec).vcos(k,j) = sec(isec).vcos(k,j-1)
      enddo
    enddo
  enddo
enddo

```

```

        enddo
    else
        !Calculate length
        soma = 0
        do k = 1,3
            soma = soma+ (sec(isec).p(k,j+1)-sec(isec).p(k,j))**2
        enddo
        !calculate cossine
        do k = 1,3
            sec(isec).vcos(k,j)=(sec(isec).p(k,j+1)-sec(isec).p(k,j))
&            /sqrt(soma)
        enddo
    endif
enddo
enddo
!Now we interpolate sectional velocity (Vs)
!...and calculate the tangential velocity q
iv = 1
ih = 1
i =1
do while(i<=Nsec)
    sc=1.0
    t = sec(i).t(1)
    if ( t<=pan(iv).nd(4).t) then
        ih=iv
        call interpol_Vs(sec(i),1,pan,Npan,ih,iv,Nv) !Vm
        do j=2,Nmax
            call interpol_Vs(sec(i),j,pan,Npan,ih,iv,Nv) !Vm
        enddo
        i = i+1
    else if (i==Nsec) then
        ih=Nv !Last section
        call interpol_Vs(sec(i),1,pan,Npan,ih,iv,Nv) !Vm
        do j=2,Nmax
            call interpol_Vs(sec(i),j,pan,Npan,ih,iv,Nv) !Vm
        enddo
        i = i+1
    else if (iv<Nv) then
        iv=iv+1
    endif
enddo
enddo

! Writing a scratch file of potential tangential velocity
open(1,file='potential_scratch.txt')
write(1,100)
do isec = 1,Nsec
    do j = 1,Nmax
        write(1,200) j,sec(isec).Up(j),sec(isec).vcos(1,j),
&        sec(isec).vcos(2,j),sec(isec).vcos(3,j)
    enddo
    write(1,*)
enddo
close(1)

return
100 format('node , velocity , cos x , cos y , cos z')
200 format(i4,1x,f8.3,1x,f8.3,1x,f8.3,1x,f8.3)
end subroutine

subroutine interpol_sec(sec,j,pan,Npan,ih,iv,Nv,Nflag)
include 'section.inc'
include 'panel.inc'
! This subroutine gets the current section with s,t variables and,
! marching along chordwise panels, interpolates x,y,z using Coons
! bicubic spline surface
! As the solution marches, ih is changed to the next panel on
! chordwise direction.
! Nflag chooses what is going to be interpolated
! 1 -> x coordinate
! 2 -> y coordinate
! 3 -> z coordinate
! 4 -> cp

```

```

!
!      Bd1, Bd2 -> first derivative coefficients
!*****
!*      (C) Augusto Veiga, University of Southampton, 2003      *
!*****

      integer j,ih,iv,Nv,Nflag
      type(section):: sec
      type(panel),dimension(Npan)::pan
      real::s,w,T,U,A(4),B1(4),B2(4),soma
      real:: delta1,delta2,Vn(4),C(4),D(4),dxs2(4),dxt2(4),x(4)
      real:: tiny

      tiny=1.0E-10
      soma=0;A=0;B1=0;B2=0;Bd1=0;Bd2=0
      s=sec.s(j)
      w=sec.t(j)
      i=ih
      lpl: do
      select case (Nflag)
      case(1) !options to interpolate variables
      do k=1,4
      dxs2(k)=pan(i).nd(k).dxs2
      dxt2(k)=pan(i).nd(k).dxt2
      x(k)=pan(i).nd(k).x
      enddo
      case(2)
      do k=1,4
      dxs2(k)=pan(i).nd(k).dys2
      dxt2(k)=pan(i).nd(k).dyt2
      x(k)=pan(i).nd(k).y
      enddo
      case(3)
      do k=1,4
      dxs2(k)=pan(i).nd(k).dzs2
      dxt2(k)=pan(i).nd(k).dzt2
      x(k)=pan(i).nd(k).z
      enddo
      case(4)
      do k=1,4
      dxs2(k)=0
      dxt2(k)=0
      x(k)=pan(i).nd(k).up
      enddo
      end select
      if (s<pan(i).nd(2).s) then
      !calculate normal coefficients T and U
      T=(s-pan(i).nd(1).s)/(pan(i).nd(2).s-pan(i).nd(1).s)
      U=(w-pan(i).nd(1).t)/(pan(i).nd(4).t-pan(i).nd(1).t)
      !calculate coefficients A
      A(1)=(1-T)*(1-U)
      A(2)=T*(1-U)
      A(3)=T*U
      A(4)=(1-T)*U
      !calculate coefficients B
      delta1=(pan(i).nd(2).s-pan(i).nd(1).s)
      delta2=(pan(i).nd(4).t-pan(i).nd(1).t)
      do k=1,4
      B1(k)=(A(k)**3-A(k))/6.*delta1
      B2(k)=(A(k)**3-A(k))/6.*delta2
      enddo
      !calculating normal vectors of 2nd derivatives
      C(1)=B1(2)*dxs2(2)-B1(1)*dxs2(1)
      C(2)=B1(3)*dxs2(3)-B1(2)*dxs2(2)
      C(3)=B1(4)*dxs2(4)-B1(3)*dxs2(3)
      C(4)=B1(4)*dxs2(4)-B1(1)*dxs2(1)
      D(1)=B2(2)*dxt2(2)-B2(1)*dxt2(1)
      D(2)=B2(3)*dxt2(3)-B2(2)*dxt2(2)
      D(3)=B2(4)*dxt2(4)-B2(3)*dxt2(3)
      D(4)=B2(4)*dxt2(4)-B2(1)*dxt2(1)
      Vn(1)=c(1)*d(4)-d(1)*c(4)
      Vn(2)=c(1)*d(2)-d(1)*c(2)
      Vn(3)=c(3)*d(2)-d(3)*c(2)

```

```

Vn(4)=c(3)*d(4)-d(3)*c(4)

!calculate final value for interpolation
soma=0
do k=1,4
    soma=soma+A(k)*x(k)+Vn(k)
enddo
if (abs(soma)<tiny) then
    soma=0
endif
exit lp1
else if (s==pan(Npan-(Nv-iv)).nd(2).s) then
    !calculate normal coefficients T and U
    T=(s-pan(i).nd(1).s)/(pan(i).nd(2).s-pan(i).nd(1).s)
    U=(w-pan(i).nd(1).t)/(pan(i).nd(4).t-pan(i).nd(1).t)
    !calculate coefficients A
    A(1)=(1-T)*(1-U)
    A(2)=T*(1-U)
    A(3)=T*U
    A(4)=(1-T)*U
    !calculate coefficients B
    delta1=(pan(i).nd(2).s-pan(i).nd(1).s)
    delta2=(pan(i).nd(4).t-pan(i).nd(1).t)
    do k=1,4
        B1(k)=(A(k)**3-A(k))*delta1
        B2(k)=(A(k)**3-A(k))*delta2
    enddo
    !calculating normal vectors of 2nd derivatives
    C(1)=B1(2)*dxs2(2)-B1(1)*dxs2(1)
    C(2)=B1(3)*dxs2(3)-B1(2)*dxs2(2)
    C(3)=B1(4)*dxs2(4)-B1(3)*dxs2(3)
    C(4)=B1(4)*dxs2(4)-B1(1)*dxs2(1)
    D(1)=B2(2)*dxt2(2)-B2(1)*dxt2(1)
    D(2)=B2(3)*dxt2(3)-B2(2)*dxt2(2)
    D(3)=B2(4)*dxt2(4)-B2(3)*dxt2(3)
    D(4)=B2(4)*dxt2(4)-B2(1)*dxt2(1)
    Vn(1)=c(1)*d(4)-d(1)*c(4)
    Vn(2)=c(1)*d(2)-d(1)*c(2)
    Vn(3)=c(3)*d(2)-d(3)*c(2)
    Vn(4)=c(3)*d(4)-d(3)*c(4)
    !calculate final value for interpolation
    soma=0
    do k=1,4
        soma=soma+A(k)*x(k)+Vn(k)
    enddo
    if (abs(soma)<tiny) then
        soma=0
    endif
    exit lp1
else if (ih>Npan-(Nv-iv)) then
    exit lp1
else if (ih<=Npan-(Nv-iv)) then
    ih=ih+Nv
    i=ih
endif
enddo lp1

select case (Nflag)
case(1) !options to interpolate variables
    sec.p(1,j)=soma
case(2)
    sec.p(2,j)=soma
case(3)
    sec.p(3,j)=soma
! case(4)
!     sec.cp(j)=soma
end select

return
end subroutine

```

```

!-----
      subroutine interpol_Vs(sec,j,pan,Npan,ih,iv,Nv)
      include 'section.inc'
      include 'panel.inc'
!      This subroutine gets the current section with s,t variables and,
!      marching along chordwise panels, interpolates x,y,z using Coons
!      bicubic spline surface
!      As the solution marches, ih is changed to the next panel on
!      chordwise direction.
!      Nflag chooses what is going to be interpolated
!      1 -> x coordinate
!      2 -> y coordinate
!      3 -> z coordinate
!      4 -> cp
!*****
!*      (C) Augusto Veiga, University of Southampton, 2003      *
!*****
      integer j,ih,iv,Nv,Nflag
      type(section):: sec
      type(panel),dimension(Npan):: pan
      real::s,w,T,U,A(4),B1(4),B2(4),soma,V(3)
      real:: delta1,delta2,Vn(4),C(4),D(4),dxs2(4),dxt2(4),x(4)
      real:: tiny

      tiny=1.0E-10
      soma=0;A=0;B1=0;B2=0
      s=sec.s(j)
      w=sec.t(j)
      i=ih
      do ind =1,3
      lp1: do
        do k=1,4
          dxs2(k)=0
          dxt2(k)=0
          x(k)=pan(i).nd(k).vm(ind)
        enddo
        if (s<pan(i).nd(2).s) then
          !calculate normal coefficients T and U
          T=(s-pan(i).nd(1).s)/(pan(i).nd(2).s-pan(i).nd(1).s)
          U=(w-pan(i).nd(1).t)/(pan(i).nd(4).t-pan(i).nd(1).t)
          !calculate coefficients A
          A(1)=(1-T)*(1-U)
          A(2)=T*(1-U)
          A(3)=T*U
          A(4)=(1-T)*U
          !calculate coefficients B
          !calculating normal vectors of 2nd derivatives
          !calculate final value for interpolation
          soma=0
          do k=1,4
            soma=soma+A(k)*x(k)
          enddo
          if (abs(soma)<tiny) then
            soma=0
          endif
          exit lp1
        else if (s==pan(Npan-(Nv-iv)).nd(2).s) then
          !calculate normal coefficients T and U
          T=(s-pan(i).nd(1).s)/(pan(i).nd(2).s-pan(i).nd(1).s)
          U=(w-pan(i).nd(1).t)/(pan(i).nd(4).t-pan(i).nd(1).t)
          !calculate coefficients A
          A(1)=(1-T)*(1-U)
          A(2)=T*(1-U)
          A(3)=T*U
          A(4)=(1-T)*U
          !calculate coefficients B
          !calculating normal vectors of 2nd derivatives
          !calculate final value for interpolation
          soma=0
          do k=1,4
            soma=soma+A(k)*x(k)
          enddo
        endif
      enddo
    enddo
  enddo
enddo

```



```

        enddo
        if (abs(soma)<tiny) then
            soma=0
        endif
        exit lp1
    else if (ih>Npan-(Nv-iv)) then
        exit lp1
    else if (ih<=Npan-(Nv-iv)) then
        ih=ih+Nv
        i=ih
    endif
enddo lp1
V(ind) = soma
enddo !indexes

soma = 0
do k = 1,3
    sec.vpot(k,j) = v(k) !potential 3D velocity
    soma = soma+ (sec.vcos(k,j)*v(k))*2 !velocity on section
    !projects velocity on each segment of section
enddo

sec.Up(j) = sqrt(soma) !tangential velocity on section pt

return
end subroutine

```

## XFOIL Routines for Initial Boundary Layer Solution that Were Added to VIX

