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

cliche param
parameter (lzx=60, jrx=40, lzx2=lzx-2 )
parameter ( lhw=2*lzx-1 )
parameter (kxp=(lzx-2)*(jrx-2), lhw=lzx-1 )
parameter (npl=3000, nps=100 )
parameter (lxpf=4*(lzx-2),nfourxf=150)
parameter (lxpg=2*lzx2)
parameter (neng=500)
parameter (ksim=51)
endcliche
cliche fstor
use param
common/fun/f1(lzx,jrx),f2(lzx,jrx),f3(lzx,jrx),f4(lzx,jrx)
c ,f5(lzx,jrx),f7(lzx,jrx),
c g1(lzx,jrx),g2(lzx,jrx),g3(lzx,jrx),g4(lzx,jrx)
c ,swg1,swg2,swg3,swg4

common/equil/b(lzx,jrx),rho(lzx,jrx),qub(lzx,jrx),r(lzx,jrx)
c ,phi(lzx,jrx),yyy(lzx,jrx),xxx(lzx,jrx),qv(lzx,jrx)
common/pertur/x100(kxp),x10(kxp),x101(kxp)
c ,xron(kxp),xro(kxp),xrol(kxp)
endcliche

cliche matrix
use param
common/couff/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,lzxp,kxx,nmax,lmax,lsw,lhw
c ,fac1,fac2,bias,du,dv,dt,ndiag,ex0,b0,rho0,ex1,f1,fizx,fj1
c ,fjrx,kplot,npn,fpst,fz,fu,fv,azm,apsim,u0,v0,amass,nengx
c ,fourpi,omegst,omegr,omegexb,fir,sf6,sf8,kplotm,kzs,zedge
c ,cpuo,clo,syso,valfk,xu,xv,n,pl,vw,psiw,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,lp1
c ,bcen,prin,epspp,phicen,phiplg,kin,xpot,ypot,wpot,pfudge,rpx
c ,phice,phipl,betslsh,betcnt,z0,z1,z2,z3,z4,psi1,psi2,psi0
c ,betcne,betslse,pslosho,pcentee,bmax,alsi,bm1,psi1,psi1,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 ,psi0rel,psi0e,psime,p2wide,wp2e,p2floor,p1floor,p2flag
c ,fring,long,no3d,no1d,dphi,dip,psistr,psisip,psihrel,psih
c ,dpsihsrel,dpsihs,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,bmn1,bmn2,ppas1,ppas2,ppas3,p1trap
c ,z1c,z2c,z3c,z1min,z2min,as(3),als(3),zs(3),bs(3),dpas1,d1trap
c ,betrap,betp,s1,bvx2,bvx3,ncoil,z2ct
common/mesh/psi(jrx),z(lzx),u(lzx),v(jrx),dpsi(jrx),dz(lzx)

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c ,vpsi(jrx),uuz(izx),vpsi1(jrx),uuzh(izx)
common/graf/ xrtimel(npli),xrspz(izx,nps),xrspst(jrx,2*nps)
c ,time(npli),xflute(izx,nps),enpot(neng),tenergy(neng)
x ,enkin(neng),timengy(neng),tenrel(neng),enbend(neng)
c ,encurve(neng),enflr(neng)
common/curvec/cr,lb,rw,beta0,delrho,stable,en0,coo,r0,
c ,echarg,omeg1,omeg2,en1,besarg,z0l,dtrel,p0,omego
c ,omana1,omana2,groana,theta0
common/tmcon/h12(izx),hzt0(izx),h3d(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),cq(izx),ebp(izx)
c ,fbp(izx),gbp(izx),betring,hp0(jrx),hpm(jrx),hpmel(jrx),hf1r(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),dp1dps1(jrx)
c ,p1(jrx),p1k(ksim,jrx),hpk12(ksim),hpk0(ksim),del1(ksim)
c ,del12(ksim),del13(ksim),del14(ksim),rzz(izx,jrx),dbdps1(izx,jrx)
c ,phi1(izx),phi2(izx),pperp(izx,jrx),ppar(izx,jrx),dflute3(izx)
c ,qubv(izx,jrx),p2(jrx),dp2dps1(jrx),dflute1(izx),dflute2(izx)
c ,flute1(jrx),flute2(jrx),flute3(jrx),p2k(ksim,jrx),del15(ksim)
c ,pperps(izx,jrx),errprp(izx,jrx),errprl(izx2,jrx-1)
c ,pperpel(izx,jrx),eps1(izx,jrx),omeg1wkb(jrx),omeg2wkb(jrx)
c ,gamwkb(jrx),dflute4(izx),rhoavel(jrx),xxxavel(jrx),yyyavel(jrx)
c ,p2t(jrx),dp2dps1(jrx),p3(jrx),dp3dps1(jrx),hpkm(ksim)
c ,hpkme(ksim),droavel(jrx),droterm(izx),ering(izx,jrx)
common/forced/nfour,nfourx,nfourmax,nfourp,jfour,ixp,locv

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

return
end

c...., the main routine

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=u*xu**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 ) .

