

| Course Number | AE8112 |
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| Course Title | Computational Fluid Dynamics and Heat Transfer |
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| Instructor | Dr. Seth Dworkin |

Problem Set 5

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|--------------------------|---------------|
| Programing Language Used | Fortran90 |

| Student Name | Student Number | |
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Q1a] using explicit finite difference, we can discretize the given vorticity - velocity eqn, with the below Tylor's series expansion centered at On

$$\frac{\partial^{2} \emptyset}{\partial x^{2}} \Big|_{i} = \frac{\phi_{i-1} - 2\phi_{i} + \phi_{i+1}}{\Delta x^{2}}$$
and
$$\frac{\partial \emptyset}{\partial x} \Big|_{i} = \frac{\phi_{i+1} - \phi_{i-1}}{Z\Delta x}$$

$$(1.1)$$

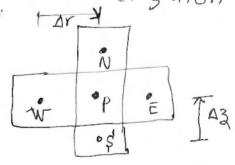
Applying this to our governing equations, we will have, $\emptyset = W$, $V_r \in V_z$ and for our z-D coordinate

Now considering vorticity transport eqn., so that after applying the product rule, we have,

$$g\frac{\partial w}{\partial t} = \mu \frac{\partial^2 w}{\partial r^2} + \mu \frac{\partial^2 w}{\partial z^2} + \mu \left(-\frac{1}{2}w + \frac{1}{7} \frac{\partial w}{\partial r} \right) - \beta v_r \frac{\partial w}{\partial r}$$

$$-\beta v_z \frac{\partial w}{\partial z} + \beta v_r w$$

so that applying the standard forward difference for the time discretization and eqn (1.1), (1.2), we get,



Q1 a] contid.

$$\mathcal{S}\left(\frac{\omega_{p}^{n+1}-\omega_{p}^{n}}{\Delta t}\right) = \mathcal{M}\left(\frac{\omega_{w}^{2}-2\omega_{p}^{2}+\omega_{E}^{2}}{\Delta r^{2}}\right) + \mathcal{M}\left(\frac{\omega_{s}^{2}-2\omega_{p}^{2}+\omega_{n}^{2}}{\Delta 3^{2}}\right) - \frac{\mathcal{M}\omega_{p}^{2}}{r_{p}^{2}} + \frac{\mathcal{M}\left(\frac{\omega_{s}^{2}-\omega_{w}^{2}}{\Delta r^{2}}\right) - \mathcal{S}v_{r}^{2}\left(\frac{\omega_{w}^{2}-\omega_{w}^{2}}{2\Delta r^{2}}\right) - \mathcal{S}v_{s}^{2}\left(\frac{\omega_{w}^{2}-\omega_{w}^{2}}{2\Delta s^{2}}\right) + \frac{\mathcal{S}v_{r}^{2}\omega_{p}^{2}}{r_{p}^{2}} + \frac{\mathcal{S}v_{r}^{2}\omega_{p}^{2}}{r_{p}^{2}}$$

now re-arranging for \mathcal{W}_{p}^{r+1} since we will be computing it's values from an we get, initial state \mathcal{W}_{p}^{r} .

$$\begin{split} \omega_{p}^{n+1} &= \frac{\mathcal{U}\Delta t}{\mathcal{P}\Delta r^{2}}(\omega_{w}^{2} - 2\omega_{p}^{2} + \omega_{E}^{n}) + \frac{\mathcal{U}\Delta t}{\mathcal{P}\Delta z^{2}}(\omega_{s}^{2} - 2\omega_{p}^{2} + \omega_{N}^{n}) - \frac{\mathcal{U}\Delta t}{\mathcal{P}r_{p}^{2}} \\ &+ \frac{\mathcal{U}\Delta t}{2\mathcal{P}r_{p}\Delta r}(\omega_{E}^{2} - \omega_{w}^{2}) - \frac{\mathcal{V}r\Delta t}{2\Delta r}(\omega_{E}^{2} - \omega_{w}^{2}) - \frac{\mathcal{V}r\Delta t}{2\Delta z}(\omega_{N}^{2} - \omega_{s}^{2}) \\ &+ \frac{\mathcal{V}r\omega_{p}^{2}\Delta t}{r_{p}} + \omega_{p} \end{split}$$