```

SUBROUTINE SETBL
C-----
C    Sets up the BL Newton system coefficients
C    for the current BL variables and the edge
C    velocities received from SETUP. The local
C    BL system coefficients are then
C    incorporated into the global Newton system.
C-----
    INCLUDE 'XFOIL.INC'
    INCLUDE 'XBL.INC'
    REAL USAV(IVX,2)
    REAL U1_M(2*IVX), U2_M(2*IVX)
    REAL D1_M(2*IVX), D2_M(2*IVX)
    REAL ULE1_M(2*IVX), ULE2_M(2*IVX)
    REAL UTE1_M(2*IVX), UTE2_M(2*IVX)
    REAL MA_CLMR, MSQ_CLMR, MDI

C
C---- set the CL used to define Mach, Reynolds numbers
    IF (LALFA) THEN
        CLMR = CL
    ELSE
        CLMR = CLSPEC
    ENDIF

C
C---- set current MINF(CL)
    !CALL MRCL(CLMR,MA_CLMR,RE_CLMR)
    MINF = 0
    RE_CLMR = 0
    MA_CLMR = 0
    CLMR = 0.000001
    MSQ_CLMR = 2.0*MINF*MA_CLMR

C
C---- set compressibility parameter TKLAM and derivative TK_MSQ
    !CALL COMSET

C
C---- set gas constant (= Cp/Cv)
    GAMBL = GAMMA
    GM1BL = GAMM1

C
C---- set parameters for compressibility correction
    QINFBL = QINF
    TKBL = TKLAM
    TKBL_MS = TKL_MSQ

```

```

C
C---- stagnation density and 1/enthalpy
RSTBL = (1.0 + 0.5*GM1BL*MINF**2) ** (1.0/GM1BL)
RSTBL_MS = 0.5*RSTBL/(1.0 + 0.5*GM1BL*MINF**2)
C
HSTINV = GM1BL*(MINF/QINFBL)**2 / (1.0 + 0.5*GM1BL*MINF**2)
HSTINV_MS = GM1BL*( 1.0/QINFBL)**2 / (1.0 + 0.5*GM1BL*MINF**2)
& - 0.5*GM1BL*HSTINV / (1.0 + 0.5*GM1BL*MINF**2)
C
C---- Sutherland's const./To (assumes stagnation conditions are at STP)
HVRAT = 0.35
C
C---- set Reynolds number based on freestream density, velocity, viscosity
HERAT = 1.0 - 0.5*QINFBL**2*HSTINV
HERAT_MS = - 0.5*QINFBL**2*HSTINV_MS
C
REYBL = REINF * SQRT(HERAT**3) * (1.0+HVRAT)/(HERAT+HVRAT)
REYBL_RE = SQRT(HERAT**3) * (1.0+HVRAT)/(HERAT+HVRAT)
REYBL_MS = REYBL * (1.5/HERAT - 1.0/(HERAT+HVRAT))*HERAT_MS
C
AMCRIT = ACRIT
C
C---- save TE thickness
DWTE = WGAP(1)
C
IF(.NOT.LBLINI) THEN
C----- initialize BL by marching with Ue (fudge at separation)
WRITE(*,*)
WRITE(*,*) 'Initializing BL ...'
CALL MRCHUE
LBLINI = .TRUE.
ENDIF
C
WRITE(*,*)
C
C---- march BL with current Ue and Ds to establish transition
CALL MRCHDU
C
DO 5 IS=1, 2
DO 6 IBL=2, NBL(IS)
USAV(IBL, IS) = UEDG(IBL, IS)
6 CONTINUE
5 CONTINUE
C
CALL UESET
C
DO 7 IS=1, 2
DO 8 IBL=2, NBL(IS)
TEMP = USAV(IBL, IS)
USAV(IBL, IS) = UEDG(IBL, IS)
UEDG(IBL, IS) = TEMP
8 CONTINUE
7 CONTINUE
C
ILE1 = IPAN(2,1)
ILE2 = IPAN(2,2)
ITE1 = IPAN(IBLTE(1),1)
ITE2 = IPAN(IBLTE(2),2)
C
JVTE1 = ISYS(IBLTE(1),1)
JVTE2 = ISYS(IBLTE(2),2)
C
DULE1 = UEDG(2,1) - USAV(2,1)
DULE2 = UEDG(2,2) - USAV(2,2)
C
C---- set LE and TE Ue sensitivities wrt all m values
DO 10 JS=1, 2
DO 110 JBL=2, NBL(JS)
J = IPAN(JBL, JS)
JV = ISYS(JBL, JS)
ULE1_M(JV) = -VTI( 2,1)*VTI(JBL, JS)*DIJ(ILE1, J)
ULE2_M(JV) = -VTI( 2,2)*VTI(JBL, JS)*DIJ(ILE2, J)

```

```

          UTE1_M(JV) = -VTI( IBLTE(1),1)*VTI( JBL,JS)*DIJ( ITE1,J)
          UTE2_M(JV) = -VTI( IBLTE(2),2)*VTI( JBL,JS)*DIJ( ITE2,J)
110    CONTINUE
10    CONTINUE
C
      ULE1_A = UINV_A(2,1)
      ULE2_A = UINV_A(2,2)
C
C**** Go over each boundary layer/wake
      DO 2000 IS=1, 2
C
C---- there is no station "1" at similarity, so zero everything out
      DO 20 JS=1, 2
        DO 210 JBL=2, NBL(JS)
          JV = ISYS( JBL,JS)
          U1_M(JV) = 0.
          D1_M(JV) = 0.
210    CONTINUE
20    CONTINUE
      U1_A = 0.
      D1_A = 0.
C
      DUE1 = 0.
      DDS1 = 0.
C
C---- similarity station pressure gradient parameter  x/u du/dx
      IBL = 2
      BULE = 1.0
C
C---- set forced transition arc length position
      CALL XIFSET(IS)
C
      TRAN = .FALSE.
      TURB = .FALSE.
C
C**** Sweep downstream setting up BL equation linearizations
      DO 1000 IBL=2, NBL(IS)
C
      IV = ISYS( IBL,IS)
C
      SIMI = IBL.EQ.2
      WAKE = IBL.GT.IBLTE(IS)
      TRAN = IBL.EQ.ITRAN(IS)
      TURB = IBL.GT.ITRAN(IS)
C
      I = IPAN( IBL,IS)
C
C---- set primary variables for current station
      XSI = XSSI( IBL,IS)
      IF( IBL.LT.ITRAN(IS)) AMI = CTAU( IBL,IS)
      IF( IBL.GE.ITRAN(IS)) CTI = CTAU( IBL,IS)
      UEI = UEDG( IBL,IS)
      THI = THET( IBL,IS)
      MDI = MASS( IBL,IS)
C
      DSI = MDI/UEI
C
      IF( WAKE) THEN
        IW = IBL - IBLTE(IS)
        DSWAKI = WGAP( IW)
      ELSE
        DSWAKI = 0.
      ENDIF
C
C---- set derivatives of DSI (= D2)
      D2_M2 = 1.0/UEI
      D2_U2 = -DSI/UEI
C
      DO 30 JS=1, 2
        DO 310 JBL=2, NBL(JS)
          J = IPAN( JBL,JS)
          JV = ISYS( JBL,JS)

```

```

          U2_M(JV) = -VTI( IBL, IS) *VTI( JBL, JS) *DIJ( I, J)
          D2_M(JV) = D2_U2*U2_M(JV)
310    CONTINUE
30    CONTINUE
      D2_M(IV) = D2_M(IV) + D2_M2
C
      U2_A = UINV_A( IBL, IS)
      D2_A = D2_U2*U2_A
C
C----- "forced" changes due to mismatch between UEDG and USAV=UINV+dij*MASS
      DUE2 = UEDG( IBL, IS) - USAV( IBL, IS)
      DDS2 = D2_U2*DUE2
C
      CALL BLPRV( XSI, AMI, CTI, THI, DSI, DSWAKI, UEI)
      CALL BLKIN
C
C----- check for transition and set TRAN, XT, etc. if found
      IF(TRAN) THEN
          CALL TRCHEK
          AMI = AMPL2
      ENDIF
      IF( IBL.EQ.ITRAN( IS) .AND. .NOT.TRAN) THEN
          WRITE(*,*) 'SETBL: Xtr??? n1 n2: ', AMPL1, AMPL2
      ENDIF
C
C----- assemble 10x4 linearized system for dCtau, dTh, dDs, dUe, dXi
C      at the previous "1" station and the current "2" station
C
      IF( IBL.EQ.IBLTE( IS)+1) THEN
C
C----- define quantities at start of wake, adding TE base thickness to Dstar
      TTE = THET( IBLTE(1),1) + THET( IBLTE(2),2)
      DTE = DSTR( IBLTE(1),1) + DSTR( IBLTE(2),2) + ANTE
      CTE = ( CTAU( IBLTE(1),1) *THET( IBLTE(1),1)
&          + CTAU( IBLTE(2),2) *THET( IBLTE(2),2) ) / TTE
      CALL TESYS( CTE, TTE, DTE)
C
      TTE_TTE1 = 1.0
      TTE_TTE2 = 1.0
      DTE_MTE1 = 1.0 / UEDG( IBLTE(1),1)
      DTE_UTE1 = -DSTR( IBLTE(1),1) / UEDG( IBLTE(1),1)
      DTE_MTE2 = 1.0 / UEDG( IBLTE(2),2)
      DTE_UTE2 = -DSTR( IBLTE(2),2) / UEDG( IBLTE(2),2)
      CTE_CTE1 = THET( IBLTE(1),1) / TTE
      CTE_CTE2 = THET( IBLTE(2),2) / TTE
      CTE_TTE1 = (CTAU( IBLTE(1),1) - CTE) / TTE
      CTE_TTE2 = (CTAU( IBLTE(2),2) - CTE) / TTE
C
C----- re-define D1 sensitivities wrt m since D1 depends on both TE Ds values
      DO 35 JS=1, 2
          DO 350 JBL=2, NBL( JS)
              J = IPAN( JBL, JS)
              JV = ISYS( JBL, JS)
              D1_M(JV) = DTE_UTE1*UTE1_M(JV) + DTE_UTE2*UTE2_M(JV)
350    CONTINUE
35    CONTINUE
      D1_M(JVTE1) = D1_M(JVTE1) + DTE_MTE1
      D1_M(JVTE2) = D1_M(JVTE2) + DTE_MTE2
C
C----- "forced" changes from UEDG --- USAV=UINV+dij*MASS mismatch
      DUE1 = 0.
      DDS1 = DTE_UTE1*(UEDG( IBLTE(1),1) - USAV( IBLTE(1),1))
&          + DTE_UTE2*(UEDG( IBLTE(2),2) - USAV( IBLTE(2),2))
C
      ELSE
C
      CALL BLSYS
C
      ENDIF
C
C
C----- Save wall shear and equil. max shear coefficient for plotting output

```

```

        TAU( IBL, IS ) = 0.5*R2*U2*U2*CF2
        DIS( IBL, IS ) =      R2*U2*U2*U2*DI2*HS2*0.5
        CTQ( IBL, IS ) = CQ2
        DELT( IBL, IS ) = DE2
        USLP( IBL, IS ) = 1.60/(1.0+US2)