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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 cliche 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.....floratm, tandem mirror equilibrium

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

c.....,flortex, tandem mirror equilib. 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.....(e-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..... parallel pressure check, and e-psi (= -dphi/dpsi) . Phi2 modified
c..... to " 1.-(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=a1+b1*(
c..... psi/psime)+c1*(psi/psime)**2, and pe2=.5*(1-tanh((psi-ps10e)/p2ewide))
c

c..... flortm6, modified flortm5 as follows; for p2(psille, 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, no1d=1, prevents graflib plots, no3d=1
c..... prevents tv80lib 3d plots.
c

c..... flortm8 (18 for larger psi grid) modified flortm6 as follows:
c..... the analytic calculation of gammwb 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..... phicent+dphi in plug, dphi a new input variable). Also b calculated
c..... with expansion in low beta regions.

c..... flortn8, like flort18 with first order energy check added .

c..... firm1 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..... firm2, 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..... firm3, like firm2 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..... firm4, cold plasma halo modeled by changing zmax boundary conditions
c..... for psi > psih (psihrel an input parameter)

c..... firm6, like firm4 and firm5, (mixed boundary condition at zmax,
c..... higher order b, c.) with special vyy to force "rigid mode" in psi.

c..... firrot, modified firm6 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/npte
data tally/1/

c.....call link call here
call link('unit59=terminal,unit2=(inflm4,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 locally=ii
locally=14
if(locally,eq,0)go to 299
call timer(locally,'ztally00',tallyb,2000b,floratim,1)
299 itally=-1
endif

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lsw=1
if(lzx.gt.jrx) lsw=-1
jtbw=.5*(1+lsw)*ltbw+.5*(1-lsw)*(2*jrx-1)
lhbw=.5*(1+lsw)*lbw+.5*(1-lsw)*(jrx-1)
nn=jtbw*kxp
nn1=jtbw+kxp
call memory(ww,nn-1)
call memory(ww1(nn),nn1)
namelist/noplot/nold,no3d
call ddi(noplot,2,0,1)
if(nold.ne.1)call pstart(dev,4rplot,1,'box u21$',1)
c
npote=1
if(nold.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
c call constant
c call tmcon2
c call equiltm
c call equilrot
c call equilcur
c call fito11
c call amat
c call comat(ww,jtbw)
c 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(xiio(1),xiol(1),kxx)
    call banfac(kxp,lhbw,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 zmoveurd(ww1(nn),rhs1,kxx)
call bansol(kxp,lhbw,ww,1,-(kxp-1),ww1(nn))
do 10 j=2,jx-1
kp=1+izxp*(j-2)
call zmoveurd(xrol,ww1(nn),kxx)
10 continue
call zmoveurd(ww1(nn),rhs2,kxx)
call bansol(kxp,lhbw,ww,1,-(kxp-1),ww1(nn))
do 20 j=2,jx-1
kp=1+kxp*(j-2)
call zmoveurd(xiol,ww1(nn),kxx)
20 continue
fac1=-.5/dt
90 fac2=.5/dt

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call zmovewrd(x100,x10,kxx)
call zmovewrd(x10,x10,kxx)
call zmovewrd(xr00,xr0,kxx)
call zmovewrd(xr0,xr0,kxx)

c..... time array
xrtim(n)*xr0(kplot)
if(mod(n,ndiag).eq.0)call diagno
if(mod(n,nfourp).eq.0)call fourier
if((mod(n,nent).eq.0).or.(lnt.ne.0).and.(lce.le.neng))call energy
100 continue
call cldsk(lacv,0)
call timeused(lcp,lo,isy)
cpuo=lcp*tim
clo=lo*tim
syso=sy*tim
if(nold.ne.1)
c call picshar
call close(100)
if(no3d.eq.1)go to 300
call keep80(1,3)
call fr80id
call threed
call plot
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 fito11 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=izx
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))*.

