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      subroutine adummy
c,,,,, cliché storage set up here

      cliché param
      parameter (lzx=66, jrx=30, lzx2=lzx-2 )
      parameter ( ltbw=2*lzx-1 )
      parameter (kxp=(lzx-2)*(jrx-2), lbw=lzx-1 )
      parameter (nplt=3000, nps=100 )
      parameter (lxf=4*(lzx-2),nfourxf=150)
      parameter (lxpg=2*lzx2)
      parameter (ksim=51)
      endcliché
      cliché 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/xloo(kxp),xlo(kxp),xlo1(kxp)
c ,xroo(kxp),xro(kxp),xro1(kxp)
      endcliché

      cliché 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)
      endcliché

      cliché const
      use param
      common/title/aname(5)
      common/con/gam1,gam2,ix,jx,mm,lzxp,kxx,nmax,lmax,lsw,lhbw
c ,fac1,fac2,bias,du,dv,dt,ndlag,ex0,b0,rho0,ex1,f11,f1zx,fj1
c ,fjrx,kplot,npm,fps1,fz,fu,fv,azm,apsim,u0,v0,amass
c ,fourpi,omegst,omegr,omegexb,flr,sf6,sf8,kplotm,kzs,zedge
c ,cpuo,clo,syso,va1fk,xu,xv,n,pl,vw,slw,dvin,dvout
      common/contm/
c psi0rel,psi1rel,psi2rel,z1rel,z2rel,z3rel,z0rel,nslosh,bmg
c ,ncenter,pslosh,pcenter,rp,ztrans,ltrans,bm,ltran,ztran,rp1
c ,bcen,pring,psp,phicen,phiplg,kin,xpot,ypot,wpot,pfudge,rpx
c ,phice,phipl,betslsh,betcent,z0,z1,z2,z3,z4,psi1,psi2,psi0
c ,betcene,betslse,psloshe,pcentee,bmax,alsi,bm1,psls1,cold
c ,p2wide,psi3rel,psi3,p1max,bv0,bv3,bv4,bceng,psloshin,psloshen
c ,nsloshin,pxp1,pxp2,p3a,p3b,p3c,p3d,psim,pe10,ael,bel,cel
c ,psi0erel,psi0,psime,p2ewide,wp2e
      common/mesh/psi(jrx),z(lzx),u(lzx),v(jrx),dpsi(jrx),dz(lzx)
c ,vpsi(jrx),uuz(lzx),vpsi1(jrx),uuzh(lzx)
      common/graf/ xrtime(nplt),xrspz(lzx,nps),xrpppsi(jrx,2*nps)
c ,time(nplt),xflute(lzx,nps)
      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(lzx),h1(lzx),h34(lzx),abp(lzx),bbp(lzx)
c ,cbp(lzx),abf(lzx),bbf(lzx),cbf(lzx),hp3(jrx),hp12(jrx)
c ,htrans(lzx),abq(lzx),bbq(lzx),cbq(lzx),h56(lzx),abr(lzx)

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c , bbr(izx), cbr(izx), betring, hp0(jrx), hpm(jrx), hpme(jrx)
c , 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), dp2dpsi(jrx), p3(jrx), dp3dpsi(jrx), hpkm(ksim)
c , hpme(ksim)
c , common/forced/nfour, nfourx, nfourmax, nfourp, jfour, lxp, locv

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

return
end

c , , , , the main routine

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

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

c , , , , flora1 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, fpsl*psimax=fv*vmax ,
c , , , , fz, fu, fpsl, fv, input, xu=ln fz / ln fu, xv=ln fpsl / 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

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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.....lequillrot 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.....le-ring pperp in b field only, and additional term in curvature
c.....drive ),

c.....flortm3, like flortm2 with corrections to pressusre 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 ( $=-d\phi/d\psi$ ) , 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=ae1+be1*(
c.....psi/psime)+ce1*(psi/psime)**2, and pe2=.5*(1-tanh((psi-psi0e)/p2ewide))
      use param
      use fstor
      use matrix
      use const

      data t' , e6/
      integer tallyb(2000b)
      common / q8iocs/iocf(0:15)
      common/pic100/npete
      data itally/1/

c.....call link call here
      call link('unit59=terminal,unit2=(inflt4,open),unit3=(output,
c      create) //')

      if(itally.gt.0) then
      do 200 ii=1,15
200   if(iocf(ii).eq.0)go to 210
      ii=0
210   ioc tally=ii
      ioc tally=14
      if(ioc tally.eq.0)go to 299
      call timer(ioc tally,'ztally00',tallyb,2000b,floralm,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

