

```

subroutine adummy
c..... cliche storage set up here

```

```

cliche param
parameter (izx=60, jrx=40, izx2=izx-2 )
parameter ( ithw=2*izx-1 )
parameter (kxp=(izx-2)*(jrx-2), ibw=izx-1 )
parameter (nplt=3000, nps=100 )
parameter (ixpf=4*(izx-2),nfourxf=150)
parameter (ixpg=2*izx2)
parameter (neng=500)
parameter (ksim=51)
endcliche

```

```

cliche fstor
use param
common/fun/f1(izx,jrx),f2(izx,jrx),f3(izx,jrx),f4(izx,jrx)
c ,f5(izx,jrx),f7(izx,jrx),
c g1(izx,jrx),g2(izx,jrx),g3(izx,jrx),g4(izx,jrx)
c ,swg1,swg2,swg3,swg4

```

```

common/equil/b(izx,jrx),rho(izx,jrx),qub(izx,jrx),r(izx,jrx)
c ,phi(izx,jrx),yyy(izx,jrx),xxx(izx,jrx),qv(izx,jrx)
common/pertur/xloo(kxp),xlo(kxp),xlo1(kxp)
c ,xron(kxp),xro(kxp),xro1(kxp)
endcliche

```

```

cliche matrix
use param
common/coeff/a1(kxp,9),a2(kxp,9),a3(kxp,9),b1(kxp,3)
c rhs1(kxp),rhs2(kxp)
c.....unnamed common for dynamic memory expansion
common ww(1), ww1(1)
endcliche

```

```

cliche const
use param
common/title/aname(5)
common/con/gam1,gam2,ix,jx,mm,izxp,kxx,nmax,lmax,lsz,ithw
c ,fac1,fac2,bias,du,dv,dt,ndiag,ex0,b0,rho0,ex1,f11,fizx,fj1
c ,fjrx,kplot,npm,fpsi,fz,fu,fv,azm,apsim,u0,v0,amass,nengx
c ,fourpi,omegst,omegr,omegexb,flr,sf6,sf8,kplotm,kzs,zedge
c ,cpuo,cio,syso,va1fk,xu,xv,n,pl,vw,pslw,dvin,dvout,ltt,nen,lee
common/contm/
c psi0rel,psi1rel,psi2rel,z1rel,z2rel,z3rel,z0rel,nslosh,bmg
c ,ncenter,pslosh,pcenter,rp,ztrans,ltrans,bm,ltran,ztran,rp1
c ,buen,pring,epsp,phicen,phiplg,kin,xpot,ypot,wpot,pfudge,rpx
c ,phice,phipl,betslsh,betcent,z0,z1,z2,z3,z4,psi1,psi2,psi0
c ,betcone,betslse,psloshe,pcentee,bmax,alsi,bm1,psls1,cold
c ,p2wide,psi3rel,psi3,p1max,bv0,bv1,bv4,bceng,psloshin,psloshen
c ,nsloshin,pxp1,pxp2,p3a,p3b,p3c,p3d,psim,pe10,ae1,be1,ce1,de1
c ,psi0erel,psi0e,psime,p2ewide,wp2e,p2floor,p1floor,p2flag
c ,fring,long,no3d,no1d,dphi,dip,psistr,psisl,psihrel,psih
c ,dpsihrel,dpsih,er,rcwall,exrho
common/pcons/at0,bt0,ct0,cp0
c ,ap1,ap2,ap3,at1,at2,at3,bp1,bp2,bp3,bt1,bt2,bt3,cp1,cp2,cp3
c ,ct1,ct2,ct3,bmx1,bmx2,bmx3,bm1,bm2,ppas1,ppas2,ppas3,p1trap
c ,z1c,z2c,z3c,z1min,z2min,as(3),als(3),zs(3),bs(3),dpas1,d1trap
c ,betrap,betp,z1,bvx2,bvx3,ncoil,z2ct
common/mesh/psi(jrx),z(izx),u(izx),v(jrx),dpsi(jrx),dz(izx)

```

```

c , vpsi(jrx), uuz(izx), vpsi(jrx), uuz(izx)
common/graf/ xrtime(nplt), xrspsz(izx,nps), xrsppsi(jrx,2*nps)
c , time(nplt), xflute(izx,nps), enpot(neng), tenergy(neng)
x , enkin(neng), timengy(neng), tenrel(neng), enbend(neng)
c , encurve(neng), enfir(neng)
common/curvco/cr, lb, rw, beta0, delrho, stable, en0, cee, r0,
c echarg, omeg1, omeg2, en1, besarg, zol, dtrel, p0, omeg0
c , omana1, omana2, groana, theta0
common/tmcon/h12(izx), hzt0(izx), h34(izx), abp(izx), bbp(izx)
c , cbp(izx), abf(izx), bbf(izx), cbf(izx), hp3(jrx), hp12(jrx)
c , htrans(izx), abq(izx), bbq(izx), cbq(izx), ebp(izx)
c , fbp(izx), gbp(izx), betring, hp0(jrx), hpm(jrx), hpm0(jrx), hfir(jrx)
c , hzp0(izx), hzp1(izx), hzp2(izx), hzp3(izx), hzt1(izx), hzt2(izx)
c , hzt3(izx)
common/tmfield/bvac(izx), dbvdz(izx), d2bvdz2(izx), dp1dpsi(jrx)
c , p1(jrx), p1k(ksim, jrx), hpk12(ksim), hpk0(ksim), deli1(ksim)
c , deli2(ksim), deli3(ksim), deli4(ksim), rzz(izx, jrx), dbdpsi(izx, jrx)
c , phi1(izx), phi2(izx), pperp(izx, jrx), ppar(izx, jrx), dflute3(izx)
c , qubv(izx, jrx), p2(jrx), dp2dpsi(jrx), dflute1(izx), dflute2(izx)
c , flute1(jrx), flute2(jrx), flute3(jrx), p2k(ksim, jrx), deli5(ksim)
c , pperps(izx, jrx), errprp(izx, jrx), errprl(izx2, jrx-1)
c , pperpe(izx, jrx), epsi(izx, jrx), omeg1wkb(jrx), omeg2wkb(jrx)
c , gamwkb(jrx), dflute4(izx), rhoave(jrx), xxxave(jrx), yyyave(jrx)
c , p2t(jrx), dp2dpst(jrx), p3(jrx), dp3dpsi(jrx), hpkm(ksim)
c , hpkm0(ksim), droave(jrx), droterm(izx), ering(izx, jrx)
common/forced/nfour, nfourx, nfourmax, nfourp, jfour, ixp, locv

real lb, ltrans, ltran, nslosh, ncenter, nsloshin
endliche

return
end

```

c..... the main routine

```

c *****
c *
c   FLORA is an initial value stability code developed by R. Freis and
c   B. Cohen, based on Newcomb's long thin axisymmetric formalism, including
c   finite larmor radius effects. FLORA calculates the linear response to
c   low frequency perturbations of the equilibrium magnetic field .

```

```

c *****
c..... notice of 4/8/82, this version runs correctly for isw=1, and
c..... runs correctly for isw=-1 .

```

```

c.....5/12/82, flora runs testcase 1 , 0 beta, 0 pressure, homogeneous
c..... plasma, correctly.

```

```

c..... floral transforms variables z,psi to u,v which are always equally
c..... spaced. transformation: z=au*u**xu, and psi=apsi*v**xv, where
c..... zmax=umax, psimax=vmax, and fz*zmax=fu*umax, fpsi*psimax=fv*vmax .
c..... fz, fu, fpsi, fv, input, xu=ln fz / ln fu, xv=ln fpsi / ln fv .
c..... au=umax*(-xx+1) , apsi=vmax*(-yy+1) .

```

```

c..... flora2 solves test case 2 , rotating rigid rotor stability. ref:
c..... freidberg and pearlstein, phys fluids 21(7) july 1978 1207

```

c.....flora4 includes background constant density, enbar (as does flora3),

c..... and kzs switch which when set to zero, generates initial perturbations
c..... independent of z in random spatial generator (ex0=1.) .

c..... flora5
c..... is vectorized version of flora4, (calls rightvec instead of
c..... right), also has timing routine from b. langdon (requires
c..... bzohar loaded as a binary).
c....., insert cliché storage here

c..... flora7 is mod. flora5, with psi stretching function
c..... exactly centered in amat. (flora5 used linear interpolation
c..... to get vpsi(j+1/2)). Also revised diagnostic plots included.

c.....flora12 is flora11 (rigid rotor with corrected equil. and
c..... corrected curvature terms (flora10!)) with fourier mode analyses
c..... added (using cpft and rpft) and data for zed post processing.
c.....,additional input data: jfour (v index at which xr is analyzed in
c.....,z), nfourp (analyze xr every nfour'th time step),nfourmax
c..... (number of times the buffer is read to the history file), note
c.....,xr is extended a factor of 4 to look like a periodic full wave
c.....,for cpft. If jfour is input 0, code sets it to jx/4 .
c

c.....flora13 is flora12 with curvature driven flute mode equilibrium
c.....,(equilrot replaced by equilcur, rigidcon replaced by curvecon)

c.....,floralm, tandem mirror equilibrium

c
c.....,flortm1, tandem mirror equilibrium, with 3-d plot of equilb.
c quantities added. (uses tv80 and graflib)

c.....,flortex, tandem mirror equilb. with corrections to flortm1. In-
c.....,put switches swg1, swg2, swg3, swg4 added.

c.....,flortm2, like flortex with revised electron ring, a la D'ippolito
c.....,(-ring pperp in b field only, and additional term in curvature
c.....,drive),

c.....,flortm3, like flortm2 with corrections to pressure normalization,
c.....,and additional diagnostics. (3-d plots of curvature drive-e ring
c.....,term, and perp. pressure balance check) , also 3-d plots of
c.....,pparallel pressure check, and e-psi (=-dphi/dpsi) . Phi2 modified
c.....,to ' , -(psi/psi3)**ypot ,

c.....,flortm4, modified plasma pperp with addition of p3(j) to give
c.....,a positive slope near the center.

c.....,flortm5, modified p1 in flortm4 to be two functions, pe1 and
c.....,pe2, joined at psime with equal slope and value, pe1=a*el+be1*(
c.....,psi/psime)+ce1*(psi/psime)**2, and pe2=.5*(1-tanh((psi-psime)/p2ewide))
c

c.....,flortm6, modified flortm5 as follows: for p2(psile, to p2flag (an
c.....,input value), p2 set to p2floor (an input value) and p1 set to
c.....,p1floor (an input value). Long-thin ering option added. This modifies
c.....,b dependence of ering pperp to look longer (by changing abf, bbf, cbf)
c.....,if long (an input value) = 1, otherwise leaves pperp of ring un-
c.....,changed. Plot output options, nold=1. prevents graflib plots, no3d=1
c.....,prevents tv80lib 3d plots.
c

```

c..... flortm8 (l18 for larger psi grid) modified flortm6 as follows;
c..... the analytic calculation of gamwkb corrected to include drho/dpsi
c..... term (important in the limit of large exb rotation), also phi1
c..... changed to be constants in core and plug ( phicen in core, and
c..... phicen+dphi in plug, dphi a new input variable ). Also b calculated
c..... with expansion in low beta regions.