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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*uuzh(i)
g4iphjmh=(g4(i+1,j-1)+g4(i,j)+g4(i+1,j)+g4(i,j-1))*25*uuzh(i)
g4imhjph=(g4(i-1,j+1)+g4(i,j)+g4(i-1,j)+g4(i,j+1))*25*uuzh(i)
g4imhjmh=(g4(i-1,j-1)+g4(i,j)+g4(i-1,j)+g4(i,j-1))*25*uuzh(i)
g2iphj=(g2(i+1,j)+g2(i,j))*5*uuzh(i)
g2imhj=(g2(i-1,j)+g2(i,j))*5*uuzh(i-1)
g3iphj=(g3(i+1,j)+g3(i,j))*5*uuzh(i)
g3imhj=(g3(i-1,j)+g3(i,j))*5*uuzh(i-1)

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

if(j.gt.2)go to 60
f1ijmh=0,
f6ijmh=0,
60 continue
uzbar=-uuz(i)*r(i,j)/(du2*dv2)
a1(k,1)=-gam1*b1m1jmh*g4imhjmh*bijmh*uzbar*vpsihi(j-1)
c *r(i-1,j-1)
a2(k,1)=-gam2*b1m1jmh*g4imhjmh*bijmh*uzbar*vpsihi(j-1)
c *r(i-1,j-1)
a3(k,1)=-gam3*b1m1jmh*g4imhjmh*bijmh*uzbar*vpsihi(j-1)
c *r(i-1,j-1)

a1(k,2)=-f1ijmh/((dt*dv)**2)+gam1*(f5ijmh/dv2+bijmh**2*uzbar
c *vpsihi(j-1)*r(i,j-1)*(g4imhjmh+g4iphjmh))
a2(k,2)=-f1ijmh/((dt*dv)**2)+gam2*(f5ijmh/dv2+bijmh**2*uzbar
c *vpsihi(j-1)*r(i,j-1)*(g4imhjmh+g4iphjmh))
a3(k,2)=-f1ijmh/((dt*dv)**2)+gam3*(f5ijmh/dv2+bijmh**2*uzbar
c *vpsihi(j-1)*r(i,j-1)*(g4imhjmh+g4iphjmh))

a1(k,3)=-gam1*b1p1jmh*g4iphjmh*bijmh*uzbar*vpsihi(j-1)*r(i+1,j-1)
a2(k,3)=-gam2*b1p1jmh*g4iphjmh*bijmh*uzbar*vpsihi(j-1)*r(i+1,j-1)
a3(k,3)=-gam3*b1p1jmh*g4iphjmh*bijmh*uzbar*vpsihi(j-1)*r(i+1,j-1)
a1(k,4)=gam1*((b1m1jmh*g4imhjmh*vpsihi(j-1)*bijmh+b1m1jph*g4imhjph
c *vpsihi(j)*bijph)*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*vpsihi(j-1)*bijmh+b1m1jph*g4imhjph
c *vpsihi(j)*bijph)*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*vpsihi(j-1)*bijmh+b1m1jph*g4imhjph
c *vpsihi(j)*bijph)*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)+(-bijmh**2*(g4imhjmh+g4iphjmh)*vpsihi(j-1)
c -bijph**2*(g4imhjph+g4iphjph)*vpsihi(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)+(-bijmh**2*(g4imhjmh+g4iphjmh)*vpsihi(j-1)
c -bijph**2*(g4imhjph+g4iphjph)*vpsihi(j))*r(i,j)*uzbar-
c mm**2*b(i,j)*uzbar*(g2imhj+g2iphj)*g3(i,j)*dv2/vpsi(j))

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a3(k,5)=((f1;jph+f1;jmh)/dv2-f2(i,j))/dt2+gam3*(-(f5;jph+f5;jmh
c 1/dv2+f7(i,j)+g1(i,j)+(-b1;jmh**2*(g4imh;jmh+g4iph;jmh)*vpsi(j-1)
c -b1;jph**2*(g4imh;jph+g4iph;jph)*vpsi(j)*r(i,j)*uzbar-
c mm**2*b(i,j)*uzbar*(g2imh;j+g2iph;j)*g3(i,j)*dv2/vpsi(j))