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nn1=jtbw+kxp
call memory(ww,nn-1)
call memory(ww1(nn),nn1)
call pstart(dev,4rplot,1,'box u21$',1)
c npete=1
call p100
call input
call inputtm
c call rigidcon
c call curvecon
call grid
call constant
call tmcon2
call equiltm
c call equilrot
c call equilcur
call fltol1
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(xloo(1),xio(1),kxx)
call banfac(kxp,ihbw,ww,1,-(kxp-1))
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 zmove wrd(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 zmove wrd(xrol,ww1(nn),kxx)
10 continue
call zmove wrd(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 zmove wrd(xiol,ww1(nn),kxx)
20 continue
fac1=-.5/dt
90 fac2=.5/dt
call zmove wrd(xloo,xio,kxx)
call zmove wrd(xio,xiol,kxx)
call zmove wrd(xroo,xro,kxx)
call zmove wrd(xro,xrol,kxx)
c,,,,, time array
xrtime(n)=xro(kplot)
if(mod(n,ndiag).eq.0)call diagno
if(mod(n,nfourp).eq.0)call fourier
100 continue

```

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call clsdsk(1ocv,0)
call timeused(1cp,1o,1sy)
cpuo=1cp*tim
cio=1o*tim
syso=1sy*tim
call picsher
call close(100)
call keep80(1,3)
call fr80id
call threed
call plote
call timend
call exit(1)
end
subroutine constant

```

```

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

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

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subroutine curvecon

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

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c,,,,, insert storage clichés here
use param
use const
real klbsq

```

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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,lb,rw,xm
c , zol,dtrel,theta0
call ddi(curve,2,3,1)
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

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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 inputtm

c..... calculates grid quantities needed by sub. grid
c..... in floratm version

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

c.....glossary of input parameters
c    psi0rel,..value of psi (in relative units) where plasma pperp
c    profile is half the maximum.
c    p2wide,..parameter inversely proportional to "ramp width" of p2
c    psi0erel,..relative value of psi where ering pperp (pe2) is half
c    the maximum
c    p2ewide,..parameter inversely proportional to "ramp width" of pe2
c    psi1rel,..lower limit of psi where pring =0.
c    psi2,..upper limit of psi where pring=0.
c    z0rel,..axial center of plug
c    z1rel,..axial position of plug inward mirror
c    z2rel,..axial position of plug outboard mirror
c    z3rel,..inner axial position where pring=0.
c    z4rel,..outer axial position where pring=0.
c    nslosh,..peak sloshing ion number density
c    ncenter,..peak center cell ion density
c    betslsh,..peak sloshing ion beta-perp w.r.t. bvac(z0)
c    betcent,..peak center cell ion beta-perp w.r.t. bvac(z=0)
c    betslse,..peak sloshing electron beta-perp w.r.t. bvac(z0)
c    betcene,..peak center cell electron beta-perp w.r.t. bvac(z=0)
c    betring,..peak electron ring beta perp w.r.t. bvac(z0)

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c      rp,...plug vacuum mirror ratio
c      ztrans,...begining of center cell transition region
c      ltrans,...transition length
c      bmg,...vacuum mirror peak (approx.) in gauss
c      bceng,...center cell bmax in gauss
c      epsp,...minimum pressure, below which b=bvac
c      phicen,...center cell potential relative to plug ion t perp
c      phiplg,...plug potential relative to plug ion t perp
c      kin,...number of points in slmsons rule quadratures
c      xpot,...exponent coefficient in center cell potential
c      wpot,...exponent coefficient in plug potential
c      ypot,...power of polynomial in potential psi dependence
c      pxp1,...power of polynomial in e ring pperp(psi) (p1(psi))
c      pxp2,...power of polynomial in e ring pperp(psi) (p1(psi))
c      zmax,...axial end of system in cm.
c....e ring pperp at psi=0,

c..... input for tandem mirror equilb,
      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      , xpot/1./, ypot/2./, wpot/2./, kin/5/psi1rel/2./, psi2rel/1.5/
c      , psi0rel/.50/, zmax/100./, z1rel/.5/, z2rel/1./, z0rel/.75/
c      , z3rel/.625/, z4rel/.875/, ztrans/.4/, ltrans/.05/, p2wide/.1/
c      , bmg/1.e4/, bceng/1.e3/, nsloshin/2.e13/, ncenter/1.e13/,
c      , betslsh/.25/, betcent/.10/, rp/4./, epsp/1.e-6/, rp1/2./
c      , phicen/.1/, phiplg/.1/, betcene/.1/, betslse/.1/
c      , dt/1.e-5/, pfudge/0./, cold/1.e-3/, betring/0./, psi3rel/1.05/
c      , pxp1/2./, pxp2/2./, safe/.9/, pe10/0./, psi0erel/.5/, p2ewide/.2/
      namelist/curve/echarg, lb, rw, xm, rw1
c      , zmax, dt, theta0, pfudge
c      , psi0rel, z1rel, z2rel, z0rel, nsloshin, cold, z3rel, z4rel
c      , ncenter, betslsh, betcent, rp, ztrans, ltrans, bmg, rp1
c      , bceng, epsp, phicen, phiplg, kin, xpot, ypot, wpot, betcene, betslse
c      , betring, psi1rel, psi2rel, p2wide, psi3rel, pxp1, pxp2, safe, pe10
c      , p2ewide, psi0erel
      call ddi(curve,2,3,1)
      call pframe
      call p100
      call ddo(curve,100,0,1)
c..... transform input z into physical units
      z0=z0rel*zmax
      z1=z1rel*zmax
      z2=z2rel*zmax
      z3=z3rel*zmax
      z4=z4rel*zmax
      ztran=ztrans*zmax
      ltran=ltrans*zmax
      bm=bmg/sqrt(4.*pi)
      bcen=bceng/sqrt(4.*pi)
c..... generate zmax, psimax, u(ix), v(jx), du, dv
      asq=.25/((2.*rp)**(2./3.)-1.)*(z2-z1)**2
      argt1=+(z1-ztran)/ltran
      bv0=bm*(1./((1.+(z0-z1)**2/asq)**1.5+1./((1.+(z0-z2)**2/asq)
c      **1.5)
      bmax=bm*(1.+1./((1.+(z1-z2)**2/asq)**1.5) +bcen*.5*
c      (1.-tanh(argt1))
      bv4=bm*(1./((1.+(z1)**2/asq)**1.5+1./((1.+(z2)**2/asq)**1.5)
      psimax=rw**2*bmax*.5
      psiw=rw1**2*bmax*.5
      psi0=psimax*psi0rel