```

```

c..... flortn8, like flortl8 with first order energy check added .

```

```

c..... flrm1 like flortn8 with multi region equilibrium.
c..... Bvac is generated from 3 solenoids ( 1 choke coil and 2
c..... mirror coils). Pressures are the sum of passing and trapped
c..... components. See the glossary of input parameters in subroutine
c..... inputtm for the revised list .

```

```

c..... flrm2, either 2 or 3 regions, depending on the number of solenoids
c..... specified (ncoil=2, or 3 ) in the input. For ncoil=2, no passing
c..... pressures are allowed, and the situation is similar to earlier
c..... versions, except that the vacuum fields are generated by solenoids
c..... instead of circular filaments.

```

```

c..... flrm3, like flrm2 with corrections to energy subroutine. Also
c..... option to remove hollowness from pperp psi profile (dip=0.), and
c..... ering psi profile changed from quadratic to cubic in inner region.

```

```

c..... flrm4, cold plasma halo modeled by changing zmax boundary conditions
c..... for psi > psih (psihrel an input parameter )

```

```

c..... flrm6, like flrm4 and flrm5, (mixed boundary condition at zmax,
c..... higher order b, c.) with special yyy to force "rigid mode" in psi.

```

```

c..... flrrot, modified flrm6 to study rotational stability with pos-
c..... itive density gradients. Set phi2(psi)=er (a constant), add
c..... rhoc = a + b*psi to rho .

```

```

      use param
      use fstor
      use matrix
      use const

```

```

      data tim/1.e6/
      integer tallyb(2000b)
      common / q8locs/locf(0:15)
      common/pic100/npete
      data itally/1/

```

```

c..... call link call here
      call link('unit59=terminal,unit2=(infil4,open),unit3=(output,
c create) //')

```

```

      if(itally.gt.0) then
      do 200 ii=1,15
200   if(locf(ii).eq.0)go to 210
      ii=0
210   loctally=ii
      loctally=14
      if(loctally.eq.0)go to 299
      call timer(loctally,'ztally00',tallyb,2000b,florati,1)
299   itally=-1
      endif

```

```

      isw=1
      if(izx.gt,jrx)isw=-1
      jtbw=.5*(1+isw)*itbw+.5*(1-isw)*(2*jrx-1)
      ihbw=.5*(1+isw)*ibw+.5*(1-isw)*(jrx-1)
      nn=jtbw*kxp
      nn1=jtbw+kxp
      call memory(ww,nn-1)
      call memory(ww1(nn),nn1)
      namelist/noplot/no1d,no3d
      call ddi(noplot,2,0,1)
      if(no1d.ne.1)call pstart(dev,4rplot,1,'box u21$',1)
c      npete=1
      if(no1d.ne.1)call p100
      call input
      call inputtm
c      .... temporary input to test three region model
c      call inptemp
c      call rigidcon
c      call curvecon
      call constant
      call tmcon2
      call equilrm
c      call equilrot
c      call equilcur
      call fltoll
      call amat
      call comat(ww,jtbw)
      call initial
c      .... special version for testing fourier analysis and zed file
c      .... maker
      call fourplay
      call fourier
      call mymove(xrol(1),xro(1),kxx)
      call mymove(xiol(1),xio(1),kxx)
c      call mymove(xroo(1),xro(1),kxx)
c      call mymove(xioo(1),xio(1),kxx)
      call banfac(kxp,ihbw,ww,1,-(kxp-1))
      call energy
      t=0.
      do 100 n=1,nmax
      t=t+dt
      time(n)=t
      fac1=-1./dt
      fac2=1./dt
      do 90 l=0,lmax
      call rightvec
      call zmovewrd(ww1(nn),rhs1,kxx)
      call bansol(kxp,ihbw,ww,1,-(kxp-1),ww1(nn))
      do 10 j=2,jx-1
      kp=1+izxp*(j-2)
      call zmovewrd(xrol,ww1(nn),kxx)
10      continue
      call zmovewrd(ww1(nn),rhs2,kxx)
      call bansol(kxp,ihbw,ww,1,-(kxp-1),ww1(nn))
      do 20 j=2,jx-1
      kp=1+kxp*(j-2)
      call zmovewrd(xiol,ww1(nn),kxx)
20      continue
      fac1=-.5/dt
90      fac2=.5/dt

```

```

        call zmovewrd(xloo,xlo,kxx)
        call zmovewrd(xlo,xlo1,kxx)
        call zmovewrd(xroo,xro,kxx)
        call zmovewrd(xro,xro1,kxx)
c..... time array
        xrtime(n)=xro(kplot)
        if(mod(n,ndiag).eq.0)call diagno
        if(mod(n,nfourp).eq.0)call fourier
        if(((mod(n,nen1)).eq.0).or.((ltt.ne.0).and.((lee.le.neng)))call energy
100    continue
        call clsdsk(iocv,0)
        call timeused(icp,io,isy)
        cpuo=icp*tim
        cio=io*tim
        syso=isy*tim
        if(nold.ne.1)
c call picsher
        call close(100)
        if(no3d.eq.1)go to 300
        call keep80(1,3)
        call fr80id
        call threed
        call plote
300    continue
        call timend
        call exit(1)
        end
        subroutine amat

```

c..... calculates the matrix coefficients for a1, a2, a3, b1, b2
c..... in the equation $a1*x(n+1)=a2*x(n)+a3*x(n-1)+b1*y(n)+b2*y(n-1)$.
c..... uses f1 to f11 from subroutine fltol1 and equilibrium quantities.
c..... cliché storage here

```

        use param
        use fstor
        use matrix
        use const

        dimension bc(jrx),delco1(jrx),delco2(jrx)
        data unit/1./

        gam3=-gam2
        du2=du**2
        dt2=dt**2
        dv2=dv**2
        dvt=2.*dv
        m2=mm**2
        jx=jrx
        ix=ixx
        do 10 i=2,ix-1
        do 10 j=2,jx-1
        k1=i-1+(j-2)*(ix-2)
        k2=j-1+(jx-2)*(i-2)
        k=.5*(1+isw)*k1+.5*(1-isw)*k2
        r2=r(i,j)**2
        vp=vpsi(j)
        uz=uuz(i)
        bijmh=(b(i,j)+b(i,j-1))*5

```

```

b1jph=(b(i,j)+b(i,j+1))*5
b1p1jph=(b(i+1,j+1)+b(i+1,j))*5
b1p1jmh=(b(i+1,j-1)+b(i+1,j))*5
b1m1jph=(b(i-1,j+1)+b(i-1,j))*5
b1m1jmh=(b(i-1,j-1)+b(i-1,j))*5
g4iphjph=(g4(i+1,j+1)+g4(i,j)+g4(i+1,j)+g4(i,j+1))*25*uzzh(i)
g4iphjmh=(g4(i+1,j-1)+g4(i,j)+g4(i+1,j)+g4(i,j-1))*25*uzzh(i)
g4imhjph=(g4(i-1,j+1)+g4(i,j)+g4(i-1,j)+g4(i,j+1))*25*uzzh(i)
g4imhjmh=(g4(i-1,j-1)+g4(i,j)+g4(i-1,j)+g4(i,j-1))*25*uzzh(i)
g2iphj=(g2(i+1,j)+g2(i,j))*5*uzzh(i)
g2imhj=(g2(i-1,j)+g2(i,j))*5*uzzh(i-1)
g3iphj=(g3(i+1,j)+g3(i,j))*5*uzzh(i)
g3imhj=(g3(i-1,j)+g3(i,j))*5*uzzh(i-1)

```

```

f1ijph=(f1(i,j)+f1(i,j+1))*5*vpsi h(j)
f1ijmh=(f1(i,j)+f1(i,j-1))*5*vpsi h(j-1)
f5ijph=(f5(i,j)+f5(i,j+1))*5*vpsi h(j)
f5ijmh=(f5(i,j)+f5(i,j-1))*5*vpsi h(j-1)

```

```

60 if(j.gt.2)go to 60

```

```

f1ijmh=0,
f6ijmh=0,

```

```

continue

```

```

uzbar=-uuz(i)*r(i,j)/(du2*dv2)

```

```

a1(k,1)=-gam1*b1m1jmh*g4imhjmh*b1jmh*uzbar*vpsi h(j-1)

```

```

c *r(i-1,j-1)

```

```

a2(k,1)=-gam2*b1m1jmh*g4imhjmh*b1jmh*uzbar*vpsi h(j-1)

```

```

c *r(i-1,j-1)

```

```

a3(k,1)=-gam3*b1m1jmh*g4imhjmh*b1jmh*uzbar*vpsi h(j-1)

```

```

c *r(i-1,j-1)

```

```

a1(k,2)=-f1ijmh/((dt*dv)**2)+gam1*(f5ijmh/dv2+b1jmh**2*uzbar

```

```

c *vpsi h(j-1)*r(i,j-1)*(g4imhjmh+g4iphjmh)

```

```

a2(k,2)=-f1ijmh/((dt*dv)**2)+gam2*(f5ijmh/dv2+b1jmh**2*uzbar

```

```

c *vpsi h(j-1)*r(i,j-1)*(g4imhjmh+g4iphjmh)

```

```

a3(k,2)=-f1ijmh/((dt*dv)**2)+gam3*(f5ijmh/dv2+b1jmh**2*uzbar

```

```

c *vpsi h(j-1)*r(i,j-1)*(g4imhjmh+g4iphjmh)

```

```

a1(k,3)=-gam1*b1p1jmh*g4iphjmh*b1jmh*uzbar*vpsi h(j-1)*r(i+1,j-1)