a1(k,6)=gam1*((b1p1;jmh*g4iph;jmh*b1;jmh*vpsi(j-1)+b1p1;jph*g4iph;jph
c *b1;jph*vpsi(j))*uzbar*r(i+1,j)+mm**2*b(i,j)*uzbar*
c g2iph;j*g3(i+1,j)*dv2/vpsi(j))
a2(k,6)=gam2*((b1p1;jmh*g4iph;jmh*b1;jmh*vpsi(j-1)+b1p1;jph*g4iph;jph
c *b1;jph*vpsi(j))*uzbar*r(i+1,j)+mm**2*b(i,j)*uzbar*
c g2iph;j*g3(i+1,j)*dv2/vpsi(j))
a3(k,6)=gam3*((b1p1;jmh*g4iph;jmh*b1;jmh*vpsi(j-1)+b1p1;jph*g4iph;jph
c *b1;jph*vpsi(j))*uzbar*r(i+1,j)+mm**2*b(i,j)*uzbar*
c g2iph;j*g3(i+1,j)*dv2/vpsi(j))

a1(k,7)=gam1*(-b1m1;jph*g4imh;jph*b1;jph*vpsi(j)*uzbar*r(i-1,j+1))
a2(k,7)=gam2*(-b1m1;jph*g4imh;jph*b1;jph*vpsi(j))*uzbar*r(i-1,j+1)
a3(k,7)=gam3*(-b1m1;jph*g4imh;jph*b1;jph*vpsi(j))*uzbar*r(i-1,j+1)
a1(k,8)=-f11;jph/(dt2*dv2)+gam1*(f5;jph/dv2+(b1;jph**2*(g4imh;jph
c +g4iph;jph)*vpsi(j)*r(i,j+1)*uzbar))
a2(k,8)=-f11;jph/(dt2*dv2)+gam2*(f5;jph/dv2+(b1;jph**2*(g4imh;jph
c +g4iph;jph)*vpsi(j)*r(i,j+1)*uzbar))
a3(k,8)=-f11;jph/(dt2*dv2)+gam3*(f5;jph/dv2+(b1;jph**2*(g4imh;jph
c +g4iph;jph)*vpsi(j)*r(i,j+1)*uzbar))
a1(k,9)=gam1*(-b1p1;jph*g4iph;jph*b1;jph*vpsi(j)*uzbar*r(i+1,j+1))
a2(k,9)=gam2*(-b1p1;jph*g4iph;jph*b1;jph*vpsi(j))*uzbar*r(i+1,j+1)
a3(k,9)=gam3*(-b1p1;jph*g4iph;jph*b1;jph*vpsi(j))*uzbar*r(i+1,j+1)

c.....,b1 array for rhs
    f3ijmh=(f3(i,j)+f3(i,j-1))*,.5*vpsi(j-1)
    f3ijph=(f3(i,j)+f3(i,j+1))*,.5*vpsi(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.sfl1.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.sflzx.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(sfj1.eq.1.)sfj1=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)+sfj1*a1(k,m-1)
a2(k,m)=a2(k,m)+sfj1*a2(k,m-1)
a3(k,m)=a3(k,m)+sfj1*a3(k,m-1)
11 continue
13 continue
i=ix-1
do 12 j=2,jx-1
bc(j)=0,
c if(psi(j).lt.(psi-h-dpsi*h))bc(j)=1.
c if(psi(j).ge.(psi+h-dpsi*h).and.psi(j).le.(psi+h+dpsi*h))
bc(j)=.5*(tanh((-psi(j)+psi*h)/dpsi*h)+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)-1,1))
down=(bc(j)*r(ix,j)*b(ix,j)*uuzh(ix-1)/du+rbt*(-bc(j)+1,1))
sfizx=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,

```

```

      a2(k,m-3)=0,
      a3(k,m-3)=0,
44    continue
      b1(k,1)=0,
45    continue
      j=jx-1
      do 48 i=2,jx-1
      k1=i-1+(j-2)*(jx-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
      bcorrp=0,
      bcorpp=0,
      tanhyp=tanh((z(i)-zcc)/znorm)
      bcorrp=-bcc*.5*(1.-tanhyp**2)/znorm
      bcorpp=+bcc*tanhyp*(1.-tanhyp**2)/znorm**2
      bcorr=bcc*.5*(1.-tanhyp)
18    bv(i)=bv(i)+bcorr
      bvp(i)=bvp(i)+bcorrp
      bvpp(i)=bvpp(i)+bcorpp
      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)
50 continue
```

```
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),1,14,amin)
z4=z(14)
if(ncoil.eq.2) return
b5=aminaf(bv,is(2),is(3),1,15,amin)
z5=z(15)