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    psi0e=psimax*psi0ere1
    psi3=psimax*psi3rel
    psi1=psimax*psi1rel
    psi2=psimax*psi2rel
    v(jx)=psimax
    u(ix)=zmax
    du=u(ix)/(ix-1.5)
    dv=v(jx)/(jx-1.5)
c.....transform input constants from relative to physical units
    pmag0=(bv0)**2*.5
    pmag1=(bv4+bcen)**2*.5
    psloshin=(betslsh)*pmag0
    pcenter=(betcent)*pmag1
    psloshen=betslse*pmag0
    pcentee=betcene*pmag1
    pring=betring*pmag0
    phice=phicen*psloshin/(nsloshin*echarg)
    phipl=phiplg*psloshin/(nsloshin*echarg)
c.....calculate maximum p1(psi)
    psidif=psi2-psi1
    plmax=(pxp1*psidif/((pxp1+pxp2)*psi1))*pxp1*(pxp2*psidif/
    c ((pxp1+pxp2)*psi2))*pxp2
c.....check that kin is not too big
    ksimp=ksim
    if(kin.gt.ksimpl)then
        kin=ksimp
        write(59,200)kin
200    format('kin initially too large, reduced to',i4)
        end if
c..... check that kin is odd for simpsons quadratures
    kch=mod(kin,2)
    if (kch.eq.0)then
        kin=kin-1
        write(59,201)kin
201    format('kin initially even, changed to',i4)
        end if
c..... adjust epsp to accomodate zero pressure case
    epsp=epsp*(1.e-4+psloshin+pcenter)/(1.e-8*epsp+psloshin+pcenter)
c.....check that bcen and betcen are physically consistent
    asq=.25/((2.*rp)**(2./3.)-1.)*(z2-z1)**2
    bv0=bm*(1./(1.+(z0-z1)**2/asq)**1.5+1./(1.+(z0-z2)**2/asq)
    c**1.5)
    argt3=(z3-ztran)/ltran
    bv3=bm*(1./(1.+(z3-z1)**2/asq)**1.5+1./(1.+(z3-z2)**2/asq)
    c**1.5)+bcen*.5*(1.-tanh(argt3))
    rpx=bmax/(bv0)
    rp3=bv3/bv0
    afac=(rp*(1.-1./rp**2))**2/bm**2
    afac1=(rp*(1.-1./rp3**2))**2/bm**2
    tcon6=pring/((1.-1./rp3**2)**2)
    tmin=afac*(bcen+bv4)**2
    if(tmin.le.betcent+betcene)then
        bcen1=sqrt((betcent+betcene)/afac)-bv4
        tmine=tmin-betcene
        write(59,210)bcen1,tmine
210    format('initial central cell beta and b are inconsistent for',
    c ' equilibrium',/'either increase bcen to more than',e14.6/
    c 'or decrease betcent to less than',e14.6/'use namelist',
    c ' data format')
        namelist/better/betcent,bcen

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        read(59,better)
        pcenter=betcent*pmag0
        end if
c..... check plug ion beta and e ring beta
        tmin1=afac*bv0**2
        if (tmin1.le.(betslsh+betslse))then
            tbetsl=tmin1-betslse
            tbetr=(-betslsh+tmin1)
            delbet=betslsh+betslse-tmin1
            write(59,220)delbet,tbetsl,tbetr
            namelist/fix/betslsh,betslse
220    format('the sum of (betslsh+betslse) is too large by',
c    'e14.6/'either decrease betslsh to',e14.6/'or decrease
c    betslse to',e14.6/'or something (use namelist data format)')
            read(59,fix)
            psloshin=betslsh*pmag0
            psloshen=betslse*pmag0
            end if

c..... check betring in plug
        if(tcon6.ge.(safe*.5*bv0**2))then
            ratio=safe*.5*bv0**2/tcon6
            pring=pring*ratio
            ratinv=1./ratio
            write(59,231)ratinv,pring
231    format('pring was too large by a factor ('e14.6,') and was',
c    'reduced to ',e14.6 )
            end if