```

```

a2(k,3)=-gam2*b1p1jmh*g4iphjmh*b1jmh*uzbar*vpsi h(j-1)*r(i+1,j-1)

```

```

a3(k,3)=-gam3*b1p1jmh*g4iphjmh*b1jmh*uzbar*vpsi h(j-1)*r(i+1,j-1)

```

```

a1(k,4)=gam1*((b1m1jmh*g4imhjmh*vpsi h(j-1)*b1jmh+b1m1jph*g4imhjph

```

```

c *vpsi h(j)*b1jph)*uzbar*r(i-1,j)+mm**2*b(i,j)*uzbar*

```

```

c g2imhj*g3(i-1,j)*dv2/vpsi(j))

```

```

a2(k,4)=gam2*((b1m1jmh*g4imhjmh*vpsi h(j-1)*b1jmh+b1m1jph*g4imhjph

```

```

c *vpsi h(j)*b1jph)*uzbar*r(i-1,j)+mm**2*b(i,j)*uzbar*

```

```

c g2imhj*g3(i-1,j)*dv2/vpsi(j))

```

```

a3(k,4)=gam3*((b1m1jmh*g4imhjmh*vpsi h(j-1)*b1jmh+b1m1jph*g4imhjph

```

```

c *vpsi h(j)*b1jph)*uzbar*r(i-1,j)+mm**2*b(i,j)*uzbar*

```

```

c g2imhj*g3(i-1,j)*dv2/vpsi(j))

```

```

a1(k,5)=((f1ijph+f1ijmh)/dv2-f2(i,j))/dt2+gam1*(-(f5ijph+f5ijmh

```

```

c )/dv2+f7(i,j)+g1(i,j)+(-b1jmh**2*(g4imhjmh+g4iphjmh)*vpsi h(j-1)

```

```

c -b1jph**2*(g4imhjph+g4iphjph)*vpsi h(j))*r(i,j)*uzbar-

```

```

c mm**2*b(i,j)*uzbar*(g2imhj+g2iphj)*g3(i,j)*dv2/vpsi(j))

```

```

a2(k,5)=((f1ijph+f1ijmh)/dv2-f2(i,j))/dt2+gam2*(-(f5ijph+f5ijmh

```

```

c )/dv2+f7(i,j)+g1(i,j)+(-b1jmh**2*(g4imhjmh+g4iphjmh)*vpsi h(j-1)

```

```

c -b1jph**2*(g4imhjph+g4iphjph)*vpsi h(j))*r(i,j)*uzbar-

```

```

c mm**2*b(i,j)*uzbar*(g2imhj+g2iphj)*g3(i,j)*dv2/vpsi(j))

```

```

a3(k,5)=(f1ijph+f1ijmh)/dv2-f2(i,j)/dt2+gam3*(-(f5ijph+f5ijmh
c 1/dv2+f7(i,j)+g1(i,j)+(-bijmh**2*(g4imhjmh+g4iphjmh)*vpsih(j-1)
c -bijph**2*(g4imhjph+g4iphjph)*vpsih(j))*r(i,j)*uzbar-
c mm**2*b(i,j)*uzbar*(g2imhj+g2iphj)*g3(i,j)*dv2/vps(i,j))

a1(k,6)=gam1*((bip1jmh*g4iphjmh*bijmh*vpsih(j-1)+bip1jph*g4iphjph
c *bijph*vpsih(j))*uzbar*r(i+1,j)+mm**2*b(i,j)*uzbar*
c g2iphj*g3(i+1,j)*dv2/vps(i,j))
a2(k,6)=gam2*((bip1jmh*g4iphjmh*bijmh*vpsih(j-1)+bip1jph*g4iphjph
c *bijph*vpsih(j))*uzbar*r(i+1,j)+mm**2*b(i,j)*uzbar*
c g2iphj*g3(i+1,j)*dv2/vps(i,j))
a3(k,6)=gam3*((bip1jmh*g4iphjmh*bijmh*vpsih(j-1)+bip1jph*g4iphjph
c *bijph*vpsih(j))*uzbar*r(i+1,j)+mm**2*b(i,j)*uzbar*
c g2iphj*g3(i+1,j)*dv2/vps(i,j))

a1(k,7)=gam1*(-bim1jph*g4imhjph*bijph*vpsih(j)*uzbar*r(i-1,j+1))
a2(k,7)=gam2*(-bim1jph*g4imhjph*bijph*vpsih(j)*uzbar*r(i-1,j+1))
a3(k,7)=gam3*(-bim1jph*g4imhjph*bijph*vpsih(j)*uzbar*r(i-1,j+1))
a1(k,8)=-f1ijph/(dt2*dv2)+gam1*(f5ijph/dv2+(bijph**2*(g4imhjph
c +g4iphjph)*vpsih(j))*r(i,j+1)*uzbar)
a2(k,8)=-f1ijph/(dt2*dv2)+gam2*(f5ijph/dv2+(bijph**2*(g4imhjph
c +g4iphjph)*vpsih(j))*r(i,j+1)*uzbar)
a3(k,8)=-f1ijph/(dt2*dv2)+gam3*(f5ijph/dv2+(bijph**2*(g4imhjph
c +g4iphjph)*vpsih(j))*r(i,j+1)*uzbar)
a1(k,9)=gam1*(-bip1jph*g4iphjph*bijph*vpsih(j)*uzbar*r(i+1,j+1))
a2(k,9)=gam2*(-bip1jph*g4iphjph*bijph*vpsih(j)*uzbar*r(i+1,j+1))
a3(k,9)=gam3*(-bip1jph*g4iphjph*bijph*vpsih(j)*uzbar*r(i+1,j+1))
c.....,b1 array for rhs
f3ijmh=(f3(i,j)+f3(i,j-1))*0.5*vpsih(j-1)
f3ijph=(f3(i,j)+f3(i,j+1))*0.5*vpsih(j)
denom=1./(dv2)
b1(k,1)=f3ijmh*denom
b1(k,3)=f3ijph*denom
b1(k,2)=-(f3ijmh+f3ijph-f4(i,j)*dv2/vp)*denom
10 continue
c.....,correct coefficients on boundaries
sf11=sign(unit,f11)
sfj1=sign(unit,fj1)
sfjrx=sign(unit,fjrx)
sfizx=sign(unit,fizx)
c....., set corners to 0
k1i=ix-2
k1j=1+(ix-3)*(jx-2)
k1=.5*(1+isw)*k1i+.5*(1-isw)*k1j
fac1=-1.
if(sfj1.eq.1.and.sfizx.eq.1)fac1=r(ix-1,2)*b(ix-1,2)/
c (r(ix,2)*b(ix,2))
a1(k1,5)=a1(k1,5)+fac1*a1(k1,3)
a2(k1,5)=a2(k1,5)+fac1*a2(k1,3)
a3(k1,5)=a3(k1,5)+fac1*a3(k1,3)
a1(k1,3)=0.
a2(k1,3)=0.
a3(k1,3)=0.
k2i=1+(ix-2)*(jx-3)
k2j=jx-2
k2=.5*(1+isw)*k2i+.5*(1-isw)*k2j
fac3=-dvout/dvin
if(sfjrx.eq.1.and.sf11.eq.1)fac3=r(2,jx-1)*b(2,jx-1)/
c (r(1,jx-1)*b(1,jx-1))
a1(k2,5)=a1(k2,5)+fac3*a1(k2,7)

```



```

    a2(k2,5)=a2(k2,5)+fac3*a2(k2,7)
    a3(k2,5)=a3(k2,5)+fac3*a3(k2,7)
    a1(k2,7)=0,
    a2(k2,7)=0,
    a3(k2,7)=0,
    fac2=-1,
    if(sfj1,eq,1,and,sf11,eq,1)fac2=r(2,2)*b(2,2)/(r(1,2)*b(1,2))
    a1(1,5)=a1(1,5)+fac2*a1(1,1)
    a2(1,5)=a2(1,5)+fac2*a2(1,1)
    a3(1,5)=a3(1,5)+fac2*a3(1,1)
    a1(1,1)=0,
    a2(1,1)=0,
    a3(1,1)=0,
    fac4=-dvout/dvin
c    if(sfjrx,eq,1,and,sf1zx,eq,1)fac4=r(ix-1,jx-1)*b(ix-1,jx-1)/
c    (r(ix,jx-1)*b(ix,jx-1))
    a1(kxp,5)=a1(kxp,5)+fac4*a1(kxp,9)
    a2(kxp,5)=a2(kxp,5)+fac4*a2(kxp,9)
    a3(kxp,5)=a3(kxp,5)+fac4*a3(kxp,9)
    a1(kxp,9)=0,
    a2(kxp,9)=0,
    a3(kxp,9)=0,
    i=2
    do 11 j=2,jx-1
    if(sf11,eq,1,)sf11=r(2,j)*b(2,j)/(r(1,j)*b(1,j))
    k1=i-1+(j-2)*(ix-2)
    k2=j-1+(jx-2)*(i-2)
    k=.5*(1+isw)*k1+.5*(1-isw)*k2
    do 11 m=2,8,3
    a1(k,m)=a1(k,m)+sf11*a1(k,m-1)
    a2(k,m)=a2(k,m)+sf11*a2(k,m-1)
    a3(k,m)=a3(k,m)+sf11*a3(k,m-1)
11    continue
13    continue
    i=ix-1
    do 12 j=2,jx-1
c    bc(j)=0,
c    if(psi(j).lt.(psih-dpsih))bc(j)=1,
c    if(psi(j).ge.(psih-dpsih),and,psi(j).le.(psih+dpsih))
c    bc(j)=.5*(tanh((-psi(j)+psih)/dpsih)+1)
    rbt=r(ix-1,j)*b(ix-1,j)
    rbt1=r(ix,j)*b(ix,j)
    down2=8.*bc(j)*uuzh(ix-1)/du+3.*(1.-bc(j))
    up2=1.-bc(j)
    gm2=up2/down2
    up1=-up2*rbt1*.75+bc(j)*rbt*uuzh(ix-1)/du
    down1=(bc(j)*uuzh(ix-1)/du+up2*3./8.)*rbt1
    gm1=up1/down1
    delco1(j)=gm1
    delco2(j)=gm2
    up=(bc(j)*r(ix-1,j)*b(ix-1,j)*uuzh(ix-1)/du+rbt*(bc(j)-1.))
    down=(bc(j)*r(ix,j)*b(ix,j)*uuzh(ix-1)/du+rbt*(-bc(j)+1.))
    sf1zx=up/down
    k1=i-1+(j-2)*(ix-2)
    k2=j-1+(jx-2)*(i-2)
    k=.5*(1+isw)*k1+.5*(1-isw)*k2
    do 12 m=2,8,3
    a1(k,m)=a1(k,m)+gm1*a1(k,m+1)
    a1(k,m-1)=a1(k,m-1)+gm2*a1(k,m+1)
    a2(k,m)=a2(k,m)+gm1*a2(k,m+1)