C
C---- set XI sensitivities wrt LE Ue changes
      IF( IS.EQ.1 ) THEN
        XI_ULE1 = SST_GO
        XI_ULE2 = -SST_GP
      ELSE
        XI_ULE1 = -SST_GO
        XI_ULE2 = SST_GP
      ENDIF

C
C---- stuff BL system coefficients into main Jacobian matrix
C
      DO 40 JV=1, NSYS
        VM(1,JV,IV) = VS1(1,3)*D1_M(JV) + VS1(1,4)*U1_M(JV)
&      + VS2(1,3)*D2_M(JV) + VS2(1,4)*U2_M(JV)
&      + (VS1(1,5) + VS2(1,5) + VSX(1))
&      * (XI_ULE1*ULE1_M(JV) + XI_ULE2*ULE2_M(JV))
40 CONTINUE

C
      VB(1,1,IV) = VS1(1,1)
      VB(1,2,IV) = VS1(1,2)

C
      VA(1,1,IV) = VS2(1,1)
      VA(1,2,IV) = VS2(1,2)

C
      IF( LALFA ) THEN
        VDEL(1,2,IV) = VSR(1)*RE_CLMR + VSM(1)*MSQ_CLMR
      ELSE
        VDEL(1,2,IV) =
&      (VS1(1,4)*U1_A + VS1(1,3)*D1_A)
&      + (VS2(1,4)*U2_A + VS2(1,3)*D2_A)
&      + (VS1(1,5) + VS2(1,5) + VSX(1))
&      * (XI_ULE1*ULE1_A + XI_ULE2*ULE2_A)
      ENDIF

C
      VDEL(1,1,IV) = VSREZ(1)
&      + (VS1(1,4)*DUE1 + VS1(1,3)*DDS1)
&      + (VS2(1,4)*DUE2 + VS2(1,3)*DDS2)
&      + (VS1(1,5) + VS2(1,5) + VSX(1))
&      * (XI_ULE1*DULE1 + XI_ULE2*DULE2)

C
C
      DO 50 JV=1, NSYS
        VM(2,JV,IV) = VS1(2,3)*D1_M(JV) + VS1(2,4)*U1_M(JV)
&      + VS2(2,3)*D2_M(JV) + VS2(2,4)*U2_M(JV)
&      + (VS1(2,5) + VS2(2,5) + VSX(2))
&      * (XI_ULE1*ULE1_M(JV) + XI_ULE2*ULE2_M(JV))
50 CONTINUE

C
      VB(2,1,IV) = VS1(2,1)
      VB(2,2,IV) = VS1(2,2)

C
      VA(2,1,IV) = VS2(2,1)
      VA(2,2,IV) = VS2(2,2)

C
      IF( LALFA ) THEN
        VDEL(2,2,IV) = VSR(2)*RE_CLMR + VSM(2)*MSQ_CLMR
      ELSE
        VDEL(2,2,IV) =
&      (VS1(2,4)*U1_A + VS1(2,3)*D1_A)
&      + (VS2(2,4)*U2_A + VS2(2,3)*D2_A)
&      + (VS1(2,5) + VS2(2,5) + VSX(2))
&      * (XI_ULE1*ULE1_A + XI_ULE2*ULE2_A)
      ENDIF

C
      VDEL(2,1,IV) = VSREZ(2)
&      + (VS1(2,4)*DUE1 + VS1(2,3)*DDS1)

```

```

& + (VS2(2,4)*DUE2 + VS2(2,3)*DDS2)
& + (VS1(2,5) + VS2(2,5) + VSX(2))
& * (XI_ULE1*DULE1 + XI_ULE2*DULE2)
C
C
DO 60 JV=1, NSYS
  VM(3,JV,IV) = VS1(3,3)*D1_M(JV) + VS1(3,4)*U1_M(JV)
& + VS2(3,3)*D2_M(JV) + VS2(3,4)*U2_M(JV)
& + (VS1(3,5) + VS2(3,5) + VSX(3))
& * (XI_ULE1*ULE1_M(JV) + XI_ULE2*ULE2_M(JV))
60 CONTINUE
C
VB(3,1,IV) = VS1(3,1)
VB(3,2,IV) = VS1(3,2)
C
VA(3,1,IV) = VS2(3,1)
VA(3,2,IV) = VS2(3,2)
C
IF(LALFA) THEN
  VDEL(3,2,IV) = VSR(3)*RE_CLMR + VSM(3)*MSQ_CLMR
ELSE
  VDEL(3,2,IV) =
& (VS1(3,4)*U1_A + VS1(3,3)*D1_A)
& + (VS2(3,4)*U2_A + VS2(3,3)*D2_A)
& + (VS1(3,5) + VS2(3,5) + VSX(3))
& * (XI_ULE1*ULE1_A + XI_ULE2*ULE2_A)
ENDIF
C
VDEL(3,1,IV) = VSREZ(3)
& + (VS1(3,4)*DUE1 + VS1(3,3)*DDS1)
& + (VS2(3,4)*DUE2 + VS2(3,3)*DDS2)
& + (VS1(3,5) + VS2(3,5) + VSX(3))
& * (XI_ULE1*DULE1 + XI_ULE2*DULE2)
C
C
IF( IBL.EQ.IBLTE(IS)+1) THEN
C
C----- redefine coefficients for TTE, DTE, etc
VZ(1,1) = VS1(1,1)*CTE_CTE1
VZ(1,2) = VS1(1,1)*CTE_TTE1 + VS1(1,2)*TTE_TTE1
VB(1,1,IV) = VS1(1,1)*CTE_CTE2
VB(1,2,IV) = VS1(1,1)*CTE_TTE2 + VS1(1,2)*TTE_TTE2
C
VZ(2,1) = VS1(2,1)*CTE_CTE1
VZ(2,2) = VS1(2,1)*CTE_TTE1 + VS1(2,2)*TTE_TTE1
VB(2,1,IV) = VS1(2,1)*CTE_CTE2
VB(2,2,IV) = VS1(2,1)*CTE_TTE2 + VS1(2,2)*TTE_TTE2
C
VZ(3,1) = VS1(3,1)*CTE_CTE1
VZ(3,2) = VS1(3,1)*CTE_TTE1 + VS1(3,2)*TTE_TTE1
VB(3,1,IV) = VS1(3,1)*CTE_CTE2
VB(3,2,IV) = VS1(3,1)*CTE_TTE2 + VS1(3,2)*TTE_TTE2
C
ENDIF
C
C---- turbulent intervals will follow if currently at transition interval
IF(TRAN) THEN
  TURB = .TRUE.
C
C----- save transition location
ITRAN(IS) = IBL
TFORCE(IS) = TRFORC
XSSITR(IS) = XT
C
C----- interpolate airfoil geometry to find transition x/c
C- (for user output)
IF(IS.EQ.1) THEN
  STR = SST - XT
ELSE
  STR = SST + XT
ENDIF
CHX = XTE - XLE

```

```

        CHY = YTE - YLE
        CHSQ = CHX**2 + CHY**2
        XTR = SEVAL(STR,X,XP,S,N)
        YTR = SEVAL(STR,Y,YP,S,N)
        XOCTR(IS) = ((XTR-XLE)*CHX + (YTR-YLE)*CHY)/CHSQ
        YOCTR(IS) = ((YTR-YLE)*CHX - (XTR-XLE)*CHY)/CHSQ
    ENDIF
C
    TRAN = .FALSE.
C
    IF( IBL.EQ.IBLTE(IS) ) THEN
C----- set "2" variables at TE to wake correlations for next station
C
        TURB = .TRUE.
        WAKE = .TRUE.
        CALL BLVAR(3)
        CALL BLMID(3)
    ENDIF
C
    DO 80 JS=1, 2
        DO 810 JBL=2, NBL(JS)
            JV = ISYS(JBL,JS)
            U1_M(JV) = U2_M(JV)
            D1_M(JV) = D2_M(JV)
810     CONTINUE
        80 CONTINUE
C
        U1_A = U2_A
        D1_A = D2_A
C
        DUE1 = DUE2
        DDS1 = DDS2
C
C----- set BL variables for next station
        DO 190 ICOM=1, NCOM
            COM1(ICOM) = COM2(ICOM)
        190 CONTINUE
C
C----- next streamwise station
        1000 CONTINUE
C
        IF(TFORCE(IS)) THEN
            WRITE(*,9100) IS,XOCTR(IS),ITRAN(IS)
9100     FORMAT(1X,'Side',I2,' forced transition at x/c = ',F7.4,I5)
        ELSE
            WRITE(*,9200) IS,XOCTR(IS),ITRAN(IS)
9200     FORMAT(1X,'Side',I2,' free transition at x/c = ',F7.4,I5)
        ENDIF
C
C----- next airfoil side
        2000 CONTINUE
C
        RETURN
    END

    SUBROUTINE IBLSYS
C-----
C    Sets the BL Newton system line number
C    corresponding to each BL station.
C-----
        INCLUDE 'XFOIL.INC'
        INCLUDE 'XBL.INC'
C
        IV = 0
        DO 10 IS=1, 2
            DO 110 IBL=2, NBL(IS)
                IV = IV+1
                ISYS(IBL,IS) = IV
            110 CONTINUE
        10 CONTINUE
C

```

```

      NSYS = IV
      IF(NSYS.GT.2*IVX) STOP '*** IBLSYS: BL system array overflow. ***'
C
      RETURN
      END

      SUBROUTINE MRCHUE
C-----
C      Marches the BLs and wake in direct mode using
C      the UEDG array. If separation is encountered,
C      a plausible value of Hk extrapolated from
C      upstream is prescribed instead. Continuous
C      checking of transition onset is performed.
C-----
      INCLUDE 'XFOIL.INC'
      INCLUDE 'XBL.INC'
      LOGICAL DIRECT
      REAL MSQ
C
C---- shape parameters for separation criteria
      HLMAX = 3.8
      HTMAX = 2.5
C
      DO 2000 IS=1, 2
C
        WRITE(*,*) '   side ', IS, ' ...'
C
C---- set forced transition arc length position
        CALL XIFSET(IS)
C
C---- initialize similarity station with Thwaites' formula
        IBL = 2
        XSI = XSSI(IBL,IS)
        UEI = UEDG(IBL,IS)
C
        BULE = LOG(UEDG(IBL+1,IS)/UEI) / LOG(XSSI(IBL+1,IS)/XSI)
C
        BULE = MAX( -.08 , BULE )
        BULE = 1.0
        UCON = UEI/XSI**BULE
        TSQ = 0.45/(UCON*(5.0*BULE+1.0)*REYBL) * XSI**(1.0-BULE)
        THI = SQRT(TSQ)
        DSI = 2.2*THI
        AMI = 0.0
C
C---- initialize Ctau for first turbulent station
        CTI = 0.03
C
        TRAN = .FALSE.
        TURB = .FALSE.
        ITRAN(IS) = IBLTE(IS)
C
C---- march downstream
        DO 1000 IBL=2, NBL(IS)
            IBM = IBL-1
C
            IW = IBL - IBLTE(IS)
C
            SIMI = IBL.EQ.2
            WAKE = IBL.GT.IBLTE(IS)
C
C----- prescribed quantities
            XSI = XSSI(IBL,IS)
            UEI = UEDG(IBL,IS)
C
            IF(WAKE) THEN
                IW = IBL - IBLTE(IS)
                DSWAKI = WGAP(IW)
            ELSE
                DSWAKI = 0.
            ENDIF
C
            DIRECT = .TRUE.

```