c..... check that rp1 is less than rpx
        if(rp1.ge.rpx)then
            rp1t=rp1
            rp1=.5*rpx
            write(59,230)rp1t,rpx,rp1
230    format('rp1 ('e12.6,') was larger than rpx ('e12.6,') and was',
c    'reduced to ',e12.6 )
            end if
        end
        subroutine tmcon2

c..... calculates equilibrium quantities for tandem mirror

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

c.....calculate heaveside step functions h12, h1, h34
        call bcast(h12(1),0.,ix)
        call bcast(h1(1),0.,ix)
        call bcast(h34(1),0.,ix)
        call bcast(htrans(1),0.,ix)
        do 2 i=1,ix
            if(z(i).ge.z1.and.z(i).le.z2)h12(i)=1.
            if(z(i).le.z1)h1(i)=1.
            if(z(i).ge.ztran)htrans(i)=1.
2        if(z(i).ge.z3.and.z(i).le.z4)h34(i)=1.

        psim=psi0*(1.-p2wide)
        psime=psi0e*(1.-p2ewide)

```

```

c.....calculate heaveside step functions hp0(j), hp3(j) and hp12(j)
  call bcast(hp12(1),0.,jx)
  call bcast(hp3(1),0.,jx)
  call bcast(hp0(1),0.,jx)
  call bcast(hpm(1),0.,jx)
  call bcast(hpme(1),0.,jx)
  do 4 j=1,jx
    if(psi(j).le.psi0)hp0(j)=1.
    if(psi(j).le.psi3)hp3(j)=1.
    if(psi(j).le.psim)hpm(j)=1.
    if(psi(j).le.psime)hpme(j)=1.
4    if(psi(j).ge.psi1.and.psi(j).le.psi2)hp12(j)=1.

c..... calculate vacuum b field (bvac) and derivatives dbv/dz, d2bv/dz2
  asq=.25/((2.*rp)**(2./3.)-1.)*(z2-z1)**2
  bv0=bm*(1./(1.+(z0-z1)**2/asq)**1.5+1./(1.+(z0-z2)**2/asq)
  c**1.5)
  do 5 i=1,ix
    eb1=(z(i)-z1)/asq
    eb2=(z(i)-z2)/asq
    fb1=sqrt(1.+eb1*(z(i)-z1))
    fb2=sqrt(1.+eb2*(z(i)-z2))
    argt=(z(i)-ztran)/ltran
    dbvdz(i)=-3*bm*(eb1/fb1**5+eb2/fb2**5)-bcen*.5*(1.-
  c tanh(argt)**2)/ltran
    d2bvdz(i)=-3*bm*(1./(asq*fb1**5)+1./(asq*fb2**5)-eb1**2*5/
  cfb1**7-eb2**2*5/fb2**7)+bcen*tanh(argt)*(1.-tanh(argt)**2)/
  c ltran**2
5    bvac(i)=bm*(1./fb1**3+1./fb2**3)+bcen*.5*(1.-tanh(argt))
    argt3=(z3-ztran)/ltran
    bv3=bm*(1./(1.+(z3-z1)**2/asq)**1.5+1./(1.+(z3-z2)**2/asq)
  c**1.5)+bcen*.5*(1.-tanh(argt3))
    rp3=bv3/(+bv0)
    bm1=rp1*(bv0)
    alsit=((1.-1./rp1**2)/(1.-1./rp3**2)*(rp1/rp3))**2
    als1=alsit+.5*(1.-alsit)*prudge

c.....calculate heaveside function h56(i) for second component of
c..... sloshing pperp
  do 10 i=1,ix
    eb2=(z(i)-z2)/asq
    if(z(i).gt.z1.and.z(i).lt.z2.and.bm1.ge.bvac(i))
  c h56(i)=1.
10  continue

c..... calculate pressure coeff. including electron ring (pr)
  con1=1./((1.-1./rp3**2)**2*bmax**4)
  con2=con1*(-2*bmax**2)
  con3=-.5*con2*bmax**2
  con4=pring/((1.-1./rp3**2)**2*bv3**4)
  con5=con4*(-2*bv3**2)
  con6=-.5*con5*bv3**2
  don1=1./((1.-1./rp1**2)**2*bm1**4)
  don2=don1*(-2*bm1**2)
  don3=-.5*don2*bm1**2

c..... calculate constants for p3 pressure componetnt
  data p30/.01/
  wp2=p2wide*.5*psi0
  wp2e=p2ewide*.5*psi0e
  wp3=wp2/psim

```