```

```

a2(k,m-1)=a2(k,m-1)+gm2*a2(k,m+1)
a3(k,m)=a3(k,m)+gm1*a3(k,m+1)
a3(k,m-1)=a3(k,m-1)+gm2*a3(k,m+1)
12 continue
20 continue
i=2
do 21 j=2,jx-1
k1=i-1+(j-2)*(ix-2)
k2=j-1+(jx-2)*(i-2)
k=.5*(1+isw)*k1+.5*(1-isw)*k2
do 21 m=1,7,3
a1(k,m)=0.
a2(k,m)=0.
a3(k,m)=0.
21 continue
i=ix-1
do 22 j=2,jx-1
k1=i-1+(j-2)*(ix-2)
k2=j-1+(jx-2)*(i-2)
k=.5*(1+isw)*k1+.5*(1-isw)*k2
do 22 m=3,9,3
a1(k,m)=0.
a2(k,m)=0.
a3(k,m)=0.
22 continue
j=2
do 31 i=2,ix-1
k1=i-1+(j-2)*(ix-2)
k2=j-1+(jx-2)*(i-2)
k=.5*(1+isw)*k1+.5*(1-isw)*k2
do 30 m=4,6
a1(k,m)=a1(k,m)+sfj1*a1(k,m-3)
a2(k,m)=a2(k,m)+sfj1*a2(k,m-3)
a3(k,m)=a3(k,m)+sfj1*a3(k,m-3)
30 continue
b1(k,2)=b1(k,2)+sfj1*b1(k,1)
31 continue
32 continue
j=jx-1
fac5=sfjrx
if(sfjrx.eq.-1) fac5=fac5*dvout/dvin
do 35 i=2,ix-1
k1=i-1+(j-2)*(ix-2)
k2=j-1+(jx-2)*(i-2)
k=.5*(1+isw)*k1+.5*(1-isw)*k2
do 34 m=4,6
a1(k,m)=a1(k,m)+fac5*a1(k,m+3)
a2(k,m)=a2(k,m)+fac5*a2(k,m+3)
a3(k,m)=a3(k,m)+fac5*a3(k,m+3)
34 continue
b1(k,2)=b1(k,2)+fac5*b1(k,3)
35 continue
40 continue
j=2
do 45 i=2,ix-1
k1=i-1+(j-2)*(ix-2)
k2=j-1+(jx-2)*(i-2)
k=.5*(1+isw)*k1+.5*(1-isw)*k2
do 44 m=4,6
a1(k,m-3)=0.

```

```
a1(k,m-3)=0.
```

```

a3(k,m-3)=0,
44 continue
b1(k,1)=0,
45 continue
j=jx-1
do 48 i=2,ix-1
k1=i-1+(j-2)*(ix-2)
k2=j-1+(jx-2)*(i-2)
k=.5*(1+isw)*k1+.5*(1-isw)*k2
do 47 m=4,6
a1(k,m+3)=0,
a3(k,m+3)=0,
a2(k,m+3)=0,
47 continue
b1(k,3)=0,
48 continue
return
end

```

```
c*****
```

```

subroutine bvcal(bs,zs,as,als,z,n,zcc,bcc,znorm,bv,bvp,bvpp,
1 b4,b5,z4,z5,nc)
dimension bs(1),zs(1),as(1),als(1),z(1),bv(1),bvp(1),bvpp(1)
call bcccal(bs,zs,as,als,z,n,bv,bvp,bvpp,b4,b5,z4,z5,nc)
do 15 i=1,n
bcorr=bcc
bcorrpp=0.
bcorrpp=0.
tanhyp=tanh((z(i)-zcc)/znorm)
bcorrpp=-bcc*.5*(1.-tanhyp**2)/znorm
bcorrpp=+bcc*tanhyp*(1.-tanhyp**2)/znorm**2
18 bcorr=bcc*.5*(1.-tanhyp)
bv(i)=bv(i)+bcorr
bvp(i)=bvp(i)+bcorrpp
bvpp(i)=bvpp(i)+bcorrpp
if(z(i).eq.z4)b4=b4+bcorr
if(z(i).eq.z5)b5=b5+bcorr
15 continue
return
end

```

```
c*****
```

```

subroutine bcccal(bs,zs,as,als,z,n,bv,bvp,bvpp,b4,b5,z4,z5,nc)
dimension z(n),bv(n),bvp(n),bvpp(n)
dimension bs(3),zs(3),as(3),als(3),alpha(3,3),ak(3),is(3)

```

```
15
```

```
c....compute matrix elements
```

```

do 10 j=1,nc
do 10 i=1,nc
alpha(i,j)=bfun(zs(i),zs(j),als(j),as(j),0)
10 continue

```

```
c....determinant
```

```
if(nc.eq.2)alpha(3,3)=1.
```

```

det=alpha(1,1)*(alpha(2,2)*alpha(3,3)-alpha(3,2)*alpha(2,3))
1 -alpha(1,2)*(alpha(2,1)*alpha(3,3)-alpha(3,1)*alpha(2,3))
2 +alpha(1,3)*(alpha(2,1)*alpha(3,2)-alpha(3,1)*alpha(2,2))

```

c...solution

```

ak(1)=(bs(1)*(alpha(2,2)*alpha(3,3)-alpha(3,2)*alpha(2,3))
1      -bs(2)*(alpha(1,2)*alpha(3,3)-alpha(3,2)*alpha(1,3))
2      +bs(3)*(alpha(1,2)*alpha(2,3)-alpha(2,2)*alpha(1,3)))
3      /det
ak(2)=(-bs(1)*(alpha(2,1)*alpha(3,3)-alpha(3,1)*alpha(2,3))
1      +bs(2)*(alpha(1,1)*alpha(3,3)-alpha(3,1)*alpha(1,3))
2      -bs(3)*(alpha(1,1)*alpha(2,3)-alpha(2,1)*alpha(1,3)))
3      /det
ak(3)=(bs(1)*(alpha(1,2)*alpha(2,3)-alpha(1,3)*alpha(2,2))
1      -bs(2)*(alpha(1,1)*alpha(2,3)-alpha(1,3)*alpha(2,1))
2      +bs(3)*(alpha(1,1)*alpha(2,2)-alpha(1,2)*alpha(2,1)))
3      /det

```

c....fields and derivatives

```

do 50 i=1,n
  bv(i)=0.
  bvp(i)=0.
  bvpp(i)=0.
  do 50 j=1,nc
    bv(i)=bv(i)+ak(j)*bfun(z(i),zs(j),als(j),as(j),0)
    bvp(i)=bvp(i)+ak(j)*bfun(z(i),zs(j),als(j),as(j),1)
    bvpp(i)=bvpp(i)+ak(j)*bfun(z(i),zs(j),als(j),as(j),2)
  continue
50

```

c....minima and their positions

```

do 70 j=1,nc
do 60 i=2,n
if(zs(j).lt.z(i)) go to 63
60 continue
63 is(j)=i-1
70 continue
b4=aminaf(bv,is(1),is(2),1,i4,amin)
z4=z(i4)
if(ncoll.eq.2) return
b5=aminaf(bv,is(2),is(3),1,i5,amin)
z5=z(i5)

```

return
end

```

c *****
      function bfun(z,zx,a1,a,ind)
      ind1=ind+1
      up=zx+.5*a1
      um=zx-.5*a1
      go to (10,20,30),ind1
10    t1=gfun(z,up,a)
      t2=gfun(z,um,a)
      go to 40
20    t1=gfun1(z,up,a)
      t2=gfun1(z,um,a)
      go to 40
30    t1=gfun2(z,up,a)
      t2=gfun2(z,um,a)
40    bfun=(t1-t2)/a**2
      return

```

```

      end
c*****
      function gfun(z,u,a)
      x1=u-z
      x2=sqrt(x1**2+a**2)
      gfun=x1/x2
      return
      end
c*****
      function gfun1(z,u,a)
      x1=u-z
      x2=(x1**2+a**2)**1.5
      gfun1=-a**2/x2
      return
      end
c*****
      function gfun2(z,u,a)
      x1=u-z
      x2=(x1**2+a**2)**2.5
      gfun2=-3.*x1*a**2/x2
      return
      end

      subroutine comat(abar,nd)

c.....transforms the elements of the a1(k,m) array into into the
c..... elements of the compressed column matrix abar which will be
c..... operated upon by banfac and bansol.

c..... insert storage cliches here
      use param
      use fstor
      use matrix
      use const

      dimension abar(kxp,1)

      kxx=kxp
      len=nd*kxp
      call bcast(abar(1,1),0.,len)
      do 10 k=1,kxx
      do 10 m=1,9
      lp1=m+((m-1)/3)*(ihbw-4)
      lp2=1+mod(m-1,3)*(ihbw-1)+(m-1)/3
      lp=.5*(1+isw)*lp1+.5*(1-isw)*lp2
      abar(k,lp)=a1(k,m)
1? continue
      return
      end

      subroutine constant

c..... insert storage cliches here
      use param
      use fstor
      use matrix
      use const

      gam1=.25*(3*bias+1)

```

```

gam2=.25*(1-bias)
ip=.5*(ix-2)
jp=.5*(jx-2)
kp1=ip-1+(jp-2)*(ix-2)
kp2=jp-1+(ix-2)*(ip-2)
kplot=.5*(1+isw)*kp1+.5*(1-isw)*kp2
if(kplotm.ne.0)kplot=kplotm
return
end

```

```

subroutine curvecon

```

```

c.... calculates constants necessary for curvature driven
c.... flute mode case.