expanding and grouping like terms, assuming M, &=continue we get,

$$\begin{split} \omega_{p}^{n+1} &= \frac{\mu_{\Delta t}}{P_{\Delta r^{2}}} - \frac{\mu_{\Delta t}}{z_{P}r_{p}\Delta r} + \frac{\nu_{r_{p}}^{r_{\Delta}}}{v_{\Delta r}} \omega_{w}^{n} + \left(\frac{\mu_{\Delta t}}{P_{\Delta 3}^{2}} + \frac{\nu_{3p}^{r_{\Delta}}}{z_{\Delta 3}}\right) \omega_{s}^{n} \\ &- \left(\frac{2\mu_{\Delta t}}{P_{\Delta r^{2}}} + \frac{2\mu_{\Delta t}}{P_{\Delta 3}^{2}} + \frac{\mu_{\Delta t}}{P_{r_{p}}^{2}} - \frac{\nu_{r_{p}}^{r_{p}}\Delta t}{r_{p}} - 1\right) \omega_{p}^{n} \\ &+ \left(\frac{\mu_{\Delta t}}{P_{\Delta r^{2}}} + \frac{\mu_{\Delta t}}{z_{P}r_{p}\Delta r} - \frac{\nu_{r_{p}}^{r_{p}}\Delta t}}{z_{P}r_{p}\Delta r}\right) \omega_{E}^{n} + \left(\frac{\mu_{\Delta t}}{P_{\Delta 3}^{2}} - \frac{\nu_{3p}^{r_{p}}\Delta t}}{z_{\Delta 3}}\right) \omega_{N}^{n} \end{split}$$

Q1a contid.

For ease of computational setup, let have the above as,

$$\left[\omega_{p}^{n+1} = a_{w}^{n} \omega_{w}^{n} + a_{s}^{n} \omega_{s}^{n} - a_{p}^{n} \omega_{p}^{n} + a_{e}^{n} \omega_{e}^{n} + a_{w}^{n} \omega_{n}^{n} \right] - (2.2)$$

This can be solved directly from given initial wop

-> Now we discretize the poisson radial velocity eqn in similar way,

$$\frac{\partial^{2} V_{r}}{\partial r^{2}} + \frac{\partial^{2} V_{r}}{\partial z^{2}} = \frac{\partial W}{\partial z} + \frac{V_{r}}{V^{2}} - \frac{1}{7} \frac{\partial V_{r}}{\partial r} - \frac{(2.3)}{2}$$

we can apply ego (1.1) e(1.2) to discretize as,

$$\frac{\sqrt{r_{W}} - 2\sqrt{r_{p}} + \sqrt{r_{E}} + \sqrt{r_{g}} - 2\sqrt{r_{p}} + \sqrt{r_{N}} = \sqrt{r_{N}} + \sqrt{r_{p}}}{43^{2}} = \frac{\sqrt{r_{p}} + \sqrt{r_{p}}}{243} + \frac{\sqrt{r_{p}}}{\sqrt{r_{p}}}$$

multiplying by rp2 and re-arranging we ge,

$$\left(\frac{v_{p}^{2}}{\Delta r^{2}} - \frac{v_{p}}{z_{\Delta r}}\right) v_{rw}^{n+1} + \frac{v_{p}^{2}}{\Delta z^{2}} v_{rs}^{n+1} - \left(\frac{z_{r}^{2}}{\Delta r^{2}} + \frac{z_{r}^{2}}{\Delta z^{2}} + 1\right) v_{rp}^{n+1} + \left(\frac{v_{p}^{2}}{\Delta r^{2}} + \frac{v_{p}}{z_{\Delta r}}\right) v_{re} + \left(\frac{v_{p}^{2}}{\Delta r^{2}} + \frac{v_{p}^{2}}{z_{\Delta r}}\right) v$$

Assuming that won't from ego (2.2), we can re-write the above as, (for ease of computational setup).

Q1a/cont'd.;

we will use preconditioned Bi-GGSTAB to some the above pentadiagonal linear system for Ur!