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
  ip1=m+((m-1)/3)*(ihbw-4)
  ip2=1+mod(m-1,3)*(ihbw-1)+(m-1)/3
  ip=.5*t1+isw)*ip1+.5*(1-isw)*ip2
  abar(k,ip)=a1(k,m)
10 continue
return
end

subroutine constant

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

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

```

```

gdm2=.25*(1-bias)
ip=.5*(ix-2)
jp=.5*(jx-2)
kp1=ip-1+(jp-2)*(ix-2)
kp2=jp-1+(jx-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 echarge/4.8e-10/, en0/1.00e+12/, b0/1.e4/, amass/3.34e-24/
  c , coe/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,stabl,en0,echarge,lb,rw,xm
  c , z0l,dtrel,theta0
  call ddi(curve,2,3,1)
  if(nold.ne.1)  call ddo(curve,100,0,1)
  zmax=z0l*lb
  p0=beta0*b0**2*.5
  psimax=rw**2*b0*.5
  omeg0sq=b0**2/(en0*amass*b0**2)
  omeg0=sqrt(omeg0sq)
  ag1=1.-2.*delrho/xm**2
  klbsq=0.
  if(kzs.ne.0)klbsq=(pi/(2.*z0l))**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 trapezoidal quadrature) over the volume,
c...   every non'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 rxl(jrx),rxf(jrx),drxr(lzx,jrx),drxl(lzx,jrx),quad(12)
c ,quad1(4),psum1(jrx,12),psumb(jrx,4),dum1(lzx),dum2(lzx)
c ,xrg(lzx),xig(lzx),xrgo(lzx),xigo(lzx),drxro(lzx,jrx)
c ,drxlo(lzx,jrx),drxz2(lzx),dxiz2(lzx)

ltt=ltt+1
if((ltt,eq,2),or,(nen,eq,1))then
lmove=jx*lx
call zmovewd(drxro,drxr,lmove)
call zmovewd(drxlo,drxl,lmove)
tquad0=tquad
tquad200=tquad20
tquad2o=tquad2
tquad300=tquad30
tquad3o=tquad3
tquad400=tquad40
tquad4o=tquad4
end if
do 9 i=2,ix-1
do 5 j=2,jx-1
k1=i-1+(j-2)*(lx-2)
k2=j-1+(jx-2)*(i-2)
k=.5*((1+isw)*k1+(1-isw)*k2)
rxf(j)=r(i,j)*xro(k)
rxl(j)=r(i,j)*xlo(k)
5    continue
call ddpsi(rxf,drxr,1)
call ddpsi(rxl,drxl,1)
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)*(lx-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)=xloo(k)
end if

```