```

    psi0r=psi0/psim
    p3b=p30
    p3c=.5*(1.-(tanh(2.))**2)/wp3-p3b*p2wide*(2.-p2wide)
c *.5/p2wide
    p3d=-(p3b+2.*p3c*psi0r)/(3.*psi0r**2)
    p3a=-(p3b*psi0r+p3c*psi0r**2+p3d*psi0r**3)

c.....calculate constants for pe1
    g2p=-.5*(1.-(tanh(-2.))**2)/wp2e
    g2f=.5*(1.-tanh(-2.))
    ae1=pe10
    ce1=psime*g2p-g2f+ae1
    be1=2.*(g2f-ae1)-psime*g2p
c.....readjust peak plug pressure for sloshing
    rtemp=(bm1/bmax)**2
    rpxi2=1./rpx**2
    ter1=(rtemp-rpxi2)**2
    ter2=(1.-rpxi2)**2
    bstr2=bmax**2*(-ter1+alsi*rtemp*ter2)/(-ter1+alsi*ter2)
    pfloat=((con1-don1*alsi)*bstr2**2+(con2-don2*alsi)*bstr2
c +con3-don3*alsi)*(1.+p3a)
    pslosh=psloshin/pfloat
    psloshe=psloshen/pfloat
    nslosh=nsloshin/pfloat
    psls1=pslosh+psloshe
    pcen1=pcenter+pcentee
    do 6 i=1,ix
    eon1=con1-don1*alsi*h56(i)
    eon2=con2-don2*alsi*h56(i)
    eon3=con3-don3*alsi*h56(i)
    abf(i)=con4*h34(i)
    bbf(i)=con5*h34(i)
    cbf(i)=con6*h34(i)

    abp(i)=eon1*(psls1*h12(i)+pcen1*h1(i))
    bbp(i)=eon2*(psls1*h12(i)+pcen1*h1(i))
    cbp(i)=eon3*(psls1*h12(i)+pcen1*h1(i))

    abr(i)=eon1*(nslosh*h12(i)+ncenter*h1(i))
    bbr(i)=eon2*(nslosh*h12(i)+ncenter*h1(i))
    cbr(i)=eon3*(nslosh*h12(i)+ncenter*h1(i))

    abq(i)=eon1*(pslosh*h12(i)+pcenter*h1(i))
    bbq(i)=eon2*(pslosh*h12(i)+pcenter*h1(i))
6   cbq(i)=eon3*(pslosh*h12(i)+pcenter*h1(i))
    return
end

subroutine equiltm

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/ce1
plmax=ae1+be1*psibar+ce1*psibar**2

do 10 j=1,jx
psir=psi(j)/psim
psire=psi(j)/psime
pe1=(ae1+be1*psire+ce1*psire**2)*hpme(j)
pe2=.5*(1.-tanh((psi(j)-psi0e)/wp2e))*(1.-hpme(j))
p1(j)=(pe1+pe2)/plmax
dp1dpsi(j)=(be1+2.*ce1*psire)/(psime*plmax)*hpme(j)-.5*(1.-tanh
c ((psi(j)-psi0e)/wp2e)**2)/(wp2e*plmax)*(1.-hpme(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)
dp2dpsi(j)=dp2dpst(j)+dp3dpsi(j)
if((p2(j)).lt.epsp)go to 50
do 11 i=1,ix
if(abp(i).eq.0)then
b(i,j)=bvac(i)
go to 11
end if
u1=p2(j)*abp(i)+p1(j)*abf(i)
u2=p2(j)*bbp(i)+.5*p1(j)*bbf(i)
t1=-u2*.5/u1
u3=p2(j)*cbp(i)-bvac(i)**2*.5+p1(j)*cbf(i)
radic=sqrt(t1**2-u3/u1)
bsq1=t1+radic
bsq2=t1-radic
bsq=bsq1
if(t1.gt.radic)bsq=bsq2
b(i,j)=sqrt(bsq)
11 continue
go to 10
c.... b=bvac if p2.lt.epsp
50 do 51 i=1,ix
51 b(i,j)=bvac(i)
10 continue