```

```

c..... insert storage cliches here
use param
use const
real klbsq

```

```

c..... input for curvature driven flute case
data echarg/4.8e-10/, en0/1.00e+12/, b0/1.e4/, amass/3.34e-24/
c , ceo/3.e10/, stable/.4/, fourpi/12.56637/, pi/3.1415926/
c , delrho/.05/,dtrel/.02/,xm/3.8317/, theta0/1.570796/
namelist/curve/b0,beta0,delrho,stable,en0,echarg,ib,rw,xm
c ,zol,dtrel,theta0
call ddi(curve,2,3,1)
if(nold.ne.1) call ddo(curve,100,0,1)
zmax=zol*lb
p0=beta0*b0**2*.5
psimax=rw**2*b0*.5
omeg0sq=b0**2/(en0*amass*lb**2)
omeg0=sqrt(omeg0sq)
ag1=1.-2.*delrho/xm**2
klbsq=0.
if(kzs.ne.0)klbsq=(pi/(2.*zol))**2
delomeg=0.
if(sf8.ne.0)
1 delomeg=stable*((beta0/xm**2-klbsq/mm**2)/(ag1*sf8)+1.-sf6*ag1/
2 sf8)*omeg0sq
omeg1=omeg0+sqrt(delomeg)
omeg2=omeg1-2*sqrt(delomeg)
dt=dtrel/omeg1
en1=2.*en0*delrho/rw**2
u(ix)=zmax
v(jx)=psimax
du=u(ix)/(ix-1.5)
dv=v(jx)/(jx-1.5)
besarg=(xm/rw)
c..... calculate analytic growth rate
tsf6=tan(theta0*.5)*sf6
radical=((ag1*mm*tsf6)**2*(omeg1+omeg2)**2-4.*((omeg1*omeg2*ag1
c *sf8+omeg0sq*beta0/xm**2)*mm**2-klbsq*omeg0sq))
if(radical.lt.0)go to 5
root=sqrt(radical)
omana1=ag1*tsf6*mm*(omeg1+omeg2)*.5+root*.5
omana2=omana1-root
groana=0.
return

```

```

5      continue
      omana1=ag1*tsf6*mm*(omeg1+omeg2)*.5
      groana=sqrt(-radical)*.5
      omana2=0,
      return
    end
    subroutine energy

c,...  energy calculates the total first order energy (tenergy) as
c,...  an integral (using trapazoidal quadrature) over the volume,
c,...  every nen'th time step, using two consecutive time steps to
c,...  determine the value at n+.5 ,

      use param
      use fstor
      use matrix
      use const

      dimension rxi(jrx),rxr(jrx),drxr(izx,jrx),drxi(izx,jrx),quad(12)
c ,quad(4),psuma(jrx,12),psumb(jrx,4),dum1(izx),dum2(izx)
c ,xrg(izx),xig(izx),xrgo(izx),xigo(izx),drxro(izx,jrx)
c ,drxio(izx,jrx),dxrz2(izx),dxlz2(izx)

      ltt=ltt+1
      if((ltt.eq.2).or.(nen.eq.1))then
        lmove=jx*ix
        call zmovewrd(drxro,drxr,lmove)
        call zmovewrd(drxio,drxi,lmove)
        tquadoo=tquado
        tquado=tquad
        tquad2oo=tquad2o
        tquad2o=tquad2
        tquad3oo=tquad3o
        tquad3o=tquad3
        tquad4oo=tquad4o
        tquad4o=tquad4
      end if
      do 9 i=2,ix-1
        do 5 j=2,jx-1
          k1=i-1+(j-2)*(ix-2)
          k2=j-1+(jx-2)*(i-2)
          k=.5*((1+isw)*k1+(1-isw)*k2)
          rxr(j)=r(i,j)*xro(k)
          rxi(j)=r(i,j)*xio(k)
5          continue
          call ddpsi(rxr,drxr,i)
          call ddpsi(rxi,drxi,i)
9          continue

c      loop 100 evaluates z quadrature

        do 100 j=2,jx-1
          do 10 i=2,ix-1
            k1=i-1+(j-2)*(ix-2)
            k2=j-1+(jx-2)*(i-2)
            k=.5*((1+isw)*k1+(1-isw)*k2)
            if((ltt.eq.2).or.(nen.eq.1))then
              xrgo(i)=xroo(k)
              xigo(i)=xioo(k)
            end if

```

```

      xrg(i)=xro(k)
      xlg(i)=xlo(k)
10  do 211 i=2,ix-1
211 dum2(i)=r(i,j)*b(i,j)*xrg(i)
      call ddz(dum2,dxrz2)
      do 212 i=2,ix-1
212 dum2(i)=r(i,j)*b(i,j)*xlg(i)
      call ddz(dum2,dxiz2)
      do 11 i=2,ix-1
11  dum1(i)=(1./uuz(i))*qub(i,j)*dxrz2(i)/((r(i,j)*b(i,j))**2)
      psuma(j,1)=psum(dum1,ix)/vpsi(j)
      do 12 i=2,ix-1
12  dum1(i)=(1./uuz(i))*qub(i,j)*dxiz2(i)/((r(i,j)*b(i,j))**2)
      psuma(j,2)=psum(dum1,ix)/vpsi(j)
      do 13 i=2,ix-1
13  dum1(i)=b(i,j)*drxi(i,j)
      call ddz(dum1,dum2)
      do 14 i=2,ix-1
14  dum1(i)=(1./uuz(i))*qub(i,j)*r(i,j)**2*dum2(i)
      psuma(j,3)=psum(dum1,ix)/vpsi(j)/mm**2
      do 15 i=2,ix-1
15  dum1(i)=b(i,j)*drxr(i,j)
      call ddz(dum1,dum2)
      do 16 i=2,ix-1
16  dum1(i)=(1./uuz(i))*qub(i,j)*r(i,j)**2*dum2(i)
      psuma(j,4)=psum(dum1,ix)/vpsi(j)/mm**2
      do 17 i=2,ix-1
17  dum1(i)=(1./uuz(i))*xrg(i)**2*r(i,j)*(qubv(i,j)*rzz(i,j)
c +r(i,j)*ering(i,j)/b(i,j))
      psuma(j,5)=psum(dum1,ix)/vpsi(j)
      do 18 i=2,ix-1
18  dum1(i)=(1./uuz(i))*xlg(i)**2*r(i,j)*(qubv(i,j)*rzz(i,j)
c +r(i,j)*ering(i,j)/b(i,j))
      psuma(j,6)=psum(dum1,ix)/vpsi(j)
      do 19 i=2,ix-1
19  dum1(i)=(1./uuz(i))*xrg(i)**2*yyy(i,j)/b(i,j)
      psuma(j,7)=mm**2*psum(dum1,ix)/vpsi(j)
      do 20 i=2,ix-1
20  dum1(i)=(1./uuz(i))*xlg(i)**2*yyy(i,j)/b(i,j)
      psuma(j,8)=mm**2*psum(dum1,ix)/vpsi(j)
      do 21 i=2,ix-1
21  dum1(i)=(1./uuz(i))*yyy(i,j)*xrg(i)*r(i,j)*drxr(i,j)
      psuma(j,9)=-2.*psum(dum1,ix)/vpsi(j)
      do 22 i=2,ix-1
22  dum1(i)=(1./uuz(i))*yyy(i,j)*r(i,j)**2*b(i,j)*drxr(i,j)**2
      psuma(j,10)=psum(dum1,ix)/vpsi(j)
      do 23 i=2,ix-1
23  dum1(i)=-(1./uuz(i))*yyy(i,j)*xlg(i)*r(i,j)*drxi(i,j)
      psuma(j,11)=2*psum(dum1,ix)/vpsi(j)
      do 24 i=2,ix-1
24  dum1(i)=(1./uuz(i))*yyy(i,j)*r(i,j)**2*b(i,j)*drxi(i,j)**2
      psuma(j,12)=psum(dum1,ix)/vpsi(j)
      if((lft.eq.2).or.(nen.eq.1))then
      do 30 i=2,ix-1
30  dum1(i)=(1./uuz(i))*rho(i,j)*((xrg(i)-xrgo(i))/dt)**2/b(i,j)
      psumb(j,1)=psum(dum1,ix)/vpsi(j)
      do 31 i=2,ix-1
31  dum1(i)=(1./uuz(i))*rho(i,j)*((xlg(i)-xigo(i))/dt)**2/b(i,j)
      psumb(j,2)=psum(dum1,ix)/vpsi(j)
      do 32 i=2,ix-1

```



```

32   dum1(i)=(1./uuz(i))*(r(i,j)*(drxr(i,j)-drxro(i,j))/dt)**2*b(i,j)
c   *rho(i,j)
   psumb(j,3)=psum(dum1,ix)/vpsi(j)/mm**2
   do 33 i=2,ix-1
33   dum1(i)=(1./uuz(i))*(r(i,j)*(drxi(i,j)-drxio(i,j))/dt)**2*b(i,j)
c   *rho(i,j)
   psumb(j,4)=psum(dum1,ix)/vpsi(j)/mm**2
   end if
100  continue
   do 120 l=1,12
120  quad(l)=psum(psuma(1,l),jx)
   tquad=ssum(12,quad,1)*dv*du
   tquad2=0.
   do 121 l=1,4
121  tquad2=tquad2+quad(l)
   tquad3=quad(5)+quad(6)
   tquad4=0.
   do 122 l=7,12
122  tquad4=quad(l)+tquad4
   if((l==2).or.(l==1))then
   tquad1=0.
   lee=lee+1
   do 130 l=1,4
   quad1(l)=psum(psumb(1,l),jx)
130  tquad1=tquad1+quad1(l)
   enpot(lee)=(3.*tquad+6.*tquado-tquadoo)/8.
   enkin(lee)=tquad1*dv*du
   tenergy(lee)=abs(enpot(lee)+enkin(lee))
   timengy(lee)=time(n)
   enbend(lee)=(3.*tquad2+6.*tquad2o-tquad2oo)/8.*du*dv
   encurve(lee)=abs(3.*tquad3+6.*tquad3o-tquad3oo)/8.*du*dv
   tenrel(lee)=abs(tenergy(lee))/(enkin(lee)+abs(enpot(lee)))*2
   enfir(lee)=abs(3.*tquad4+6.*tquad4o-tquad4oo)/8.*du*dv
   ltt=0
   end if
   return
   end

   function psum(f,k)

   dimension f(1)

   nq=k-4
   psum=ssum(nq,f(3),1)+.5*(f(2)+f(k-1))
   return
   end

   subroutine ddz(f1,f2)

   use const
   dimension f1(1),f2(1)

   do 10 i=4,ix-1
10  f2(i-1)=((f1(i)-f1(i-2))/(2.*du)*uuz(i-1))**2
   f2(2)=(((1.-f1(1))*f1(2)*uuzh(1)+(f1(3)-f1(2))*uuzh(2))/(du*2))**2
   f2(ix-1)=((-1.-f1(ix))*f1(ix-1)*uuzh(ix-1)+(f1(ix-1)-f1(ix-2))
c   *uuzh(ix-2))/(du*2))**2
   return
   end