```

C
C----- Newton iteration loop for current station
          DO 100 ITBL=1, 25
C
C----- assemble 10x3 linearized system for dCtau, dTh, dDs, dUe, dXi
C          at the previous "1" station and the current "2" station
C          (the "1" station coefficients will be ignored)
C
C          CALL BLPRV(XSI,AMI,CTI,THI,DSI,DSWAKI,UEI)
          CALL BLKIN
C
C----- check for transition and set appropriate flags and things
          IF((.NOT.SIMI) .AND. (.NOT.TURB)) THEN
            CALL TRCHEK
            AMI = AMPL2
C
C----- fixed BUG    MD 7 Jun 99
          IF (TRAN) THEN
            ITRAN(IS) = IBL
            IF (CTI.LE.0.0) THEN
              CTI = 0.03
              S2 = CTI
            ENDIF
          ELSE
            ITRAN(IS) = IBL+2
          ENDIF
C
C          ENDIF
C
C          IF (IBL.EQ.IBLTE(IS)+1) THEN
            TTE = THET (IBLTE(1),1) + THET (IBLTE(2),2)
            DTE = DSTR (IBLTE(1),1) + DSTR (IBLTE(2),2) + ANTE
            CTE = ( CTAU (IBLTE(1),1)*THET (IBLTE(1),1)
&                + CTAU (IBLTE(2),2)*THET (IBLTE(2),2) ) / TTE
            CALL TESYS (CTE,TTE,DTE)
          ELSE
            CALL BLSYS
          ENDIF
C
C          IF (DIRECT) THEN
C
C----- try direct mode (set dUe = 0 in currently empty 4th line)
          VS2(4,1) = 0.
          VS2(4,2) = 0.
          VS2(4,3) = 0.
          VS2(4,4) = 1.0
          VSREZ(4) = 0.
C
C----- solve Newton system for current "2" station
          CALL GAUSS(4,4,VS2,VSREZ,1)
C
C----- determine max changes and underrelax if necessary
          DMAX = MAX( ABS(VSREZ(2)/THI),
&                  ABS(VSREZ(3)/DSI) )
          IF (IBL.LT.ITRAN(IS)) DMAX = MAX(DMAX,ABS(VSREZ(1)/10.0))
          IF (IBL.GE.ITRAN(IS)) DMAX = MAX(DMAX,ABS(VSREZ(1)/CTI) )
C
          RLX = 1.0
          IF (DMAX.GT.0.3) RLX = 0.3/DMAX
C
C----- see if direct mode is not applicable
          IF (IBL .NE. IBLTE(IS)+1) THEN
C
C----- calculate resulting kinematic shape parameter Hk
          MSQ = UEI*UEI*HSTINV / (GM1BL*(1.0 - 0.5*UEI*UEI*HSTINV))
          HTEST = (DSI + RLX*VSREZ(3)) / (THI + RLX*VSREZ(2))
          CALL HKIN( HTEST, MSQ, HKTEST, DUMMY, DUMMY)
C
C----- decide whether to do direct or inverse problem based on Hk
          IF (IBL.LT.ITRAN(IS)) HMAX = HLMAX

```

```

        IF (IBL.GE.ITRAN(IS)) HMAX = HTMAX
        DIRECT = HKTEST.LT.HMAX
    ENDIF

C
    IF (DIRECT) THEN
C----- update as usual
ccc        IF (IBL.LT.ITRAN(IS)) AMI = AMI + RLX*VSREZ(1)
            IF (IBL.GE.ITRAN(IS)) CTI = CTI + RLX*VSREZ(1)
            THI = THI + RLX*VSREZ(2)
            DSI = DSI + RLX*VSREZ(3)
        ELSE
C----- set prescribed Hk for inverse calculation at the current station
            IF (IBL.LT.ITRAN(IS)) THEN
C----- laminar case: relatively slow increase in Hk downstream
                HTARG = HK1 + 0.03*(X2-X1)/T1
            ELSE IF (IBL.EQ.ITRAN(IS)) THEN
C----- transition interval: weighted laminar and turbulent case
                HTARG = HK1 + (0.03*(XT-X1) - 0.15*(X2-XT))/T1
            ELSE IF (WAKE) THEN
C----- turbulent wake case:
                asymptotic wake behavior with approximate Backward Euler
                CONST = 0.03*(X2-X1)/T1
                HK2 = HK1
                HK2 = HK2 - (HK2 + CONST*(HK2-1.0)**3 - HK1)
                & / (1.0 + 3.0*CONST*(HK2-1.0)**2)
                HK2 = HK2 - (HK2 + CONST*(HK2-1.0)**3 - HK1)
                & / (1.0 + 3.0*CONST*(HK2-1.0)**2)
                HK2 = HK2 - (HK2 + CONST*(HK2-1.0)**3 - HK1)
                & / (1.0 + 3.0*CONST*(HK2-1.0)**2)
                HTARG = HK2
            ELSE
C----- turbulent case: relatively fast decrease in Hk downstream
                HTARG = HK1 - 0.15*(X2-X1)/T1
            ENDIF
C
C----- limit specified Hk to something reasonable
            IF (WAKE) THEN
                HTARG = MAX( HTARG , 1.01 )
            ELSE
                HTARG = MAX( HTARG , HMAX )
            ENDIF
C
        WRITE(*,1300) IBL, HTARG
1300        FORMAT(' MRCHUE: Inverse mode at', I4, '      Hk =', F8.3)
C
C----- try again with prescribed Hk
        GO TO 100
C
    ENDIF

C
    ELSE
C
C----- inverse mode (force Hk to prescribed value HTARG)
        VS2(4,1) = 0.
        VS2(4,2) = HK2_T2
        VS2(4,3) = HK2_D2
        VS2(4,4) = HK2_U2
        VSREZ(4) = HTARG - HK2
C
        CALL GAUSS(4,4,VS2,VSREZ,1)
C
        DMAX = MAX( ABS(VSREZ(2)/THI),
                & ABS(VSREZ(3)/DSI) )
        IF (IBL.GE.ITRAN(IS)) DMAX = MAX( DMAX , ABS(VSREZ(1)/CTI))
C
        RLX = 1.0
        IF (DMAX.GT.0.3) RLX = 0.3/DMAX
C
C----- update variables
ccc        IF (IBL.LT.ITRAN(IS)) AMI = AMI + RLX*VSREZ(1)
            IF (IBL.GE.ITRAN(IS)) CTI = CTI + RLX*VSREZ(1)
            THI = THI + RLX*VSREZ(2)

```

```

        DSI = DSI + RLX*VSREZ(3)
        UEI = UEI + RLX*VSREZ(4)
C
        ENDIF
C
C----- eliminate absurd transients
        IF( IBL.GE.ITRAN(IS) ) THEN
            CTI = MIN(CTI , 0.30 )
            CTI = MAX(CTI , 0.0000001 )
        ENDIF
C
        IF( IBL.LE.IBLTE(IS) ) THEN
            HKLIM = 1.02
        ELSE
            HKLIM = 1.00005
        ENDIF
        MSQ = UEI*UEI*HSTINV / (GM1BL*(1.0 - 0.5*UEI*UEI*HSTINV))
        DSW = DSI - DSWAKI
        CALL DSLIM(DSW,THI,UEI,MSQ,HKLIM)
        DSI = DSW + DSWAKI
C
        IF(DMAX.LE.1.0E-5) GO TO 110
C
100    CONTINUE
        WRITE(*,1350) IBL, IS, DMAX
1350    FORMAT(' MRCHUE: Convergence failed at',I4,' side',I2,
&          ' Res =', E12.4)
C
C----- the current unconverged solution might still be reasonable...
CCC    IF(DMAX .LE. 0.1) GO TO 110
        IF(DMAX .LE. 0.1) GO TO 109
C
C----- the current solution is garbage --> extrapolate values instead
        IF( IBL.GT.3 ) THEN
            IF( IBL.LE.IBLTE(IS) ) THEN
                THI = THET(IBM,IS) * (XSSI( IBL,IS)/XSSI( IBM,IS))**0.5
                DSI = DSTR(IBM,IS) * (XSSI( IBL,IS)/XSSI( IBM,IS))**0.5
            ELSE IF( IBL.EQ.IBLTE(IS)+1 ) THEN
                CTI = CTE
                THI = TTE
                DSI = DTE
            ELSE
                THI = THET(IBM,IS)
                RATLEN = (XSSI( IBL,IS)-XSSI( IBM,IS)) / (10.0*DSTR(IBM,IS))
                DSI = (DSTR(IBM,IS) + THI*RATLEN) / (1.0 + RATLEN)
            ENDIF
            IF( IBL.EQ.ITRAN(IS) ) CTI = 0.05
            IF( IBL.GT.ITRAN(IS) ) CTI = CTAU(IBM,IS)
C
                UEI = UEDG( IBL,IS)
                IF( IBL.GT.2 .AND. IBL.LT.NBL(IS) )
&                UEI = 0.5*(UEDG( IBL-1,IS) + UEDG( IBL+1,IS) )
            ENDIF
C
109    CALL BLPRV(XSI,AMI,CTI,THI,DSI,DSWAKI,UEI)
        CALL BLKIN
C
C----- check for transition and set appropriate flags and things
        IF((.NOT.SIMI) .AND. (.NOT.TURB)) THEN
            CALL TRCHEK
            AMI = AMPL2
            IF(      TRAN) ITRAN(IS) = IBL
            IF(.NOT.TRAN) ITRAN(IS) = IBL+2
        ENDIF
C
C----- set all other extrapolated values for current station
        IF( IBL.LT.ITRAN(IS) ) CALL BLVAR(1)
        IF( IBL.GE.ITRAN(IS) ) CALL BLVAR(2)
        IF(WAKE) CALL BLVAR(3)
C
        IF( IBL.LT.ITRAN(IS) ) CALL BLMID(1)
        IF( IBL.GE.ITRAN(IS) ) CALL BLMID(2)

```

```

                IF(WAKE) CALL BLMID(3)
C
C----- pick up here after the Newton iterations
110    CONTINUE
C
C----- store primary variables
        IF( IBL.LT.ITRAN(IS)) CTAU( IBL,IS) = AMI
        IF( IBL.GE.ITRAN(IS)) CTAU( IBL,IS) = CTI
        THET( IBL,IS) = THI
        DSTR( IBL,IS) = DSI
        UEDG( IBL,IS) = UEI
        MASS( IBL,IS) = DSI*UEI
        TAU( IBL,IS) = 0.5*R2*U2*U2*CF2
        DIS( IBL,IS) = R2*U2*U2*U2*DI2*HS2*0.5
        CTQ( IBL,IS) = CQ2
        DELT( IBL,IS) = DE2
C
C----- set "1" variables to "2" variables for next streamwise station
        CALL BLPRV(XSI,AMI,CTI,THI,DSI,DSWAKI,UEI)
        CALL BLKIN
        DO 310 ICOM=1, NCOM
            COM1( ICOM) = COM2( ICOM)
310    CONTINUE
C
C----- turbulent intervals will follow transition interval or TE
        IF(TRAN .OR. IBL.EQ.IBLTE(IS)) THEN
            TURB = .TRUE.
C
C----- save transition location
            TFORCE( IS) = TRFORC
            XSSITR( IS) = XT
            ENDIF
C
            TRAN = .FALSE.
C
            IF( IBL.EQ.IBLTE(IS)) THEN
                THI = THET( IBLTE(1),1) + THET( IBLTE(2),2)
                DSI = DSTR( IBLTE(1),1) + DSTR( IBLTE(2),2) + ANTE
            ENDIF
C
1000 CONTINUE
2000 CONTINUE
C
        RETURN
        END