```

      xrg(1)=xrg0(k)
10      xig(1)=xi0(k)
      do 211 i=2,1x-1
211  dum2(i)=r(i,j)*b(i,j)*xrg(i)
           call ddz(dum2,dxrz2)
      do 212 i=2,1x-1
212  dum2(i)=r(i,j)*b(i,j)*xig(i)
           call ddz(dum2,dxit2)
      do 11 i=2,1x-1
11    dum1(i)=(1./uuz(i))*qub(i,j)*dxrz2(i)/((r(i,j)*b(i,j))**2)
           psumal(j,1)=psum(dum1,1x)/vpsi(j)
      do 12 i=2,1x-1
12    dum1(i)=(1./uuz(i))*qub(i,j)*dxitz2(i)/((r(i,j)*b(i,j))**2)
           psumal(j,2)=psum(dum1,1x)/vpsi(j)
      do 13 i=2,1x-1
13    dum1(i)=b(i,j)*drxit(i,j)
           call ddz(dum1,dum2)
      do 14 i=2,1x-1
14    dum1(i)=(1./uuz(i))*qub(i,j)*r(i,j)**2*dum2(i)
           psumal(j,3)=psum(dum1,1x)/vpsi(j)/mm**2
      do 15 i=2,1x-1
15    dum1(i)=b(i,j)*drxr(i,j)
           call ddz(dum1,dum2)
      do 16 i=2,1x-1
16    dum1(i)=(1./uuz(i))*qub(i,j)*r(i,j)**2*dum2(i)
           psumal(j,4)=psum(dum1,1x)/vpsi(j)/mm**2
      do 17 i=2,1x-1
17    dum1(i)=(1./uuz(i))*xrg(i)**2*r(i,j)*(qubv(i,j)*rzzi(j)
c   +r(i,j)*eringl(i,j)/b(i,j))
           psumal(j,5)=psum(dum1,1x)/vpsi(j)
      do 18 i=2,1x-1
18    dum1(i)=(1./uuz(i))*xig(i)**2*r(i,j)*(qubv(i,j)*rzzi(j)
c   +r(i,j)*eringl(i,j)/b(i,j))
           psumal(j,6)=psum(dum1,1x)/vpsi(j)
      do 19 i=2,1x-1
19    dum1(i)=(1./uuz(i))*xrg(i)**2*yyy(i,j)/b(i,j)
           psumal(j,7)=mm**2*psum(dum1,1x)/vpsi(j)
      do 20 i=2,1x-1
20    dum1(i)=(1./uuz(i))*xig(i)**2*yyy(i,j)/b(i,j)
           psumal(j,8)=mm**2*psum(dum1,1x)/vpsi(j)
      do 21 i=2,1x-1
21    dum1(i)=(1./uuz(i))*yyy(i,j)*xrg(i)*r(i,j)*drxr(i,j)
           psumal(j,9)=-2.*psum(dum1,1x)/vpsi(j)
      do 22 i=2,1x-1
22    dum1(i)=(1./uuz(i))*yyy(i,j)*r(i,j)**2*b(i,j)*drxr(i,j)**2
           psumal(j,10)=psum(dum1,1x)/vpsi(j)
      do 23 i=2,1x-1
23    dum1(i)=-(1./uuz(i))*yyy(i,j)*xig(i)*r(i,j)*drxit(i,j)
           psumal(j,11)=2*psum(dum1,1x)/vpsi(j)
      do 24 i=2,1x-1
24    dum1(i)=(1./uuz(i))*yyy(i,j)*r(i,j)**2*b(i,j)*drxit(i,j)**2
           psumal(j,12)=psum(dum1,1x)/vpsi(j)
           if(ltt.eq.2).or.(nen.eq.1) then
      do 30 i=2,1x-1
30    dum1(i)=(1./uuz(i))*rho(i,j)*((xrg(i)-xrg0(i))/dt)**2/b(i,j)
           psumbl(j,1)=psum(dum1,1x)/vpsi(j)
      do 31 i=2,1x-1
31    dum1(i)=(1./uuz(i))*rho(i,j)*((xig(i)-xigo(i))/dt)**2/b(i,j)
           psumbl(j,2)=psum(dum1,1x)/vpsi(j)
      do 32 i=2,1x-1

```

```

32      dum1(1)=(1./uuz(1))*(r(1,j)*(drxr(1,j)-drxro(1,j))/dt)**2*b(1,j)
c *rho(1,j)
      psumb(j,3)=psum(dum1,1x)/vpsi(j)/mm**2
      do 33 i=2,ix-1
33      dum1(1)=(1./uuz(1))*(r(1,j)*(drxi(1,j)-drxio(1,j))/dt)**2*b(1,j)
c *rho(1,j)
      psumb(j,4)=psum(dum1,1x)/vpsi(j)/mm**2
      end if
100    continue
      do 120 l=1,12
120    quad(l)=psum(psum(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(ltt.eq.2).or.(nen.eq.1) then
      tquad1=0.
      lee=lee+1
      do 130 l=1,4
130    quad1(l)=psum(psumb(1,l),jx)
      tquad1=tquad1+quad1(l)
      enpot(lee)=(3.*tquad+6.*tquad0-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
      enflr(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-1))*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(1x,1)

        do 10 j=4,jx-1
10      f2(1,j-1)=(f1(j)-f1(j-2))/(2.*dv)*vpsi(j-1)
            f2(1,2)=(2.*f1(2)*vpsi(1)+(f1(3)-f1(2))*vpsi(2))/(dv*2)
            f2(1,jx-1)=(-(1.-fjrx)*f1(jx-1)*vpsi(jx-1)+(f1(jx-1)-
c f1(jx-2))*vpsi(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
        uzz=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 cliches 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(j)
10      continue
        return
        end

```

subroutine equilrot

c..... sets up equilibrium for rigid rator, 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(fourni)

omegr=ratrod*omegst*(1.-enbar/en0)

foursq=sqrt(fourpi)

do 5 i=1,ix

uzz=uuz()

do 5 j=1,jx

fac=exp(psi(i)/psi0)/sqrt(beta0)

b(i,j)=b0*sqrt(fac**2-1.)/(fac*foursq)

rholl(i,j)=en0*amass/(beta0*fac**2)+enbar*amass

beta=1/fac**2

arg1=fac+sqrt(fac**2-1.)

acosh=log(arg1)

r(i,j)=r0*sqrt(-cr+acosh)

qub(i,j)=b(i,j)

omegstr=omegst*(1.-enbar*amass/rholl(j))

entest=enbar*amass

if(entest.ge.rholl(j))omegstr=0.

omegexb=(1.+ratrod)*omegstr

omeggb+=beta*omegstr*.5/(1.-beta)

xxx(i,j)=rholl(i,j)*(2.*omegexb+omeggb-omegst)

yyy(i,j)=-rholl(i,j)*(omegexb+omeggb)*(omegexb-omegst)

xxx(i,j)=xxx(i,j)*fir

yyy(i,j)=yyy(i,j)*fir

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 equillm

c.....equilibrium for tandem mirror electron ring plug

c....insert storage cliches 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/cel

psistr=-cel/(3.*del)

psibar=psistr+sqrt(psistr**2-be1/(3.*del))

p1max=cel+be1*psibar+cel*psibar**2+del*psibar**3

do 10 j=1,jx

psir=psi(j)/psim

psire=psi(j)/psime