```

```

      subroutine ddpsi(f1,f2,i)
      use const
      dimension f1(1),f2(1:ix,1)

      do 10 j=4,jx-1
10      f2(i,j-1)=(f1(j)-f1(j-2))/(2.*dv)*vpsi(j-1)
          f2(i,2)=(2.*f1(2)*vpsih(1)+(f1(3)-f1(2))*vpsih(2))/(dv*2)
          f2(i,jx-1)=(-(1.-fjrx)*f1(jx-1)*vpsih(jx-1)+(f1(jx-1)-
c f1(jx-2))*vpsih(jx-2))/(dv*2)
          return
      end
      subroutine equil

c.....special case equilibrium, 0 beta, 0 pressure, rho=const.
c.....test case 1
c.....set up 1/4/82 by r. freis

c.....insert cliché storage here
      use param
      use matrix
      use const
      use fstor

      data rho0/1.e12/,b0/1.e4/,azm/1./,apsim/1./

      do 10 j=1,jx
      do 10 i=1,ix
          uz=uuz(i)
          rho(i,j)=rho0
          b(i,j)=b0
          r(i,j)=sqrt(2.*abs(psi(j))/b0)
          xxx(i,j)=0.
          yyy(i,j)=0.
          qub(i,j)=b0
10      continue
      return
      end
      subroutine equilcur

c.....equilibrium for curvature driven flute mode case.

c.....insert storage clichés here.
      use param
      use const
      use fstor

      do 10 i=1,ix
      do 10 j=1,jx
c..... special b(i,j) to test b.c. on flute test case
          b(i,j)=b0
          r(i,j)=sqrt(2*psi(j)/b(i,j))
          rho(i,j)=(en0-en1*r(i,j)**2*.5)*amass
          xxx(i,j)=rho(i,j)*(omeg1+omeg2)*sf6
          yyy(i,j)=rho(i,j)*(-omeg1*omeg2)*sf8
          qub(i,j)=b(i,j)
          qv(i,j)=p0/psi(jx)
10      continue
      return
      end

```

```

      subroutine equilrot
c..... sets up equilibrium for rigid rotor, test case 2 ,
c.....flora3 adds cold plasma halo to equilibrium density
c..... insert cliché storage here
      use param
      use const
      use fstor

      psi0=b0*r0sq*.5/sqrt(fourn1)
      omegr=ratrod*omegst*(1.-enbar/en0)
      foursq=sqrt(fourn1)
      do 5 i=1,ix
      uz=uuz(i)
      do 5 j=1,jx
      fac=exp(psi(i,j)/psi0)/sqrt(beta0)
      b(i,j)=b0*sqrt(fac**2-1.)/(fac*foursq)
      rho(i,j)=en0*amass/(beta0*fac**2)+enbar*amass
      beta=1/fac**2
      arg1=fac+sqrt(fac**2-1.)
      acosh=alog(arg1)
      r(i,j)=r0*sqrt(-cr+acosh)
      qub(i,j)=b(i,j)
      omegstr=omegst*(1.-enbar*amass/rho(i,j))
      entest=enbar*amass
      if(entest.ge.rho(i,j))omegstr=0.
      omegxb=(1.+ratrod)*omegstr
      omeggb=+beta*omegstr*.5/(1.-beta)
      xxx(i,j)=rho(i,j)*(2.*omegxb+omeggb-omegst)
      yyy(i,j)=-rho(i,j)*(omegxb+omeggb)*(omegxb-omegst)
      xxx(i,j)=xxx(i,j)*flr
      yyy(i,j)=yyy(i,j)*flr
5      continue
      do 30 i=1,ix
      do 30 j=2,jx-1
      qv(i,j)=.5*(qub(i,j+1)*b(i,j+1)-qub(i,j-1)*b(i,j-1))
30      continue
      return
      end
      subroutine equilrm
c.....equilibrium for tandem mirror electron ring plug
c.....insert storage clichés here
      use param
      use const
      use fstor

      data epsb/1.e-2/, p30/.01/
c..... loop 10, calculate p2 p1 and b and dp2/dpsi and dp1/dpsi
      wp2=p2wide*.5*psi0
      psibar=-be1*.5/ce1
      psistr=-ce1/(3.*de1)
      psibar=psistr+sqrt(psistr**2-be1/(3.*de1))
      plmax=ae1+be1*psibar+ce1*psibar**2+de1*psibar**3

      do 10 j=1,jx
      psir=psi(j)/psim
      psire=psi(j)/psime