        SUBROUTINE MRCHDU
C-----
C    Marches the BLs and wake in mixed mode using
C    the current Ue and Hk. The calculated Ue
C    and Hk lie along a line quasi-normal to the
C    natural Ue-Hk characteristic line of the
C    current BL so that the Goldstein or Levy-Lees
C    singularity is never encountered. Continuous
C    checking of transition onset is performed.
C-----
        INCLUDE 'XFOIL.INC'
        INCLUDE 'XBL.INC'
        REAL VTMP(4,5), VZTMP(4)
        REAL MSQ
ccc    REAL MDI
C
        DATA DEPS / 5.0E-6 /
C
C----- constant controlling how far Hk is allowed to deviate
C-    from the specified value.
        SENSWT = 1000.0
C
        DO 2000 IS=1, 2
C
C----- set forced transition arc length position

```

```

      CALL XIFSET(IS)
C
C----- set leading edge pressure gradient parameter  x/u du/dx
      IBL = 2
      XSI = XSSI(IBL,IS)
      UEI = UEDG(IBL,IS)
CCC      BULE = LOG(UEDG(IBL+1,IS)/UEI) / LOG(XSSI(IBL+1,IS)/XSI)
CCC      BULE = MAX( -.08 , BULE )
      BULE = 1.0
C
C----- old transition station
      ITROLD = ITRAN(IS)
C
      TRAN = .FALSE.
      TURB = .FALSE.
      ITRAN(IS) = IBLTE(IS)
C
C----- march downstream
      DO 1000 IBL=2, NBL(IS)
        IBM = IBL-1
C
        SIMI = IBL.EQ.2
        WAKE = IBL.GT.IBLTE(IS)
C
C----- initialize current station to existing variables
        XSI = XSSI(IBL,IS)
        UEI = UEDG(IBL,IS)
        THI = THET(IBL,IS)
        DSI = DSTR(IBL,IS)
CCC        MDI = MASS(IBL,IS)
C
C----- fixed BUG MD 7 June 99
        IF (IBL.LT.ITROLD) THEN
          AMI = CTAU(IBL,IS)
          CTI = 0.03
        ELSE
          CTI = CTAU(IBL,IS)
          IF (CTI.LE.0.0) CTI = 0.03
        ENDIF
C
CCC        DSI = MDI/UEI
C
        IF (WAKE) THEN
          IW = IBL - IBLTE(IS)
          DSWAKI = WGAP(IW)
        ELSE
          DSWAKI = 0.
        ENDIF
C
        IF (IBL.LE.IBLTE(IS)) DSI = MAX(DSI-DSWAKI,1.02000*THI) + DSWAKI
        IF (IBL.GT.IBLTE(IS)) DSI = MAX(DSI-DSWAKI,1.00005*THI) + DSWAKI
C
C----- Newton iteration loop for current station
      DO 100 ITBL=1, 25
C
C----- assemble 10x3 linearized system for dCtau, dTh, dDs, dUe, dXi
C      at the previous "1" station and the current "2" station
C      (the "1" station coefficients will be ignored)
C
      CALL BLPRV(XSI,AMI,CTI,THI,DSI,DSWAKI,UEI)
      CALL BLKIN
C
C----- check for transition and set appropriate flags and things
      IF ((.NOT.SIMI) .AND. (.NOT.TURB)) THEN
        CALL TRCHEK
        AMI = AMPL2
        IF (TRAN) ITRAN(IS) = IBL
        IF (.NOT.TRAN) ITRAN(IS) = IBL+2
      ENDIF
C
      IF (IBL.EQ.IBLTE(IS)+1) THEN

```

```

      TTE = THET( IBLTE(1),1) + THET( IBLTE(2),2)
      DTE = DSTR( IBLTE(1),1) + DSTR( IBLTE(2),2) + ANTE
      CTE = ( CTAU( IBLTE(1),1)*THET( IBLTE(1),1)
&          + CTAU( IBLTE(2),2)*THET( IBLTE(2),2) ) / TTE
      CALL TESYS(CTE,TTE,DTE)
    ELSE
      CALL BLSYS
    ENDIF
  C
  C----- set stuff at first iteration...
  IF( ITBL.EQ.1) THEN
  C
  C----- set "baseline" Ue and Hk for forming Ue(Hk) relation
  UEREF = U2
  HKREF = HK2
  C
  C----- if current point IBL was turbulent and is now laminar, then...
  IF( IBL.LT.ITRAN( IS) .AND. IBL.GE.ITROLD ) THEN
  C----- extrapolate baseline Hk
  UEM = UEDG( IBL-1, IS)
  DSM = DSTR( IBL-1, IS)
  THM = THET( IBL-1, IS)
  MSQ = UEM*UEM*HSTINV / ( GM1BL*(1.0 - 0.5*UEM*UEM*HSTINV))
  CALL HKIN( DSM/THM, MSQ, HKREF, DUMMY, DUMMY )
  ENDIF
  C
  C----- if current point IBL was laminar, then...
  IF( IBL.LT.ITROLD) THEN
  C----- reinitialize or extrapolate Ctau if it's now turbulent
  IF(TRAN) CTAU( IBL, IS) = 0.03
  IF(TURB) CTAU( IBL, IS) = CTAU( IBL-1, IS)
  IF(TRAN .OR. TURB) THEN
    CTI = CTAU( IBL, IS)
    S2 = CTI
  ENDIF
  ENDIF
  C
  ENDIF
  C
  C
  IF( SIMI .OR. IBL.EQ.IBLTE( IS)+1) THEN
  C
  C----- for similarity station or first wake point, prescribe Ue
  VS2(4,1) = 0.
  VS2(4,2) = 0.
  VS2(4,3) = 0.
  VS2(4,4) = U2_UEI
  VSREZ(4) = UEREF - U2
  C
  ELSE
  C
  C***** calculate Ue-Hk characteristic slope
  C
  DO 20 K=1, 4
    VZTMP( K) = VSREZ( K)
    DO 201 L=1, 5
      VTMP( K,L) = VS2( K,L)
    201 CONTINUE
    20 CONTINUE
  C
  C----- set unit dHk
  VTMP(4,1) = 0.
  VTMP(4,2) = HK2_T2
  VTMP(4,3) = HK2_D2
  VTMP(4,4) = HK2_U2*U2_UEI
  VZTMP(4) = 1.0
  C
  C----- calculate dUe response
  CALL GAUSS(4,4,VTMP,VZTMP,1)
  C
  C----- set SENSWT * (normalized dUe/dHk)
  SENNEW = SENSWT * VZTMP(4) * HKREF/UEREF

```

```

        IF(ITBL.LE.5) THEN
            SENS = SENNEW
        ELSE IF(ITBL.LE.15) THEN
            SENS = 0.5*(SENS + SENNEW)
        ENDIF

C
C----- set prescribed Ue-Hk combination
VS2(4,1) = 0.
VS2(4,2) = HK2_T2 * HKREF
VS2(4,3) = HK2_D2 * HKREF
VS2(4,4) = ( HK2_U2 * HKREF + SENS/UREF ) * U2_UEI
VSREZ(4) = -(HKREF**2)*(HK2 / HKREF - 1.0)
&
        - SENS*(U2 / UREF - 1.0)

C
        ENDIF

C
C----- solve Newton system for current "2" station
CALL GAUSS(4,4,VS2,VSREZ,1)

C
C----- determine max changes and underrelax if necessary
DMAX = MAX( ABS(VSREZ(2)/THI),
&
        ABS(VSREZ(3)/DSI) )
IF( IBL.GE.ITRAN(IS) ) DMAX = MAX(DMAX,ABS(VSREZ(1)/(10.0*CTI)))

C
        RLX = 1.0
        IF(DMAX.GT.0.3) RLX = 0.3/DMAX

C
C----- update as usual
IF( IBL.LT.ITRAN(IS) ) AMI = AMI + RLX*VSREZ(1)
IF( IBL.GE.ITRAN(IS) ) CTI = CTI + RLX*VSREZ(1)
THI = THI + RLX*VSREZ(2)
DSI = DSI + RLX*VSREZ(3)
UEI = UEI + RLX*VSREZ(4)

C
C----- eliminate absurd transients
IF( IBL.GE.ITRAN(IS) ) THEN
        CTI = MIN(CTI , 0.30 )
        CTI = MAX(CTI , 0.0000001 )
ENDIF

C
        IF( IBL.LE.IBLTE(IS) ) THEN
            HKLIM = 1.02
        ELSE
            HKLIM = 1.00005
        ENDIF
        MSQ = UEI*UEI*HSTINV / (GM1BL*(1.0 - 0.5*UEI*UEI*HSTINV))
        DSW = DSI - DSWAKI
        CALL DSLIM(DSW,THI,UEI,MSQ,HKLIM)
        DSI = DSW + DSWAKI

C
        IF(DMAX.LE.DEPS) GO TO 110

C
100  CONTINUE

C
        WRITE(*,1350) IBL, IS, DMAX
1350  FORMAT(' MRCHDU: Convergence failed at',I4,' side',I2,
&
        ' Res =', E12.4)

C
C----- the current unconverged solution might still be reasonable...
CCC  IF(DMAX .LE. 0.1) GO TO 110
        IF(DMAX .LE. 0.1) GO TO 109

C
C----- the current solution is garbage --> extrapolate values instead
IF( IBL.GT.3 ) THEN
        IF( IBL.LE.IBLTE(IS) ) THEN
            THI = THET(IBM,IS) * (XSSI( IBL,IS )/XSSI( IBM,IS ))**0.5
            DSI = DSTR(IBM,IS) * (XSSI( IBL,IS )/XSSI( IBM,IS ))**0.5
            UEI = UEDG(IBM,IS)
        ELSE IF( IBL.EQ.IBLTE(IS)+1 ) THEN
            CTI = CTE
            THI = TTE
            DSI = DTE

```

```

        UEI = UEDG(IBM,IS)
    ELSE
        THI = THET(IBM,IS)
        RATLEN = (XSSI(IBM,IS)-XSSI(IBM,IS)) / (10.0*DSTR(IBM,IS))
        DSI = (DSTR(IBM,IS) + THI*RATLEN) / (1.0 + RATLEN)
        UEI = UEDG(IBM,IS)
    ENDIF
    IF(IBM.EQ.ITRAN(IS)) CTI = 0.05
    IF(IBM.GT.ITRAN(IS)) CTI = CTAU(IBM,IS)
ENDIF
C
109    CALL BLPRV(XSI,AMI,CTI,THI,DSI,DSWAKI,UEI)
    CALL BLKIN
C
C----- check for transition and set appropriate flags and things
IF((.NOT.SIMI) .AND. (.NOT.TURB)) THEN
    CALL TRCHEK
    AMI = AMPL2
    IF(      TRAN) ITRAN(IS) = IBL
    IF(.NOT.TRAN) ITRAN(IS) = IBL+2
ENDIF
C
C----- set all other extrapolated values for current station
IF(IBM.LT.ITRAN(IS)) CALL BLVAR(1)
IF(IBM.GE.ITRAN(IS)) CALL BLVAR(2)
IF(WAKE) CALL BLVAR(3)
C
    IF(IBM.LT.ITRAN(IS)) CALL BLMID(1)
    IF(IBM.GE.ITRAN(IS)) CALL BLMID(2)
    IF(WAKE) CALL BLMID(3)
C
C----- pick up here after the Newton iterations
110    CONTINUE
C
    SENS = SENNEW
C
C----- store primary variables
IF(IBM.LT.ITRAN(IS)) CTAU(IBM,IS) = AMI
IF(IBM.GE.ITRAN(IS)) CTAU(IBM,IS) = CTI
THET(IBM,IS) = THI
DSTR(IBM,IS) = DSI
UEDG(IBM,IS) = UEI
MASS(IBM,IS) = DSI*UEI
TAU(IBM,IS) = 0.5*R2*U2*U2*CF2
DIS(IBM,IS) = R2*U2*U2*U2*DI2*HS2*0.5
CTQ(IBM,IS) = CQ2
C
C----- set "1" variables to "2" variables for next streamwise station
CALL BLPRV(XSI,AMI,CTI,THI,DSI,DSWAKI,UEI)
CALL BLKIN
DO 310 ICOM=1, NCOM
    COM1(ICOM) = COM2(ICOM)
310    CONTINUE
C
C
C----- turbulent intervals will follow transition interval or TE
IF(TRAN .OR. IBL.EQ.IBLTE(IS)) THEN
    TURB = .TRUE.
C
C----- save transition location
TFORCE(IS) = TRFORC
XSSITR(IS) = XT
ENDIF
C
    TRAN = .FALSE.
C
1000 CONTINUE
C
2000 CONTINUE
C
    RETURN
END

```