c.....loop 15, calculate phi, the electric potential
do 15 i=1,ix
arg1=((z(i)-z1)/(z1-z0))**2*(-xpot)*(1.-h1(i))
arg2=wpot*((z(i)-z0)/(z0-z1))**2
15 phi1(i)=phice*exp(arg1)+phipl/cosh(arg2)
do 16 j=2,jx
phi2(j)=(1.-(psi(j)*hp3(j)/psi3)**ypot)*hp3(j)
16 continue
c.....loop 20, calculate pperp, ppar and db/dpsi
ppt=1./(1.-1./rp1**2)**2
dter1=(-8./3-4.*(bmax/bm1)**3/3.+4.*(bmax/bm1))/bm1
dter2=-alsi*ppt*dter1*psls1
do 20 i=1,ix
do 20 j=1,jx
b2=b(i,j)**2
b4=b2**2
dter3=dter2*h56(i)
dter=(4.*abp(i)*bmax**2/3.+2.*bbp(i))*bmax+dter3
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)
rho(i,j)=amass*(p2(j)**.5*(abr(i)*b4+bbr(i)*b2+cbr(i))+
c p2(j)**.5*ncenter*cold)
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.
dp2db0=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))
phang=psi(j)*hp3(j)
dphidpsi=phi1(i)*ypot*(phang/psi3)**(ypot-1.)/
c (-psi3)*hp3(j)
epsi(i,j)=-dphidpsi
dpdpsi=pperte*dp2dpsi(j)+p2(j)*dp2db0*dbdpsi(i,j)
omegci=echang*b(i,j)/(amass*cee)
omegstr=-b(i,j)*dpdpsi/omegci
omeggb=pperte*p2(j)*dbdpsi(i,j)/omegci
omegexb=-rho(i,j)*cee*dphidpsi*.5
xxx(i,j)=(omeggb-omegstr+2.*omegexb)*sf6
yyy(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)
dter3=dter2*h56(i)
dter4=(4.*abp(i)*bmax**2/3.+2.*bbp(i))*bmax+dter3
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 /dolb
      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=dpsi(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(i),0.,kin)
      call bcast(hpk12(i),0.,kin)
      call bcast(hpkm(i),0.,kin)
      call bcast(hpkme(i),0.,kin)
      do 29 k=1,kin
      psik=psib+dps*(k-1)/(kin-1)
      if(psi0.lt.psim)hpk0(k)=1.
      if(psim.lt.psime)hpkm(k)=1.
      if(psime.lt.psi2)hpkm(k)=1.
      if(psi2.and.psi1.lt.psim)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)*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)*hpk0(k)
      p2k(k,j)=(1.-tanh((psik-psi0)/wp2))*p3k
30   continue
31   continue
      do 32 k=1,kin
      if((p2k(k,j)).le.epsp.or.abp(i).eq.0) then
      bk=bvac(i)
      else
      u1=p2k(k,j)*abp(i)+p1k(k,j)*abf(i)
      u2=p2k(k,j)*bbp(i)+p1k(k,j)*bbf(i)+.5
      t1=-u2*.5/u1
      u3=p2k(k,j)*cbp(i)+p1k(k,j)*cbf(i)-bvac(i)**2*.5
      radic=sqrt(t1**2-u3/u1)
      bsq1=t1+radic
      bsq2=t1-rad ic
      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)
      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)
      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)
      flute1(j)=ans1*(ia-1)
      flute2(j)=(ans2*(mm**2-1)+ans1)*(ia-1)
      flute3(j)=-ans1/ans3
      rhoave(j)=ans3*(ia-1)
      xxave(j)=ans4*(ia-1)
      yyyave(j)=ans2*(ia-1)
60  continue

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)
      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 simpso(fin,fout,knn,df)
c.....simpsons rule quadratures, knn must be odd
      dimension fin(1)
      nse=(knn-1)/2
      se=ssum(nse,fin(2),2)
      nso=nse-1
      so=ssum(nso,fin(3),2)
      fout=df/(3.*(knn-1))*(fin(1)+fin(knn)+4.*se+2.*so)
      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(jx)
10  continue
      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 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(fourpi)
omegr=ratrod*omegst*(1.-enbar/en0)
foursq=sqrt(fourpi)
do 5 i=1,ix
uz=uuz(i)
do 5 j=1,jx
fac=exp(psi(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=log(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.
omegexb=(1.+ratrod)*omegstr
omeggb=+beta*omegstr*.5/(1.-beta)
xxx(i,j)=rho(i,j)*(2.*omegexb+omeggb-omegst)
yyy(i,j)=-rho(i,j)*(omegexb+omeggb)*(omegexb-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 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 cliché 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(kzs.eq.0)then
      rbf=1./(r(i,j)*b(i,j))
      else
      r1=ranf(b1)
      r2=ranf(b1)
      r3=ranf(b1)
      r4=ranf(b1)
      endif
5      continue
      k1=i-1+(j-2)*(ix-2)
      k2=j-1+(jx-2)*(i-2)
      k=.5*(1+isw)*k1+.5*(1-isw)*k2
      xro(k)=ex0*rbf*(r1+r2-1.)+ex1*cos(.5*pi*(z(i))/zedge)
      xio(k)=ex0*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(kzs.eq.0)then
      rbf=1./(r(i,j)*b(i,j))
      else
      r1=ranf(b1)
      r2=ranf(b1)
      r3=ranf(b1)
      r4=ranf(b1)
      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
      xroo(k)=ex0*rbf*(r1+r2-1.)+ex1*cos(.5*pi*(z(i))/zedge)
      xioo(k)=ex0*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=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.
c.....      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./

c.....,forced data loaded for testing fourier analyses and zed file
c.....,maker
c.....      data jfour/1/,nfourp/1/,nfourmax/5/
c.....      namelist/nw1/aname,mm, bias, lmax, nmax,ndiag
c.....      , f11,fizx,fj1,fjrx,ex0,ex1,fpsi,fu,fv,fz
c.....      , kplotm,kzs,jfour,nfourp,nfourmax,sf6,sf8,swg1,swg2,swg3,swg4

c.....      call ddi(nw1,2,3,1)
c.....      call ddo(nw1,100,0,1)
c.....      jx=jrx
c.....      kxx=kxp
c.....      ix=izx
c.....      return
c.....      end

c.....      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
c.....      use param
c.....      use const

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

c.....      uuzh(i)=(u(i)+.5*du)**(1.-xu)/(xu*azm)
c.....      dz(i)=z(i)-zzp
c.....      zzp=z(i)
10     continue
c.....      zedge=azm*.5*(u(ix)**xu+u(ix-1)**xu)
c.....      do 15 j=1,jx
c.....      v(j)=v0+(j-1.5)*dv
15     continue
c.....      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
      dpsl(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 f1to11
c.....calculates the f1 to f11 functions needed to generate the a and
c..... b matrices , uses the equilibrium quantities n, 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)
      g1(i,j)=+(mm*uuz(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*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 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

```