```

```

pe1=(ae1+be1*psire+ce1*psire**2+de1*psire**3)*hpmo(j)
pe2=.5*(1.-tanh((psi(j)-psi0e)/wp2e))*(1.-hpmo(j))
p1(j)=(pe1+pe2)/p1max
dp1dpsi(j)=(de1*3.*psire**2+
c be1+2.*ce1*psire)/(psime*p1max)*hpmo(j)-.5*(1.-tanh
c ((psi(j)-psi0e)/wp2e)**2)/(wp2e*p1max)*(1.-hpmo(j))
p2t(j)=(1.-tanh((psi(j)-psi0)/wp2))*5
dp2dpst(j)=-.5*(1.-tanh((psi(j)-psi0)/wp2)**2)/wp2
p3(j)=(p3a+p3b*psir+p3c*psir**2+p3d*psir**3)*hp0(j)
dp3dpsi(j)=(p3b+2.*p3c*psir+3.*p3d*psir**2)*hp0(j)/psim
p2(j)=p2t(j)+p3(j)*dip
dp2dpsi(j)=dp2dpst(j)+dp3dpsi(j)*dip
hflr(j)=1.

c..... set p2 and p3 constant at large psi
if (p2(j).le.p2flag)then
p2(j)=p2floor
p1(j)=p1floor
dp2dpsi(j)=0.
dp1dpsi(j)=0.
hflr(j)=0.
end if
do 11 i=1,ix
c if(abp(i).eq.0)then
c b(i,j)=bvac(i)
c go to 11
c end if
u1=p2(j)*abp(i)+p1(j)*abf(i)
u2=p2(j)*bbp(i)+.5+p1(j)*bbf(i)
u3=p2(j)*cbp(i)-bvac(i)**2*.5+p1(j)*cbf(i)
c..... at,bt,ct coefficients for low density expansion
at=u1
bt=u2-.5
ct=u3+bvac(i)**2*.5
bv=bvac(i)
ordera2=4.*ct*bt+4.*(bt*bv)**2+8.*at*ct*bv**2+12.*at*bt*bv**4
c +8*at**2*bv**6
bsq3=bv**2-2.*(ct+bt*bv**2+at*bv**4)+ordera2
if((ordera2/bsq3).le.epsp) then
b(i,j)=sqrt(bsq3)
else

t1=-u2*.5/u1
radic2=(t1**2-u3/u1)
round=abs(radic2)/(t1**2+abs(u3/u1))
if (round.le.1.e-6)radic2=0.
radic=sqrt(radic2)
bsq1=t1+radic
bsq2=t1-radic
bsq=bsq1
if(t1.gt.radic)bsq=bsq2
b(i,j)=sqrt(bsq)
end if
11 continue
10 continue

c.....loop 15, calculate phi, the electric potential
do 15 i=1,ix
arg1=((z(i)-z1)/(z1-z0))**2*(-xpot)*(1.-hzt0(i))
arg2=wpot*((z(i)-z0)/(z0-z2))**2

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15   phi1(i)=phice*exp(arg1)+phipl/cosh(arg2)
c.....special phi1 for high mm rotation mode test
      do 17 i=1,ix
        phi1(i)=phice*(1.-dphi*(1.-hzt0(i)))
17   continue
      do 16 j=2,jx
c       phi2(j)=(1.-(psi(j)*hp3(j)/psi3)**ypot)*hp3(j)
c..... special phi2 for constant omegexb
        phi2(j)=er*psi(j)
16   continue
c.....loop 20, calculate pperp, ppar and db/dpsi
c.....first calculate d coef. for ppar
        dp1=(ap1*bmax**2*4./3.+2.*bp1)*bmax
        dp2=(ap2-ap1)*bm1**3/3.+(bp2-bp1)*bm1+dp1+(cp1-cp2)
c       /bm1
        dp3=(ap3-ap2)*bv2**3/3.+(bp3-bp2)*bv2+dp2+(cp2-cp3)
c       /bv2
        dt0=-8.*ct0/(bmax*3.)
        dt1=-8.*ct1/(bv2*3.)
        dt2=-8.*ct2/(3.*bv2)
        dt3=-8.*ct3/(3.*bm1)
        ppt=1./(1.-1./rp1**2)**2
        dter1=(-8./3-4.*(bv3/bm2)**3/3.+4.*(bv3/bm2))/bm2
        dter2=-alsi*ppt*dter1*psisi
        do 20 i=1,ix
          do 20 j=1,jx
            if(dter2.eq.0.)dter2=0.
            b2=b(i,j)**2
            b4=b2**2
            dter=dp1*hzp1(i)+dp2*hzp2(i)+dp3*hzp3(i)+dt1*hzt1(i)
c          +dt2*hzt2(i)-dt3*hzt3(i)+dt0*hzt0(i)
            ppart=(-abp(i)/3.*b4-bbp(i)*b2+cbp(i)+dter*b(i,j))
            ppert=(abp(i)*b4+bbp(i)*b2+cbp(i))
            pperte=(abq(i)*b4+bbq(i)*b2+cbq(i))
            pperp(i,j)=p2(j)*ppert
            ppar(i,j)=p2(j)*ppart
            qub(i,j)=(b2+ppar(i,j)-pperp(i,j))/b(i,j)
c          rho(i,j)=amass*(p2(j)*(ebp(i)*b4+bbp(i)*b2+gbp(i))+
c          c ncenter*cold*(1.+(rcwall-1.)*psi(j)/psi(jx)))
c..... special rho for exb rotation case
            rho(i,j)=amass*ncenter*exp(lexrho*psi(j)/psi(jx))
            fac=b(i,j)*(1.+2.*p1(j)*(2.*abf(i)*b2+bbf(i)))
            fac1=dp1dpsi(j)*(abf(i)*b4+bbf(i)*b2+cbf(i))
            fac2=dp2dpsi(j)*(abp(i)*b4+bbp(i)*b2+cbp(i))
            fac3=p2(j)*(2.*abp(i)*b2+bbp(i))+p1(j)*(2.*abf(i)*b2+bbf(i))
            dbdpsi(i,j)=-(fac1+fac2)/(b(i,j)*(1.+fac3*2.))
            dp2db=b(i,j)*(abp(i)*b2*2.+bbp(i))*2.
            dp2dbe=b(i,j)*(abq(i)*b2*2.+bbq(i))*2.
            dp3db=dp2db-4.*abp(i)*b(i,j)**3/3.-2.*bbp(i)*b(i,j)
c          +dter
            qubv(i,j)=(-dp2dpsi(j)*(ppert+ppart)-p2(j)*dp3db*dbdpsi(i,j))
            pharg=psi(j)*hp3(j)
c          dphidpsi=phi1(i)*ypot*(pharg/psi3)**(ypot-1.)/
c          c (-psi3)*hp3(j)
c..... special dphidpsi for constant omegexb
            dphidpsi=phi1(i)*er/cee
            epsi(i,j)=-dphidpsi
            dpdpsi=pperte*dp2dpsi(j)+p2(j)*dp2dbe*dbdpsi(i,j)
            omegci=echarg*b(i,j)/(amass*cee)
            unit=1./sqrt(4.*pi)

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      omegstr=-unit*b(i,j)*dps1/(omegci*rho(i,j))
      omeggb=unit*pperp2(j)*dbdps1(i,j)/(omegci*rho(i,j))
      omegexb=-cee*dphidps1
      xxx(i,j)=hflr(j)*rho(i,j)*(omeggb-omegstr+2.*omegexb)*sf6
      yyy(i,j)=-hflr(j)*rho(i,j)*(omegexb+omeggb)*(omegexb-omegstr)*sf8

      ppertr=(abf(i)*b4+bbf(i)*b2+cbf(i))
      pperps(i,j)=pperp(i,j)+p1(j)*ppertr
      pperpe(i,j)=p1(j)*ppertr
20      continue

c.....check for negative pressure, and terminate prob. if necessary
      do 25 j=1,jx
      do 25 i=1,ix
      if(pperp(i,j).lt.0.or.pperpe(i,j).lt.0.)then
      write(59,101)
101  format('problem terminated due to negative pressure' )
      call exit(1)
      end if
25      continue

c....calculate diagnostic on perpendicular pressure balance
      do 80 i=1,ix
      do 80 j=1,jx
      errprp(i,j)=(b(i,j)**2-bvac(i)**2+2.*pperps(i,j))/bvac(i)**2
80      continue

c....calculate diagnostic on parallel pressure balance
      do 81 j=2,jx
      do 81 i=2,ix-1
      delb=b(i+1,j)-b(i-1,j)
      if(abs(delb).lt,epsb)then
      abar=-abp(i)*5./3.
      bbar=-3.*bbp(i)
      dter4=dp1*hzp1(i)+dp2*hzp2(i)+dp3*hzp3(i)+dt1*hzt1(i)
      c +dt2*hzt2(i)-dt3*hzt3(i)
      dbar=dter4*2.
      factor=-(abar*b(i,j)**4+bbar*b(i,j)**2+cbp(i)+dbar*b(i,j))
      else
      factor=(ppar(i+1,j)*b(i+1,j)-ppar(i-1,j)*b(i-1,j))
      c /delb
      end if
81  errprl(i-1,j-1)=(pperp(i,j)-2.*ppar(i,j)+factor)/bvac(i)**2*2

c.....cal. r and rzz*r
      do 40 i=1,ix
      sum1=0.
      sum2=0
      sum3=0.
      sum4=0.
      sum5=0.
      do 40 j=2,jx
      dps=dps1(j)
      psib=psi(j-1)
      if(j.eq.2)then
      dps=psi(2)
      psib=0.
      end if
      if(i.gt.2)go to 31
      call bcast(hpk0(1),0.,kin)
      call bcast(hpk12(1),0.,kin)

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call bcast(hpkm(1),0.,kin)
call bcast(hpkme(1),0.,kin)
do 29 k=1,kin
  psik=psib+dps*(k-1)/(kin-1)
  if(psik.lt.psi0)hpko(k)=1.
  if(psik.lt.psim)hpkm(k)=1.
  if(psik.lt.psim)hpkm(k)=1.
  if(psik.lt.psi2.and.psik.gt.psi1)hpkl2(k)=1.
29 continue
c..... call p1 (p1k) at intermediate points in dps1 interval
  do 30 k=1,kin
    psik=psib+dps*(k-1)/(kin-1)
    psikr=psik/psim
    psikre=psik/psime
    p1ek=(a01+b01*psikre+c01*psikre**2+d01*psikre**3)*hpkm(k)
    p2ek=.5*(1.-tanh((psik-psi0e)/wp2e))*(1.-hpkm(k))
    p1k(k,j)=(p1ek+p2ek)/p1max
    p3k=(p3a+p3b*psikr+p3c*psikr**2+p3d*psikr**3)*hpko(k)
    p2k(k,j)=(1.-tanh((psik-psi0)/wp2))*0.5 +p3k*dip
30 continue
31 continue
  do 32 k=1,kin
    if(p2k(k,j).le.p2flag)then
      p2k(k,j)=p2floor
      p1k(k,j)=p1floor
    end if
    u1=p2k(k,j)*abp(i)+p1k(k,j)*abf(i)
    u2=p2k(k,j)*bbp(i)+p1k(k,j)*bbf(i)+.5
    u3=p2k(k,j)*cbp(i)+p1k(k,j)*cbf(i)-bvac(i)**2*.5
c..... at, bt, ct coefficients for low density expansion
    at=u1
    bt=u2-.5
    ct=u3+bvac(i)**2*.5
    bv=bvac(i)
    ordera2=4.*ct*bt+4.*(bt*bv)**2+8.*at*ct*bv**2+12.*at*bt*bv**4
c +8*at**2*bv**6
    bsq3=bv**2-2.*(ct+bt*bv**2+at*bv**4)+ordera2
    if((ordera2/bsq3).le.epsp) then
      bk=sqrt(bsq3)
    else
      t1=-u2*.5/u1
      radic2=(t1**2-u3/u1)
      round=abs(radic2)/(t1**2+abs(u3/u1))
      if(round.le.1.e-6)radic2=0.
      radic=sqrt(radic2)
      bsq1=t1+radic
      bsq2=t1-radic
      bsq=bsq1
      if(t1.gt.radic)bsq=bsq2
      bk=sqrt(bsq)
    end if
    capf=1.+2.*p1k(k,j)*(2.*abf(i)*bk**2+bbf(i))+2.*p2k(k,j)*
c (2.*abp(i)*bk**2+bbp(i))
    deli1(k)=1./bk
    deli2(k)=1./(bk**3*capf)
    deli3(k)=deli2(k)/(capf*bk**2)
    deli4(k)=p1k(k,j)/(capf*bk)**3
    deli5(k)=p2k(k,j)/(capf*bk)**3
32 continue
call simps(deli1,dcapi1,kin,dps)

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      call simps(deli2,dcapi2,kin,dps)
      call simps(deli3,dcapi3,kin,dps)
c     call simps(deli4,dcapi4,kin,dps)
      call simps(deli5,dcapi5,kin,dps)
      capi1=dcapi1+sum1
      capi2=dcapi2+sum2
      capi3=dcapi3+sum3
      capi4=dcapi4+sum4
      capi5=dcapi5+sum5
      r(i,j)=sqrt(2.*capi1)
      vv1=(bvac(i)*dbvdz(i))**2
      vv2=bvac(i)*d2bvdz2(i)+dbvdz(i)**2
      rzz(i,j)=-vv1*capi2**2/r(i,j)**3+3.*vv1*capi3/r(i,j)
c     -vv2*capi2/r(i,j)+8.*bvac(i)**2*(abf(i)*capi4+abp(i)*
c     capi5)/r(i,j)
      sum1=+capi1
      sum2=capi2
      sum3=capi3
      sum4=capi4
      sum5=capi5
40    continue
c     ....set r(i,1)=r(i,2)
      do 41 i=1,ix
41    r(i,1)=r(i,2)
c     calculate diagnostic quantities flute1, flute2 and flute3
      do 60 j=2,jx-1
        is=2
        if(mod(ix-2,2).eq.0) is=3
        do 61 i=is,ix-1
          ia=i-is+1
          eterm=dp2dpsi(j)*(abp(i)*b(i,j)**4+bbp(i)*b(i,j)**2+cbp(i))
c         +p2(j)*(4*abp(i)*b(i,j)**3+2*bbp(i)*b(i,j))*dbdpsi(i,j)
          rterm=0.
          if(dp2dpsi(j).ne.0.)
c         c rterm=dp2dpsi(j)*(ebp(i)*b(i,j)**4+fbp(i)*b(i,j)**2+gbp(i))
c         c /(2.*p2(j)**.5)
c         c +p2(j)**.5*(4*ebp(i)*b(i,j)**3+2*fbp(i)*b(i,j))*dbdpsi(i,j)
c         c rterm=exrho/psi(jx)*rho(i,j)
          droterm(ia)=rterm/b(i,j)**2/uuz(i)
          ringj=dp1dpsi(j)*(abf(i)*b(i,j)**4+bbf(i)*b(i,j)**2+cbf(i))
c         +p1(j)*(4*abf(i)*b(i,j)**3+2*bbf(i)*b(i,j))*dbdpsi(i,j)
          dflute1(ia)=rzz(i,j)*qubv(i,j)/(r(i,j)*b(i,j)**2)/uuz(i)
c         +eterm*ringj/b(i,j)**3/uuz(i)
          dflute2(ia)=yyy(i,j)/(r(i,j)*b(i,j)
c         )**2*b(i,j)/uuz(i)
          dflute3(ia)=rho(i,j)/(r(i,j)*b(i,j)**2*b(i,j))/uuz(i)
          dflute4(ia)=xxx(i,j)/(r(i,j)*b(i,j)**2*b(i,j))/uuz(i)
61    continue
      call simps(dflute1,ans1,ia,du)
      call simps(dflute2,ans2,ia,du)
      call simps(dflute3,ans3,ia,du)
      call simps(dflute4,ans4,ia,du)
      call simps(droterm,ans5,ia,du)
      flute1(j)=ans1*(ia-1)
      flute2(j)=(ans2*(mm**2-1)+ans1)*(ia-1)
      flute3(j)=-ans1/ans3
      rhoave(j)=ans3*(ia-1)
      xxxave(j)=ans4*(ia-1)
      yyyave(j)=ans2*(ia-1)
      droave(j)=ans5*(ia-1)