```

      SUBROUTINE XIFSET(IS)
C-----
C      Sets forced-transition BL coordinate locations.
C-----
      INCLUDE 'XFOIL.INC'
      INCLUDE 'XBL.INC'

C      IF(XSTRIP(IS).GE.1.0) THEN
         XIFORC = XSSI(IBLTE(IS),IS)
         RETURN
      ENDIF

C      CHX = XTE - XLE
      CHY = YTE - YLE
      CHSQ = CHX**2 + CHY**2

C
C----- calculate chord-based x/c, y/c
      DO 10 I=1, N
         W1(I) = ((X(I)-XLE)*CHX + (Y(I)-YLE)*CHY) / CHSQ
         W2(I) = ((Y(I)-YLE)*CHX - (X(I)-XLE)*CHY) / CHSQ
      10 CONTINUE

C      CALL SPLIND(W1,W3,S,N,-999.0,-999.0)
      CALL SPLIND(W2,W4,S,N,-999.0,-999.0)

C      IF(IS.EQ.1) THEN

C
C----- set approximate arc length of forced transition point for SINVRT
      STR = SLE + (S(1)-SLE)*XSTRIP(IS)

C
C----- calculate actual arc length
      CALL SINVRT(STR,XSTRIP(IS),W1,W3,S,N)

C
C----- set BL coordinate value
      XIFORC = MIN( (SST - STR) , XSSI(IBLTE(IS),IS) )

C
      ELSE

C----- same for bottom side
C
      STR = SLE + (S(N)-SLE)*XSTRIP(IS)
      CALL SINVRT(STR,XSTRIP(IS),W1,W3,S,N)
      XIFORC = MIN( (STR - SST) , XSSI(IBLTE(IS),IS) )

C
      ENDIF

C      IF(XIFORC .LT. 0.0) THEN
         WRITE(*,1000) IS
      1000 FORMAT(' *** Stagnation point is past trip on side',I2,' ***')
         XIFORC = XSSI(IBLTE(IS),IS)
      ENDIF

C      RETURN
      END

      SUBROUTINE UPDATE(sec,pcor)
C-----
C      Adds on Newton deltas to boundary layer variables.
C      Checks for excessive changes and underrelaxes if necessary.
C      Calculates max and rms changes.
C      Also calculates the change in the global variable "AC".
C      If LALFA=.TRUE. , "AC" is CL
C      If LALFA=.FALSE., "AC" is alpha
C-----
      INCLUDE 'XFOIL.INC'
      include 'section.inc' !02/06/2005

      type(section):: sec

```

```

      REAL UNEW(IVX,2), U_AC(IVX,2)
      real:: pcor(1qx,2) !mass defect correction (22/05/2005)
      REAL:: QNEW(IQX), Q_AC(IQX), Qcorr(IQX) !viscous correction
      EQUIVALENCE (VA(1,1,1), UNEW(1,1)) ,
&                (VB(1,1,1), QNEW(1) )
      EQUIVALENCE (VA(1,1,IVX), U_AC(1,1)) ,
&                (VB(1,1,IVX), Q_AC(1) )
      REAL MSQ

C
C---- max allowable alpha changes per iteration
      DALMAX = 0.5*DTOR
      DALMIN = -0.5*DTOR

C
C---- max allowable CL change per iteration
      DCLMAX = 0.5
      DCLMIN = -0.5
      IF (MATYP.NE.1) DCLMIN = MAX(-0.5 , -0.9*CL)

C
      HSTINV = GAMM1*(MINF/QINF)**2 / (1.0 + 0.5*GAMM1*MINF**2)

C
C---- calculate new Ue distribution assuming no under-relaxation
C- also set the sensitivity of Ue wrt to alpha or Re
      DO 1 IS=1, 2
        DO 10 IBL=2, NBL(IS)
          I = IPAN(IBL,IS)

C
          DUI = 0.
          DUI_AC = 0.
          DO 100 JS=1, 2
            DO 1000 JBL=2, NBL(JS)
              J = IPAN(JBL,JS)
              JV = ISYS(JBL,JS)
              UE_M = -VTI(IBL,IS)*VTI(JBL,JS)*DIJ(I,J)
              DUI = DUI + UE_M*(MASS(JBL,JS)+VDEL(3,1,JV))
              DUI_AC = DUI_AC + UE_M*( -VDEL(3,2,JV))
            1000 CONTINUE
          100 CONTINUE

C
          UINV depends on "AC" only if "AC" is alpha
          IF (LALFA) THEN
            UINV_AC = 0.
          ELSE
            UINV_AC = UINV_A(IBL,IS)
          ENDIF

C
          pcor(1bl,1s) = DUI !viscous correction vector (22/05/2005)
          UNEW(IBL,IS) = UINV(IBL,IS) + DUI
          U_AC(IBL,IS) = UINV_AC + DUI_AC

C
        10 CONTINUE
      1 CONTINUE

C
C---- set new Qtan from new Ue with appropriate sign change
      DO 2 IS=1, 2
        DO 20 IBL=2, IBLTE(IS)
          I = IPAN(IBL,IS)
          Qcorr(i) = VTI(IBL,IS)*pcor(1BL,1S) !added 02/06/2005
          QNEW(I) = VTI(IBL,IS)*UNEW(IBL,IS)
          Q_AC(I) = VTI(IBL,IS)*U_AC(IBL,IS)
        20 CONTINUE
      2 CONTINUE
      QNEW(IST) = 0. !correction on 30/03/2004

C
C---- calculate new CL from this new Qtan
      SA = SIN(ALFA)
      CA = COS(ALFA)

C
      BETA = SQRT(1.0 - MINF**2)
      BETA_MSQ = -0.5/BETA

C
      BFAC = 0.5*MINF**2 / (1.0 + BETA)
      BFAC_MSQ = 0.5 / (1.0 + BETA)

```

```

&          - BFAC          / (1.0 + BETA) * BETA_MSQ
C
CLNEW = 0.
CL_A  = 0.
CL_MS = 0.
CL_AC = 0.
C
I = 1
!CGINC = 1.0 - (QNEW(I)/QINF)**2
CGINC = sec.cp(i) + Qcorr(i)**2
CPG1  = CGINC / (BETA + BFAC*CGINC)
CPG1_MS = -CPG1/(BETA + BFAC*CGINC)*(BETA_MSQ + BFAC_MSQ*CGINC)
C
!CPI_Q = -2.0*QNEW(I)/QINF**2
CPI_Q = -2.0*(1-sec.cp(i))
CPC_CPI = (1.0 - BFAC*CPG1)/(BETA + BFAC*CGINC)
CPG1_AC = CPC_CPI*CPI_Q*Q_AC(I)
C
DO 3 I=1, N
  IP = I+1
  IF(I.EQ.N) IP = 1
C
  !CGINC = 1.0 - (QNEW(IP)/QINF)**2
  CGINC = sec.cp(i) + Qcorr(i)**2
  CPG2  = CGINC / (BETA + BFAC*CGINC)
  CPG2_MS = -CPG2/(BETA + BFAC*CGINC)*(BETA_MSQ + BFAC_MSQ*CGINC)
C
  CPI_Q = -2.0*(1-sec.cp(i))
  CPC_CPI = (1.0 - BFAC*CPG2)/(BETA + BFAC*CGINC)
  CPG2_AC = CPC_CPI*CPI_Q*Q_AC(IP)
C
  DX  = (X(IP) - X(I))*CA + (Y(IP) - Y(I))*SA
  DX_A = -(X(IP) - X(I))*SA + (Y(IP) - Y(I))*CA
C
  AG  = 0.5*(CPG2 + CPG1 )
  AG_MS = 0.5*(CPG2_MS + CPG1_MS)
  AG_AC = 0.5*(CPG2_AC + CPG1_AC)
C
  CLNEW = CLNEW + DX *AG
  CL_A  = CL_A  + DX_A*AG
  CL_MS = CL_MS + DX *AG_MS
  CL_AC = CL_AC + DX *AG_AC
C
  CPG1  = CPG2
  CPG1_MS = CPG2_MS
  CPG1_AC = CPG2_AC
3 CONTINUE
C
C---- initialize under-relaxation factor
RLX = 1.0
C
  IF(LALFA) THEN
C===== alpha is prescribed: AC is CL
C
C----- set change in Re to account for CL changing, since Re = Re(CL)
DAC = (CLNEW - CL) / (1.0 - CL_AC - CL_MS*2.0*MINF*MINF_CL)
C
C----- set under-relaxation factor if Re change is too large
IF(RLX*DAC .GT. DCLMAX) RLX = DCLMAX/DAC
IF(RLX*DAC .LT. DCLMIN) RLX = DCLMIN/DAC
C
  ELSE
C===== CL is prescribed: AC is alpha
C
C----- set change in alpha to drive CL to prescribed value
DAC = (CLNEW - CLSPEC) / (0.0 - CL_AC - CL_A)
C
C----- set under-relaxation factor if alpha change is too large
IF(RLX*DAC .GT. DALMAX) RLX = DALMAX/DAC
IF(RLX*DAC .LT. DALMIN) RLX = DALMIN/DAC
C
  ENDIF

```

```

C
    RMSBL = 0.
    RMXBL = 0.
C
    DHI = 1.5
    DLO = -.5
C
C---- calculate changes in BL variables and under-relaxation if needed
    DO 4 IS=1, 2
        DO 40 IBL=2, NBL(IS)
            IV = ISYS(IBL,IS)
C
C----- set changes without underrelaxation
            DCTAU = VDEL(1,1,IV) - DAC*VDEL(1,2,IV)
            DTHET = VDEL(2,1,IV) - DAC*VDEL(2,2,IV)
            DMASS = VDEL(3,1,IV) - DAC*VDEL(3,2,IV)
            DUEDG = UNEW(IBL,IS) + DAC*U_AC(IBL,IS) - UEDG(IBL,IS)
            DDSTR = (DMASS - DSTR(IBL,IS)*DUEDG)/UEDG(IBL,IS)
C
C----- normalize changes !all corrected 27/06/2004
            IF(IBL.LT.ITRAN(IS)) DN1 = DCTAU / 10.0
            IF(IBL.GE.ITRAN(IS)) DN1 = DCTAU / CTAU(IBL,IS)
            DN2 = DTHET / THET(IBL,IS)
            DN3 = DDSTR / DSTR(IBL,IS)
            DN4 = ABS(DUEDG)/0.25

            if (iv<=iblte(is)) then !just for body 30/03/2004
C
C----- accumulate for rms change
                RMSBL = RMSBL + DN1**2 + DN2**2 + DN3**2 + DN4**2
            endif
C
C----- see if Ctau needs underrelaxation
                RDN1 = RLX*DN1
                IF(ABS(DN1) .GT. ABS(RMXBL)) THEN
                    RMXBL = DN1
                    IF(IBL.LT.ITRAN(IS)) VMXBL = 'n'
                    IF(IBL.GE.ITRAN(IS)) VMXBL = 'C'
                    IMXBL = IBL
                    ISMXBL = IS
                ENDIF
                IF(RDN1 .GT. DHI) RLX = DHI/DN1
                IF(RDN1 .LT. DLO) RLX = DLO/DN1
C
C----- see if Theta needs underrelaxation
                RDN2 = RLX*DN2
                IF(ABS(DN2) .GT. ABS(RMXBL)) THEN
                    RMXBL = DN2
                    VMXBL = 'T'
                    IMXBL = IBL
                    ISMXBL = IS
                ENDIF
                IF(RDN2 .GT. DHI) RLX = DHI/DN2
                IF(RDN2 .LT. DLO) RLX = DLO/DN2
C
C----- see if Dstar needs underrelaxation
                RDN3 = RLX*DN3
                IF(ABS(DN3) .GT. ABS(RMXBL)) THEN
                    RMXBL = DN3
                    VMXBL = 'D'
                    IMXBL = IBL
                    ISMXBL = IS
                ENDIF
                IF(RDN3 .GT. DHI) RLX = DHI/DN3
                IF(RDN3 .LT. DLO) RLX = DLO/DN3
C
C----- see if Ue needs underrelaxation
                RDN4 = RLX*DN4
                IF(ABS(DN4) .GT. ABS(RMXBL)) THEN
                    RMXBL = DUEDG
                    VMXBL = 'U'
                    IMXBL = IBL

```