```

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))* .5
bijph=(b(i,j)+b(i,j+1))* .5
bip1jph=(b(i+1,j+1)+b(i+1,j))* .5
bip1jmh=(b(i+1,j-1)+b(i+1,j))* .5
bim1jph=(b(i-1,j+1)+b(i-1,j))* .5
bim1jmh=(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*vpsi(j)

```

```

f1ijmh=(f1(i,j)+f1(i,j-1))*5*vpsih(j-1)
f5ijph=(f5(i,j)+f5(i,j+1))*5*vpsih(j)
f5ijmh=(f5(i,j)+f5(i,j-1))*5*vpsih(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*bim1jmh*g4imhjmh*bijmh*uzbar*vpsih(j-1)
c *r(i-1,j-1)
a2(k,1)=-gam2*bim1jmh*g4imhjmh*bijmh*uzbar*vpsih(j-1)
c *r(i-1,j-1)
a3(k,1)=-gam3*bim1jmh*g4imhjmh*bijmh*uzbar*vpsih(j-1)
c *r(i-1,j-1)

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

a1(k,3)=-gam1*bip1jmh*g4iphjmh*bijmh*uzbar*vpsih(j-1)*r(i+1,j-1)
a2(k,3)=-gam2*bip1jmh*g4iphjmh*bijmh*uzbar*vpsih(j-1)*r(i+1,j-1)
a3(k,3)=-gam3*bip1jmh*g4iphjmh*bijmh*uzbar*vpsih(j-1)*r(i+1,j-1)
a1(k,4)=gam1*((bim1jmh*g4imhjmh*vpsih(j-1)*bijmh+bim1jph*g4imhjph
c *vpsih(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*((bim1jmh*g4imhjmh*vpsih(j-1)*bijmh+bim1jph*g4imhjph
c *vpsih(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*((bim1jmh*g4imhjmh*vpsih(j-1)*bijmh+bim1jph*g4imhjph
c *vpsih(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)*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/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)*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/vpsi(j))
a3(k,5)=((f1ijph+f1ijmh)/dv2-f2(i,j))/dt2+gam3*(-(f5ijph+f5ijmh
c )/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/vpsi(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/vpsi(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/vpsi(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/vpsi(j))

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```

      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))*5*vpsih(j-1)
      f3ijph=(f3(i,j)+f3(i,j+1))*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
      sfi1=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.sfi1.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.sfi1.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
      if(sfjrx.eq.1.and.sfizx.eq.1)fac4=r(ix-1,jx-1)*b(ix-1,jx-1)/

```

```

c (r(1,x,jx-1)*b(1,x,jx-1))
a1(      ,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
if(sfizx.eq.1.)sfizx=r(ix-1,j)*b(ix-1,j)/(r(ix,j)*b(ix,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 12 m=2,8,3
a1(k,m)=a1(k,m)+sfizx*a1(k,m+1)
a2(k,m)=a2(k,m)+sfizx*a2(k,m+1)
a3(k,m)=a3(k,m)+sfizx*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,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

```

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)
10    continue
      return
      end

```

subroutine right

c..... calculates right hand side vector for both equations,
 c..... rhs1(k)=2*a2*xr(n)-a2*xr(n-1)+b1*(xi(1)-xi(n-1)) , and
 c..... rhs2(k)=2*a3*xi(n)-a2*xi(n-1)+b1*(xr(1)-xr(n-1)) .

c..... insert cliches for storage here

```

      use param
      use fstor
      use matrix
      use const

      do 10 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+.5*(1-isw)*k2
      t1=0.
      t2=0.
      t3=0.
      tt1=0.
      tt2=0.
      tt3=0.
      do 5 m=1,9
      ip=i-2+m-((m-1)/3)*3
      jp=j-1+(m-1)/3
      if(jp.eq.1.or.jp.eq.jrx.or.ip.eq.1.or.ip.eq.izx)go to 5
      kp1=ip-1+(ix-2)*(jp-2)
      kp2=jp-1+(jx-2)*(ip-2)
      kp=.5*(1+isw)*kp1+.5*(1-isw)*kp2
      t1=t1+a2(k,m)*xroo(kp)
      t2=t2+a3(k,m)*xro(kp)
      tt1=tt1+a2(k,m)*xioo(kp)
      tt2=tt2+a3(k,m)*xio(kp)
5    continue
      do 6 mn=1,3
      jq=j-2+mn
      if(jq.eq.1.or(jq.eq.jrx)go to 6
      kq1=i-1+(ix-2)*(jq-2)
      kq2=jq-1+(jx-2)*(i-2)
      kq=.5*(1+isw)*kq1+.5*(1-isw)*kq2
      t3=t3+b1(k,mn)*(xiol(kq)-xioo(kq))
      tt3=tt3+b1(k,mn)*(xrol(kq)-xroo(kq))
6    continue
      rhs1(k)=(2.*t2-t1+fac1*t3)
      rhs2(k)=2.*tt2-tt1+fac2*tt3