```

pe1=(ae1+be1*psire+c1*psire**2+de1*psire**3)*hpme(j)
pe2=.5*(1.-tanh((psi(j)-psi0e)/wp2e))*(1.-hpme(j))
p1(j)=(pe1+pe2)/p1max
dp1dpsi(j)=(de1*3.*psire**2+
c be1*2.*ce1*psire)/(psime*p1max)*hpme(j)-.5*(1.-(tanh
c ((psi(j)-psi0e)/wp2e)**2)/(wp2e*p1max)*(1.-hpme(j))
p2t(j)=(1.-tanh((psi(j)-psi0e)/wp2e)*.5
dp2dpsi(j)=-.5*(1.-(tanh((psi(j)-psi0e)/wp2e)**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)=dp2dpsi(j)+dp3dpsi(j)*dip
hf1r(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,
hf1r(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.epspl) 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

```

```

15  phi1(i)=phi1c*exp(arg1)+phi1l/cosh(arg2)
c.....special phi1 for high mm rotation mode test
do 17 i=1,ix
  phi1(i)=phi1c*(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)*bmn1**3/3.+((bp2-bp1)*bmn1+dp1)+(cp1-cp2)
c /bmn1
  dp3=(ap3-ap2)*bvx2**3/3.+((bp3-bp2)*bvx2+dp2)+(cp2-cp3)
c /bvx2
  dt0=-8.*ct0/(bmax*3.)
  dt1=-8.*ct1/(bvx2*3.)
  dt2=-8.*ct2/(3.*bvx3)
  dt3=-8.*ct3/(3.*bm1)
  ppt=1./(1.-1./rp1**2)**2
  dter1=(-8./3-4.*((bvx3/bmn2)**3/3.+4.*((bvx3/bmn2))/bm2
  dter2=-als1*ppt*dter1*psi1s1
  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))
    ppart=(abp(i)*b4+bbp(i)*b2+cbp(i))
    pperte=(abq(i)*b4+bbq(i)*b2+cbq(i))
    pperp(i,j)=p2(j)*ppart
    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))+ncenter*cold*(1.+(rcwall-1.)*psi(j)/psi(jx)))
c..... special rho for exb rotation case
    rho(i,j)=amass*ncenter*exp(exrho*psi(j)/psi(jx))
    fac=b(i,j)*(1.+2.*p1(j)*(2.*abf(i)*b2+bbf(i)))
    fac1=dp1dpsi(i)*(abf(i)*b4+bbf(i)*b2+cbf(i))
    fac2=dp2dpsi(i)*(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 (-psi3)*hp3(j)
c..... special dphidpsi for constant omegexb
    dphidpsi=phi1(i)*er/cos
    epsi(i,j)=-dphidpsi
    dpdpsi=pperte*dp2dpsi(j)+p2(j)*dp2dbe*dbdpsi(i,j)
    omegci=echarg*b(i,j)/(lamass*cos)
    unit=1./(sqrt(4.*pi))

```