```

        ISMXBL = IS
        ENDIF
        IF(RDN4 .GT. DHI) RLX = DHI/DN4
        IF(RDN4 .LT. DLO) RLX = DLO/DN4
C
    40    CONTINUE
    4    CONTINUE
C
C---- set true rms change
    RMSBL = SQRT( RMSBL / (4.0*FLOAT( NBL(1)+NBL(2) )) ) !/1000 !14/04/2004
C
C
        IF(LALFA) THEN
C----- set underrelaxed change in Reynolds number from change in lift
            CL = CL + RLX*DAC
        ELSE
C----- set underrelaxed change in alpha
            ALFA = ALFA + RLX*DAC
            ADEG = ALFA/DTOR
        ENDIF
C
C---- update BL variables with underrelaxed changes
    DO 5 IS=1, 2
        DO 50 IBL=2, NBL(IS)
            IV = ISYS(IBL,IS)
C
            DCTAU = VDEL(1,1,IV) - DAC*VDEL(1,2,IV)
            DTHET = VDEL(2,1,IV) - DAC*VDEL(2,2,IV)
            DMASS = VDEL(3,1,IV) - DAC*VDEL(3,2,IV)
            DUEDG = UNEW(IBL,IS) + DAC*U_AC(IBL,IS) - UEDG(IBL,IS)
            DDSTR = (DMASS - DSTR(IBL,IS)*DUEDG)/UEDG(IBL,IS)
C
            CTAU(IBL,IS) = CTAU(IBL,IS) + RLX*DCTAU
            THET(IBL,IS) = THET(IBL,IS) + RLX*DTHET
            DSTR(IBL,IS) = DSTR(IBL,IS) + RLX*DDSTR
            UEDG(IBL,IS) = UEDG(IBL,IS) + RLX*DUEDG
C
            IF(IBL.GT.IBLTE(IS)) THEN
                IW = IBL - IBLTE(IS)
                DSWAKI = WGAP(IW)
            ELSE
                DSWAKI = 0.
            ENDIF
C
C----- eliminate absurd transients
            IF(IBL.GE.ITRAN(IS))
                & CTAU(IBL,IS) = MIN( CTAU(IBL,IS) , 0.25 )
C
            IF(IBL.LE.IBLTE(IS)) THEN
                HKLIM = 1.02
            ELSE
                HKLIM = 1.00005
            ENDIF
            MSQ = UEDG(IBL,IS)**2*HSTINV
            & / (GAMM1*(1.0 - 0.5*UEDG(IBL,IS)**2*HSTINV))
            DSW = DSTR(IBL,IS) - DSWAKI
            CALL DSLIM(DSW,THET(IBL,IS),UEDG(IBL,IS),MSQ,HKLIM)
            DSTR(IBL,IS) = DSW + DSWAKI
C
C----- set new mass defect (nonlinear update)
            MASS(IBL,IS) = DSTR(IBL,IS) * UEDG(IBL,IS)
C
    50    CONTINUE
    5    CONTINUE
C
C
C---- equate upper wake arrays to lower wake arrays
    DO 6 KBL=1, NBL(2)-IBLTE(2)
        CTAU(IBLTE(1)+KBL,1) = CTAU(IBLTE(2)+KBL,2)
        THET(IBLTE(1)+KBL,1) = THET(IBLTE(2)+KBL,2)
        DSTR(IBLTE(1)+KBL,1) = DSTR(IBLTE(2)+KBL,2)
        UEDG(IBLTE(1)+KBL,1) = UEDG(IBLTE(2)+KBL,2)
    
```

```

        TAU (IBLTE (1)+KBL,1) =   TAU (IBLTE (2)+KBL,2)
        DIS (IBLTE (1)+KBL,1) =   DIS (IBLTE (2)+KBL,2)
        CTQ (IBLTE (1)+KBL,1) =   CTQ (IBLTE (2)+KBL,2)
6      CONTINUE
C
      RETURN
      END

```

```

      SUBROUTINE DSLIM(DSTR,THET,UEDG,MSQ,HKLIM)
      IMPLICIT REAL (A-H,M,O-Z)
C
      H = DSTR/THET
      CALL HKIN(H,MSQ,HK,HK_H,HK_M)
C
      DH = MAX( 0.0 , HKLIM-HK ) / HK_H
      DSTR = DSTR + DH*THET
C
      RETURN
      END

```

## VIX Subroutine Stagpoint

```

!This subroutine finds the stagnation point of
! a interpolated section. If it is a membrane,
! fsharp = true, program will set the stagnation
! point at the very leading edge
!*****
! Created by: Augusto Veiga,
! FSIIG, University of Southampton 2003
!*****

      subroutine stagpoint(Up,s,gamma,x,N,Nw,Ist,SST,Sst_GO,SST_GP,
&   fsharp)
      integer:: N,Nw,IST,IS
      real:: cpmax,h
      real:: dcp1,dcp2
      real,dimension(N+Nw):: Up,s,x,gamma,cp
      logical:: fsharp

!Calculating Cp over the wing
do i = 1,N
      cp(i) = 1.-Up(i)**2
enddo
      if (fsharp) then
            ist = int(N/2)+1
            sst = s(ist)
            is = ist
      else
            !Getting biggest Cp and position
            cpmax = 0
            is = int(N/2)+1
            k = is
            i = is
            lp1: do
                  i = i+1
                  if (cp(i)>cpmax) then
                        cpmax = cp(i)
                        is = i           !finding a possible point
                        h = s(i)-s(i-1)
                        dcp1 = (cpmax-cp(i-1))/h
                        dcp2 = (cp(i+1)-cpmax)/h
                        if ((dcp1>0 .and. dcp2<0)) then
                              ist = is !testing derivatives
                              sst = s(is)
                              exit lp1
                        endif
                  endif
            else if (i>(k+int(N/2))) then
                  is = int(N/2)+1
                  sst = s(is)
                  exit lp1
            enddo

```

```

        endif
    enddo lp1
endif
!signal for gamma
gamma(ist) = 0
do i = 1,IST-1 !IST+1,N
    gamma(i) = Up(i)
enddo
!upper part
do i = IST+1,N
    gamma(i) = -Up(i)
enddo
!wake
do i = N+1,N+Nw
    gamma(i) = Up(i)
enddo
DGAM = GAMMA(IST+1) - GAMMA(IST)
SST_GO = (SST - S(Ist+1))/DGAM
SST_GP = (S(Ist+1) - SST)/DGAM
return
end subroutine

```

## VIX Subroutine AIJCALC

```

!*****
      SUBROUTINE AIJCALC
C-----
C    Calculates two surface vorticity (gamma) distributions
C    for alpha = 0, 90 degrees. These are superimposed
C    in SPECAL or SPECCL for specified alpha or CL.
C    This subroutine was adapted from XFOIL by Augusto Veiga
C-----
      INCLUDE 'XFOIL.INC'
C
C---- distance of internal control point ahead of sharp TE
C- (fraction of smaller panel length adjacent to TE)
      BWT = 0.1
C
      WRITE(*,*) 'Calculating unit vorticity distributions ...'
C
      DO 10 I=1, N
!          GAM(I) = 0.
          GAMU(I,1) = 0.
          GAMU(I,2) = 0.
      10 CONTINUE
      PSIO = 0.
C
C---- Set up matrix system for Psi = Psio on airfoil surface.
C- The unknowns are (dGamma)i and dPsio.
      DO 20 I=1, N
C
C----- calculate Psi and dPsi/dGamma array for current node
          CALL PSILIN(I,X(I),Y(I),NX(I),NY(I),PSI,PSI_N,.FALSE.,.TRUE.)
C
          PSIINF = QINF*(COS(ALFA)*Y(I) - SIN(ALFA)*X(I))
C
C----- RES1 = PSI( 0) - PSIO
C----- RES2 = PSI(90) - PSIO
          RES1 = QINF*Y(I)
          RES2 = -QINF*X(I)
C
C----- dRes/dGamma
          DO 201 J=1, N
              AIJ(I,J) = DZDG(J)
          201 CONTINUE
C
          DO 202 J=1, N
              BIJ(I,J) = -DZDM(J)
          202 CONTINUE
C
C----- dRes/dPsio

```

```

      AIJ(I,N+1) = -1.0
C
      GAMU(I,1) = -RES1
      GAMU(I,2) = -RES2
C
20 CONTINUE
C
C----- set Kutta condition
C-   RES = GAM(1) + GAM(N)
      RES = 0.
C
      DO 30 J=1, N+1
        AIJ(N+1,J) = 0.0
30 CONTINUE
C
      AIJ(N+1,1) = 1.0
      AIJ(N+1,N) = 1.0
C
      GAMU(N+1,1) = -RES
      GAMU(N+1,2) = -RES
C
C----- set up Kutta condition (no direct source influence)
      DO 32 J=1, N
        BIJ(N+1,J) = 0.
32 CONTINUE
C
      IF(SHARP) THEN
C----- set zero internal velocity in TE corner
C
C----- set TE bisector angle
      AG1 = ATAN2(-YP(1),-XP(1))
      AG2 = ATANC( YP(N), XP(N),AG1)
      ABIS = 0.5*(AG1+AG2)
      CBIS = COS(ABIS)
      SBIS = SIN(ABIS)
C
C----- minimum panel length adjacent to TE
      DS1 = SQRT( (X(1)-X(2))**2 + (Y(1)-Y(2))**2 )
      DS2 = SQRT( (X(N)-X(N-1))**2 + (Y(N)-Y(N-1))**2 )
      DSMIN = MIN( DS1 , DS2 )
C
C----- control point on bisector just ahead of TE point
      XBIS = XTE - BWT*DSMIN*CBIS
      YBIS = YTE - BWT*DSMIN*SBIS
ccc      write(*,*) xbis, ybis
C
C----- set velocity component along bisector line
      CALL PSILIN(0,XBIS,YBIS,-SBIS,CBIS,PSI,QBIS,.FALSE.,.TRUE.)
C
CCC--- RES = DQDGj*Gammaj + DQDMj*Massj + QINF*(COSA*CBIS + SINA*SBIS)
      RES = QBIS
C
C----- dRes/dGamma
      DO J=1, N
        AIJ(N,J) = DQDG(J)
      ENDDO
C
C----- -dRes/dMass
      DO J=1, N
        BIJ(N,J) = -DQDM(J)
      ENDDO
C
C----- dRes/dPsio
      AIJ(N,N+1) = 0.
C
C----- -dRes/dUinf
      GAMU(N,1) = -CBIS
C
C----- -dRes/dVinf
      GAMU(N,2) = -SBIS
C
      ENDIF

```