```



```

60  continue

      grow=0.
      growmax=0.
      do 62 j=2,jx-1
        if(flute3(j).gt.0)grow=grow+flute3(j)
        growmax=amax1(growmax,flute3(j))
62  continue
      write(59,102)grow,growmax
102  format('grow=',e14.6,3x,'growmax=',e14.6)
c..... local growth rate for high mm, (wkb approx. )
      do 70 j=2,jx-1
        omegwkb=mm*.5*xxxave(j)/rhoave(j)
        trad=.25*mm**2*(xxxave(j)**2+4.*rhoave(j)*yyyave(j))
c -rhoave(j)**2*flute3(j)-droave(j)*yyyave(j)
        if(trad.le.0.)then
          gamwkb(j)=sqrt(-trad)/rhoave(j)
          omeg1wkb(j)=omegwkb
          omeg2wkb(j)=0.
        else
          omeg1wkb(j)=omegwkb +sqrt(trad)/rhoave(j)
          omeg2wkb(j)=omegwkb-sqrt(trad)/rhoave(j)
        end if
70  continue
      return
      end
      subroutine fltoll
c.....calculates the f1 to f11 functions needed to generate the a and
c..... b matrices . uses the equilibrium quantities r, rho, b, etc.
c..... insert cliché storage here

      use param
      use fstor
      use matrix
      use const

      m2=mm**2
      du2=du**2
      do 10 i=1,ix
        do 10 j=2,jx
          r2=r(i,j)**2
          uz=uuz(i)
          bb=b(i,j)
          vp=vpsi(j)
          r4=r2**2
          f1(i,j)=rho(i,j)*bb*r4
          f2t=(1.-m2)*rho(i,j)/bb+r2*vp*(rho(i,j+1)-rho(i,j-1))/(2.*dv)
          f2(i,j)=f2t/vp
          f3(i,j)=mm*xxx(i,j)*r4*bb
          f4(i,j)=(1.-m2)*mm*xxx(i,j)/bb
          f5(i,j)=-m2*yyy(i,j)*r4*bb
          f7(i,j)=(1.-m2)*(-m2)*yyy(i,j)/(bb*vp)
          g4(i,j)=qub(i,j)*r(i,j)**2
          g3(i,j)=r(i,j)*b(i,j)
          g2(i,j)=qub(i,j)/(r(i,j)*b(i,j))**2
          dppdpsi=dp2dpsi(j)*(abp(i)*b(i,j)**4+bbp(i)*b(i,j)**2+cbp(i))
c +p2(j)*(4*abp(i)*b(i,j)**3+2*bbp(i)*b(i,j))*dbdpsi(i,j)
          dpedpsi=dp1dpsi(j)*(abf(i)*b(i,j)**4+bbf(i)*b(i,j)**2+cbf(i))
c +p1(j)*(4*abf(i)*b(i,j)**3+2*bbf(i)*b(i,j))*dbdpsi(i,j)
          ering(i,j)=dppdpsi*dpedpsi

```

```

      g1(i,j)=+(mm*uzz(i))**2*r(i,j)*(rzz(i,j)
c*qubv(i,j)+dppdpsi*dpedpsi*r(i,j)/b(i,j))
      g1(i,j)=g1(i,j)*swg1
      g2(i,j)=g2(i,j)*swg2
      g3(i,j)=g3(i,j)*swg3
      g4(i,j)=g4(i,j)*swg4
c..... special g1 to test b.c. on flute test case
c      g1(i,j)=-(mm*uzz(i))**2*r(i,j)*r(i,j)/lb**2
c      c*qv(i,j)

```

```

10  continue
c..... fill in edge values
do 20 i=1,ix
  f1(i,1)=-f1(i,2)
  f2(i,1)=f2(i,2)
  f3(i,1)=-f3(i,2)
  f4(i,1)=f4(i,2)
  f5(i,1)=-f5(i,2)
  f7(i,1)=f7(i,2)
  g4(i,1)=-g4(i,2)
20  continue
  return
end

```

subroutine grid

c..... relates physical grid z,psi to computational grid u,v (equally
c..... spaced 1. uses input fpsl, fv, fz, fu and azm, apsim .

```

c.....insert cliche storage here
  use param
  use const

  xv=alog(fpsl)/alog(fv)
  xu=alog(fz)/alog(fu)
  zzp=0.
  psip=0.
5  do 5 i=1,ix
    u(i)=u0+du*(i-1.5)
    u(1)=-u(1)
    azm=u(ix)**(1.-xu)
    do 10 i=1,ix
      uuz(i)=u(i)**(1.-xu)/(xu*azm)
      z(i)=azm*u(i)**xu

      uuzh(i)=(u(i)+.5*du)**(1.-xu)/(xu*azm)
      dz(i)=z(i)-zzp
      zzp=z(i)
10  continue
    zedge=azm*.5*(u(ix)**xu+u(ix-1)**xu)
    do 15 j=1,jx
      v(j)=v0+(j-1.5)*dv
15  continue
    v(1)=-v(1)
    apsim=v(jx)**(1.-xv)
    do 20 j=1,jx
      vpsi(j)=v(j)**(1.-xv)/(apsim*xv)
      vpsih(j)=(v(j)+.5*dv)**(1.-xv)/(apsim*xv)

```

```

        psi(j)=apsim*v(j)**xv
        dpsi(j)=psi(j)-psip
        psip=psi(j)
20    continue
c..... calculates v at plasma edge
        vw=(psiw/apsim)**(1./xv)
        dvin=(vw-v(jx-1))/dv
        dvout=(v(jx)-vw)/dv
        return
        end
        subroutine initial

c.....set up initial displacement vectors, xro and xio
c..... test case 1, cos(kz) in z, fiat in psi
c..... set up 1/4/82 by r, freis

c.....insert cliche storage here
        use param
        use fstor
        use matrix
        use const

        data pi/3.1415926/

        rbf=1.
        do 10 j=2,jx-1
            r1=ranf(b1)
            r2=ranf(b1)
            r3=ranf(b1)
            r4=ranf(b1)
            do 10 i=2,ix-1
                if(psih.gt.psi(j))then
                    rbf=1./(r(i,j)*b(i,j))
                    kzsp=0
                else
                    r1=ranf(b1)
                    r2=ranf(b1)
                    r3=ranf(b1)
                    r4=ranf(b1)
                    kzsp=1
                endif
5            continue
            k1=i-1+(j-2)*(ix-2)
            k2=j-1+(ix-2)*(i-2)
            k=.5*(1+isw)*k1+.5*(1-isw)*k2
            cosx=cos(.5*pi*(z(i)*kzsp)/zedge)
            xro(k)=ex0*cosx*rbf*(r1+r2-1.)+ex1*cos(.5*pi*(z(i))/zedge)
            xio(k)=ex0*cosx*rbf*(r3+r4-1.)+ex1*cos(.5*pi*(z(i))/zedge)
c            xio(k)=cos(theta0)*xro(k)
10        continue
        do 20 j=2,jx-1
            r1=ranf(b1)
            r2=ranf(b1)
            r3=ranf(b1)
            r4=ranf(b1)
            do 20 i=2,ix-1
                if(psih.gt.psi(j))then
                    rbf=1./(r(i,j)*b(i,j))
                    kzsp=0
                else

```

```

        r1=ranf(b1)
        r2=ranf(b1)
        r3=ranf(b1)
        r4=ranf(b1)
        kzsp=1
        endif
15  continue
        k1=i-1+(j-2)*(ix-2)
        k2=j-1+(jx-2)*(i-2)
        k=.5*(1+isw)*k1+.5*(1-isw)*k2
        cosx=cos(.5*pi*(z(i)*kzsp)/zedge)
        xroo(k)=ex0*cosx*rbf*(r1+r2-1.)+ex1*cos(.5*pi*(z(i))/zedge)
        xloo(k)=ex0*cosx*rbf*(r3+r4-1.)+ex1*cos(.5*pi*(z(i))/zedge)
c      xloo(k)=cos(theta0)*xroo(k)
20  continue
        return
        end

        subroutine input
c.....insert storage cliches here
        use param
        use const
        use fstor

c..... boundary conditions are set as follows:
c.....          at z=z0 (i=1), f11=-1, implies x=0,
c.....                      f11=1, implies slope=0,
c.....          at z=zmax (i=ix), fizx=-1, implies x=0
c.....                      fizx=1, implies slope=0,
c.....          at psi=psi0 (j=1), fj1=-1, implies x=0,
c.....                      fj1=1, implies slope=0,
c.....          at psi=psimax (j=jx), fjrx=-1, implies x=0,
c.....                      fjrx=1, implies slope=0,
        data mm/4/, bias/.5/, lmax/2/, nmax/5/, dv/1./, du/1./,
c      ,ndiag/100/,f11/1./,fizx/1./,fj1/1./,fjrx/1./,flr/1./
c      ,sf6/1./, sf8/1./, kplotm/0/, kzs/1/, swg1/1./, swg2/1./
c      , swg3/1./, swg4/1./, nen/1/

c.....forced data loaded for testing fourier analyses and zed file
c.....,maker
        data jfour/1/,nfourp/1/,nfourmax/5/
        namelist/now1/aname,mm, bias, lmax, nmax,ndiag,nen
c      ,f11,fizx,fj1,fjrx,ex0,ex1,fpsi,fu,fv,fz
c      ,kplotm,kzs,jfour,nfourp,nfourmax,sf6,sf8,swg1,swg2,swg3,swg4

        call ddi(now1,2,3,1)
        if(nold.ne.1)call ddo(now1,100,0,1)
        jx=jrx
        kxx=kxp
        ix=izx
        return
        end
        subroutine inptemp

c.....temporary input to test three region equilib.
        use param
        use const
        real klbsq
        namelist/btemp/bmx1,bmx2,bmx3,bmn1,bmn2,betpas1,ppas2
c      ,ppas3,betrap,z1c,z2c,z3c,z1min,z2min,as,als

```