```

```

10  continue
    return
    end
    subroutine rightvec

c..... calculates right hand side vector for both equations,
c..... rhs1(k)=2*a2*xr(n)-a2*xr(n-1)+b1*(xi(1)-xi(n-1)) , and
c..... rhs2(k)=2*a3*xi(n)-a2*xi(n-1)+b1*(xr(1)-xr(n-1)) .

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

    call sscal(kxx,0.,rhs1,1)
    call sscal(kxx,0.,rhs2,1)
    do 100 m=1,9
        mdel=(m-1)/3
        m1=m-1
        koff1=-mdel*5+m+(mdel-1)*ix
        koff2=(m-(mdel+1)*3-1)*jx+1-2*m1+7*mdel
        koff=.5*(1+isw)*koff1+.5*(1-isw)*koff2
        do 110 k=1,kxx
            rhs1(k)=rhs1(k)+2.*a3(k,m)*xro(k+koff)-a2(k,m)*xroo(k+koff)
110      rhs2(k)=rhs2(k)+2.*a3(k,m)*xio(k+koff)-a2(k,m)*xioo(k+koff)
            if(m.eq.2.or.m.eq.5.or.m.eq.8)go to 119
        go to 100
119      continue
        do 120 k=1,kxx
            mbar=m-1-(m/4)*2
            rhs1(k)=rhs1(k)+fac1*b1(k,mbar)*(xio1(k+koff)-xioo(k+koff))
120      rhs2(k)=rhs2(k)+fac2*b1(k,mbar)*(xro1(k+koff)-xroo(k+koff))
100      continue
    return
    end
    subroutine rigidcon

c..... special constants needed for rigid rotor equilibrium.

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

c..... input for rigid rotor
    data echarg/4.8e-10/, en0/1.00e+12/, b0/1.e4/, amass/3.34e-24/
    c , cee/3.e10/, valfk/.,4/, fourpi/12.56637/, pi/3.1415926/
    c , enbar/0.e11/
    namelist/rotor/b0,beta0,rotrod,valfk,en0,echarg,r0,rwb,enbar
    call ddi(rotor,2,3,1)

    carg=sqrt(1.-beta0)
    r0sq=r0**2
    aasq=r0sq*(1.-carg)/beta0
    cr=.5*(alog(1.+carg)-alog(1.-carg))
    aa=sqrt(aasq)
    rw=rwb*aa
    valf=b0/(sqrt(fourpi*en0*amass))
    omegc1=echarg*b0/(amass*cee)
    omegp2=fourpi*en0*echarg**2/amass

```

```

      omegst=2.*beta0*omegc1*cee**2/(omegp2*r0sq)
      vomeg=echang*sqrt(len0*fourpi/amass)*r0sq*.5/(beta0*cee)
      u(ix)=(pi*val/(2*omegst*valfk))/(1-.5/(ix-1.5))
      du=(u(ix)-u0)/(ix-1.5)
      targ=cosh(rw**2/r0sq+cr)*sqrt(beta0)
      v(jx)=b0*r0sq*.5*alog(targ)/sqrt(fourpi)
      dv=(v(jx)-v0)/(jx-1.5)
      return
    end

    subroutine mymove(a,b,len)
      dimension a(1),b(1)
      do 10 i=1,len
        a(i)=b(i)
10      continue
      return
    end

    subroutine picsher
c...uses grafic, graflib and grafcore to make plots of xr vs. time and
c....space.

c.....insert cliché for common here
      use param
      use fstor
      use matrix
      use const

      dimens on iy(2), it(5), lab(2),dum(izx2),ymin(5),ymax(5)
c ,dum1(nplt),rplot(jrx-1)
      data epp/1.e20/
      call orgfile(ume)
      call pframe
      call p100
      write(100,102)nume
      write(100,400)aname
400  format(/10x,'problem identification : ',5a8 )
102  format(10x,'this problem run by ',a8 )
      write (100,101) dt, ix,jx,nmax,lmax,bias,
c flr,sf6,sf8,kplot,kzs,cpuo,cio,syso
c ,xu,xv,bv0,bm1,bv3
101  format(////'dt=',e16.6,4x,'ix=',i8,4x,'jx=',i8/'total time steps =
c ',i8,4x,
c 'no. of iterations =',i8,4x,'bias=',f10.5/
c 'flr=',e16.8,4x,'sf6=',e16.8,4x,
c 'sf8=',e16.8/'kplot=',i8,4x,'kzs=',i8/
c 'cpu time =',e16.8/
c 'i-o time =',e16.8/'sys time =',e16.8/3x,'u exponent (xu) =
c ',e16.8/3x,'v exponent (xv) =',e16.8/'bv0=',e16.8,4x,'bm1=',
c e16.8,4x,'bv3=',e16.8/)
c.....plot coordinate stretching

      kx='u$'
      iy(1)='z$'
      it(1)='z vs u, '
      it(2)='(z=const'
      it(3)='*u**xu)$'
      call pframe
      call pscalr(0,u(2),u(ix),z(2),z(ix),1)

```