```

C
C---- LU-factor coefficient matrix AIJ
CALL LUDCMP(IQX,N+1,AIJ,AIJPIV)
LQAIJ = .TRUE.
C
C---- solve system for the two vorticity distributions
CALL BAKSUB(IQX,N+1,AIJ,AIJPIV,GAMU(1,1))
CALL BAKSUB(IQX,N+1,AIJ,AIJPIV,GAMU(1,2))
C
C---- set inviscid alpha=0,90 surface speeds for this geometry
DO 50 I=1, N
    QINVU(I,1) = GAMU(I,1)
    QINVU(I,2) = GAMU(I,2)
50 CONTINUE
C
    LGAMU = .TRUE.
C
    RETURN
END

```

## Mesh\_Sail: Program for Creating Sail Mesh

```

Program Mesh_sail
include 'section.inc'

real:: length, height, aflow, p(3),dt
real:: org(3),x(3),y(3) !origin
real,dimension(3,3):: vr
type(section):: csec(7),tesec
real:: intquad

pi = 3.1415
open(1,file = 'dados.txt')

!reading height,footleng and flow incidence
read(1,*) height,length,aflow
!reading sections
! length, entry angle, te angle
j = 1
do i = 1,4
    read(1,*) csec(j).leng,csec(j).th1,csec(j).th2
    j = j+2
enddo
!reading trailing edge (te) section
read(1,*) trv,abat,tesec.cpos,tesec.camber
!reading foot and top section angles
read(1,*) alfa1,beta1,alfa2,beta2
close(1)
dt = 1./3
t = 0
do i=1,7,2
    csec(i).t = t
    t=t+dt
enddo
aflow = aflow*pi/180
abat = abat*pi/180
alfa1 = alfa1*pi/180
alfa2 = alfa2*pi/180
beta1 = beta1*pi/180
beta2 = beta2*pi/180

!calculation of intermediary section angles

do i = 2,6,2
    csec(i).th1 =(csec(i+1).th1+csec(i-1).th1)/2
    csec(i).th2 =(csec(i+1).th2+csec(i-1).th2)/2
    csec(i).t = (csec(i+1).t+csec(i-1).t)/2
    if (i==2) then
        k = 1
        do j=1,3
            x(j) = csec(k).t
            y(j) = csec(k).leng

```

```

        k = k+2
    enddo
else
    k=i-3
    do j=1,3
        x(j) = csec(k).t
        !t = t+tr
        y(j) = csec(k).leng
        k = k+2
    enddo
endif
csec(i).leng = intquad(x,y,csec(i).t) !quadratic interpolation
enddo

!generation of transversal sections
do k =1,7
    !xm = csec(k).cpos
    !ym = csec(k).camber
    !call solve_foil(xm,ym,a,b,c)
    ds = 1.0/10
    s = 0
    csec(k).p1 = 0
    call set_sec(csec(k))
    do i = 2,11
        s=s+ds
        csec(k).p1(2,i)= csec(k).p1(2,i)*csec(k).leng
        ! (a*s**3+b*s**2+c*s)*csec(k).leng
        csec(k).p1(1,i)= csec(k).p1(1,i)*csec(k).leng
        !s*csec(k).leng
    enddo
enddo
!foot section height
zf = tan(beta1)
b = tan(alfa1)
a = zf-b
s = 0
csec(1).p1(3,1) = 0
do i = 2,11
    s = s+ds
    csec(1).p1(3,i) = (a*s**2+b*s)*csec(1).leng
enddo

!top section height
zf = tan(beta2)
b = tan(alfa2)
a = zf-b
s = 0
csec(7).p1(3,1) = height
do i = 2,11
    s = s+ds
    csec(7).p1(3,i) = height+(a*s**2+b*s)*csec(1).leng
enddo

!intermediate sections height
dt = 1.0/6
t=0
do i = 2,6
    t = t+dt
    do j = 1,11
        !df = csec(4).p1(3,j)+csec(1).p1(3,j)
        csec(i).p1(3,j) = t*height !df
    enddo
enddo

!Generation of te section
xm = tesec.cpos
ym = tesec.camber
call solve_foil(xm,ym,a,b,c)
t = 0
!tesec.p1 = 0
do i = 1,7

```

```

        tesec.pl(2,i)= (a*t**3+b*t**2+c*t)*height
        tesec.pl(3,i)= csec(i).pl(3,10)
        tesec.pl(1,i)= csec(i).pl(1,10)
        t=t+dt
    enddo

    !rotation of te section (just y coordinate)
    a1 = aflow-abat
    xp = csec(4).pl(1,10)
    yp = csec(4).pl(2,10)
    p(1) = xp*cos(a1)-yp*sin(a1)
    p(2) = xp*sin(a1)+yp*cos(a1)
    p(3) = csec(4).pl(3,10)
    a2 = atan(p(2)/p(3))
    do i = 1,7
        zp = tesec.pl(3,i)
        yp = tesec.pl(2,i)
        !tesec.pl(3,i) = zp*cos(a1)-yp*sin(a1)
        tesec.pl(2,i) = zp*sin(a2)+yp*cos(a2)
    enddo

    !translating te section
    yt = trv*length
    do i = 1,7
        tesec.pl(2,i) = tesec.pl(2,i) +yt
    enddo

    !Calculating central and top sections twist angles
    do k = 1,7
        csec(k).asec = atan(tesec.pl(2,k)/
&          (tesec.pl(1,k)-csec(k).pl(1,1)) )
    enddo

    ! rotating sections
    do k = 1,7
        a1 = csec(k).asec
        xo = csec(k).pl(1,1)
        yo = 0
        do i = 2,11
            xp = csec(k).pl(1,i)
            yp = csec(k).pl(2,i)
            csec(k).pl(1,i) =xo+ xp*cos(a1)-yp*sin(a1)
            csec(k).pl(2,i) =yo+ xp*sin(a1)+yp*cos(a1)
        enddo
    enddo
    !writting msh file
    M = 11 !chordwise
    N = 7  !spanwise
    open(2,file = 'c:\codigos\mshuns\sail.msh')
    do i = 1,12
        write(2,10)
    enddo
    write(2,20) M,N
    t = 0
    do i = 1,N
        s = 0
        do j = 1,M
            write(2,30) csec(i).pl(1,j),csec(i).pl(2,j),csec(i).pl(3,j),
&          s,t
            s = s+ds
        enddo
        t = t+dt
    enddo
    close(2)

10    format('%',1x)
20    format(1x,i4,1x,i4)
30    format(1x,f8.5,1x,f8.5,1x,f8.5,1x,f8.5,1x,f8.5)
    end program

!*****

```

```

      subroutine set_sec(sec)
      include 'section.inc'
      !This program generates the section using the Jackson Polynomial
      !Ref: P.S. Jackson, "A Simple Model for 2D Sails
      ! AIAA Technical notes 1983
      ! Author: Augusto Veiga

      type(section):: sec
      real:: A,B, pi
      real:: delta, a1,b1,c1

      pi = 4. * atan(1.)
      sec.th1 = sec.th1*pi/180
      sec.th2 = sec.th2*pi/180
      A = sec.th1+sec.th2
      B = sec.th1-sec.th2

      !seeking maximum camber position
      a1 = -0.75*B*2
      b1 = -0.5*A
      c1 = 0.25*B

      delta = b1**2-4*a1*c1
      if (delta>=0) then
         r = (b1-sqrt(delta))/(2*a1)
         sec.cpos = (1+r)/2.
         tm = 0.25*(1-r**2)*(A+B*r)
         sec.camber = tm/2.
      endif

      !Generating sections
      ds = 2.0/10
      dx = 1.0/10
      s = -1
      x = 0
      do i=1,11
         t = 0.25*(1-s**2)*(A+B*s)
         sec.pl(2,i) = t/2
         sec.pl(1,i) = x
         s = s+ds
         x = x+dx
      enddo
      return
      end subroutine

!*****
      real function intquad(x,y,x1)
      real,dimension(3):: x,y
      real:: x1,sum

      sum = 0
      sum = sum+((x1-x(2))*(x1-x(3)))/((x(1)-x(2))*(x(1)-x(3)))*y(1)
      sum = sum+((x1-x(1))*(x1-x(3)))/((x(2)-x(1))*(x(2)-x(3)))*y(2)
      sum = sum+((x1-x(1))*(x1-x(2)))/((x(3)-x(1))*(x(3)-x(2)))*y(3)

      intquad = sum
      end function
      subroutine solve_foil(xm,ym,a,b,c)
      !makes the foil using a 3rd order polynomial
      real:: xm,ym,a,b,c
      real:: vr(3,3), v(3)

      do i=1,3
         k = 4
         if (i==2) then
            v(i) = ym/xm
         else
            v(i) = 0
         endif
         do j=1,3
            k = k-j
            if (i==1) then

```

```

        vr(i,j) = k*xm**(k-1)
      else if (i==2) then
        vr(i,j) = xm**(k-1)
      else
        vr(i,j) = 1.0
      endif
    enddo
  enddo
  !solve system using Gauss elimination
  call gauss(3,3,Vr,v,1)
  a = v(1)
  b = v(2)
  c = v(3)

return
end subroutine

SUBROUTINE GAUSS(NSIZ,NN,Z,R,NRHS)
C *****
C *
C *   Solves general NxN system in NN unknowns
C *   with arbitrary number (NRHS) of righthand sides.
C *   Assumes system is invertible...
C *   ...if it isn't, a divide by zero will result.
C *
C *   Z is the coefficient matrix...
C *   ...destroyed during solution process.
C *   R is the righthand side(s)...
C *   ...replaced by the solution vector(s).
C *
C *****
C
C   DIMENSION Z(NSIZ,NSIZ), R(NSIZ,NRHS)
C
C   DO 1 NP=1, NN-1
C     NP1 = NP+1
C
C----- find max pivot index NX
C     NX = NP
C     DO 11 N=NP1, NN
C       IF (ABS(Z(N,NP))-ABS(Z(NX,NP))) 11,11,111
111     NX = N
11    CONTINUE
C
C     PIVOT = 1.0/Z(NX,NP)
C
C----- switch pivots
C     Z(NX,NP) = Z(NP,NP)
C
C----- switch rows & normalize pivot row
C     DO 12 L=NP1, NN
C       TEMP = Z(NX,L)*PIVOT
C       Z(NX,L) = Z(NP,L)
C       Z(NP,L) = TEMP
12    CONTINUE
C
C     DO 13 L=1, NRHS
C       TEMP = R(NX,L)*PIVOT
C       R(NX,L) = R(NP,L)
C       R(NP,L) = TEMP
13    CONTINUE
C
C----- forward eliminate everything
C     DO 15 K=NP1, NN
C       ZTMP = Z(K,NP)
C
C       IF (ZTMP.EQ.0.0) GO TO 15
C
C       DO 151 L=NP1, NN
C         Z(K,L) = Z(K,L) - ZTMP*Z(NP,L)

```

```

151      CONTINUE
        DO 152 L=1, NRHS
          R(K,L) = R(K,L) - ZTMP*R(NP,L)
152      CONTINUE
15      CONTINUE
C
      1 CONTINUE
C
C---- solve for last row
      DO 2 L=1, NRHS
        R(NN,L) = R(NN,L)/Z(NN,NN)
      2 CONTINUE
C
C---- back substitute everything
      DO 3 NP=NN-1, 1, -1
        NP1 = NP+1
        DO 31 L=1, NRHS
          DO 310 K=NP1, NN
            R(NP,L) = R(NP,L) - Z(NP,K)*R(K,L)
310        CONTINUE
31      CONTINUE
      3 CONTINUE
C
      RETURN
      END ! GAUSS

```