```

omegstr=-unit*b(i,j)*dpdpsi/(omegc*i*rhol(i,j))
omeggb=unit*pporte*p2(j)*dbdpsi(i,j)/(omegc*i*rhol(i,j))
omegexb=-cee*dphidpsi
xxx(i,j)=hfir(j)*rho(i,j)*(omeggb-omegstr+2.*omegexb)*sf6
yyy(i,j)=-hfir(j)*rho(i,j)*(omegexb+omeggb)*(omegexb-omegstr)*sf8

ppertr=(abf(i)*b4+bbf(i)*b2+cbf(i))
pperps(i,j)=pperpl(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(pperpl(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 errpri(i-1,j-1)=pperpl(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=dpsi(j)
psi_b=psi(j-1)
if(j.eq.2)then
dps=psi(2)
psi_b=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)hpk0(k)=1.
if(psik.lt.psim)hpkm(k)=1.
if(psik.lt.psime)hpkme(k)=1.
if(psik.lt.psi2.and.psik.gt.psi1)hpk12(k)=1.
29 continue
c..... cal. p1 (p1k) at intermediate points in dpsi interval
      do 30 k=1,kin
        psik=psib+dps*(k-1)/(kin-1)
psikr=psik/psim
psikre=psik/psime
        p1ek=(ae1+be1*psikre+ce1*psikre**2+de1*psikre**3)*hpkme(k)
p2ek=.5*(1.-tanh((psik-psi0)/wp2e1))*(1.-hpkm(k))
        p1k(k,j)=(p1ek+p2ek)/p1max
p3k=(p3a+p3b*psikr+p3c*psikr**2+p3d*psikr**3)*hpk0(k)
p2k(k,j)=(1.-tanh((psik-psi0)/wp21))*,.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.epspl) 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
continue
call simps(deli1,dapi1,kin,dps)

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      call simps(deli2,dapi2,kin,dps)
      call simps(deli3,dapi3,kin,dps)
      call simps(deli4,dapi4,kin,dps)
      call simps(deli5,dapi5,kin,dps)
      capi1=dapi1+sum1
      capi2=dapi2+sum2
      capi3=dapi3+sum3
      capi4=dapi4+sum4
      capi5=dapi5+sum5
      (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 rterm=dp2dpsi(j)*(ebp(i)*b(i,j)**4+fbp(i)*b(i,j)**2+gbp(i))
      c /(12.*p2(j)**.5)
      c +p2(j)**.5*(4*ebp(i)*b(i,j)**3+2*fbp(i)*b(i,j))*dbdpsi(i,j)
      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)**2*b(i,j))/uuz(i))
      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=max1(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
omegwkbb=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.0.1 then
gamwkbb=sqrt(-trad)/rhoave(j)
omeg1wkbb=omegwkbb
omeg2wkbb=0.
else
omeg1wkbb=omegwkbb +sqrt(trad)/rhoave(j)
omeg2wkbb=omegwkbb-sqrt(trad)/rhoave(j)
end if
70 continue
return
end
subroutine f1to11
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 cliche 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
dppdps1=dp2dps1(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))*dbdps1(i,j)
dpedps1=dp1dps1(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))*dbdps1(i,j)
oring(i,j)=dppdps1*dpedps1

```

```

g1(i,j)=+(mm*uuz(i))**2*r(i,j)*(rzz(i,j))
c*qubv(i,j)+dppdps*i*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*uuz(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). uses input fpsi, fv, fz, fu and azm, apsim .
c..... insert cliché storage here
use param
use const

xv=alog(fpsi)/alog(fv)
xu=alog(fz)/alog(fu)
zzp=0.
psip=0.
do 5 i=1,ix
5 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, flat 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+(j-2)*(i-2)
k=.5*(1+isw)*k1+.5*(1-isw)*k2
cosx=cos(.5*pi*(z(1)*kzsp)/zedge)
xro(k)=ex0*cosx*rbf*(r1+r2-1.)+ex1*cos(.5*pi*(z(1))/zedge)
xio(k)=ex0*cosx*rbf*(r3+r4-1.)+ex1*cos(.5*pi*z(1)/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=rangf(b1)
      r2=rangf(b1)
      r3=rangf(b1)
      r4=rangf(b1)
      kzsp=1
      endif
      15 continue
      k1=i-1+(j-2)*(lx-2)
      k2=j-1+(lx-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)
      xioo(k)=ex0*cosx*rbf*(r3+r4-1.)+ex1*cos(.5*pi*z(i)/zedge)
c      xioo(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=lx), 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=max (j=jx), fjrxx=-1, implies x=0,
c.....                  fjrxx=1, implies slope=0.
c.....      data mm/4/, bias/.5/, lmax/2/, nmax/5/, dv/1./, du/1./
c ,ndiag/100/, f11/1./, fizx/1./, fj1/1./, fjrxx/1./, fir/1./
c ,sf6/1./, sf8/1./, kplotm/0/, kzs/1/, swg1/1./, swg2/1./
c , swg3/1./, swg4/1./, non/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,non
c ,f11,fizx,fj1,fjrxx,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 dd0(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

```