-> now lets discretize the continuity eqp, applying the product rule we have,

using first order backward difference for $\frac{\partial V_2}{\partial 3}$ and eyn (1.2) for $\frac{\partial V_r}{\partial r}$, we get,

$$\frac{\sqrt{3p-\sqrt{38}}}{\sqrt{3}} = -\frac{\sqrt{rp}}{r_p} - \left(\frac{\sqrt{n+1}}{\sqrt{re}} - \sqrt{rw}\right)$$

re-arranging for Jap, since vas can be known from our B.c., we will have,

for ease of computational setup, we re-write the above

$$V_{3P}^{n+1} = V_{3}^{n+1} - a_{p}V_{rp}^{n+1} - a_{p}V_{re}^{n+1} + a_{p}V_{rw}^{n+1}$$
 (2.5)

This can be solved directly for the axial velocity Jzp

Q1a] contd.]

>> Next, in order to apply the Boundary conditions (B.cs), lets Consider the four bound aries of our 2-D modell. as, 3 (out flow)
I fully Developed flow

we have the vorticity for assisymmetric cylindric coordinate as, Symmetry

$$\omega = \frac{\partial v_r}{\partial 3} - \frac{\partial v_z}{\partial r}$$

using first order one-sided difference 0,0 injet condition (inflow)

for our, and a second-order centered difference for our,

the above can be discretized as,

$$\omega_{p}^{n+1} = \frac{v_{rN}^{n+1} - v_{rp}^{n+1}}{\Delta 3} - \left(\frac{v_{3E}^{n+1} - v_{3VV}^{n+1}}{2\Delta r}\right) - (2.6)$$

no-sup boundary (wall)

and we have the half parabola egs as, x CLVZT

So that for:

Inflow;

re-written as Vap= kg (for computation and ease

using eqo(2.6), we have,

$$\omega_{p} = \frac{v_{rN}}{\Delta 3} - \left(\frac{v_{3E} - v_{3W}}{2\Delta T}\right)$$

re-written, wp=kz (for computational ease

Q1a] cont'd.!

$$V_{rp} = 0$$

$$\omega_p = \frac{\sqrt{3}w}{2\Delta r}$$

Symmetry;

since
$$\frac{\partial v_3}{\partial 3} = 0$$
, discretizing gives,

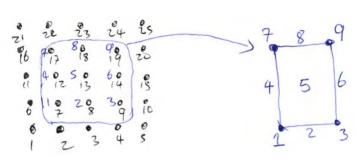
outflow;

we are give
$$\frac{\partial \omega}{\partial 3} = \frac{\partial v_r}{\partial 3} = \frac{\partial v_s}{\partial 3} = 0$$

and discretizing gives,

Q1a cont'di

>> Now because of the unique nature for solving our ur using a linear solver, lets assume we have a sample of grid mesh,



so that considering these 9 regions we will have the following boundary for Ur,

@ 1 The bottom left region:

applying both the inflow and symmetry B.Cs to eq. (2.4)

@ ITEL The bottom region:

we apply only the inflow conditions to ego (2.4),

@ 3 To 1 (aw your - apyp + a E V rE + an V rN = b)

(a) (3) The bottom right region: applying both the inflow and wall B.Cs. to equ (2.4),

we will apply only the symmetry condition to equ(2.4)

| 0 | 1 | a | cont'de! |
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@ 5 The middle region: Here we just use eqn (2.4) directly

(a) The right region:

we will only apply the wall condition to eqn(2.4)

[awdrw+asdrs-apdrp+andrn=b]

@ I The top left region applying both the outflow and Symmetry B.Cs to equ (2-4)

as vrs - apurp + a evre + an vrp = b

simply fing gives, as vrs - (ap + aw) vrp + a evre = b

The top region:

we will only apply the outflow condition to equ(z.4)

aw Vrw + as vrs - ap Urp + a E Ur = + an Urp = b simplyfying gives,

awvrw + as us - (ap-an) urp + as ure = b

@ 19 The top right region:

applying both the outflow and wall B.Cs. to egn(z.4)
awure + as urs - apurp + apurp = b
Simplyfying gives,

aw Vre + as Urs - (ap-an) Urp = b

Q1b] using the boxed equs from Q1a, we can write our program in fortran 90.

And we have our numeric integrated mass flow rate as,

>>> First I ran the Simulation with Tol=1e-5, and it seemed to be taking approximatly more than Thous and more than 5 million iterations, yet the tolerance was mit small enough.

>> So I reduced the tolerance to Tol= , which suffuciently small enough for a balance in both computational cost and accuracy.

>> Also I tried adaptive time step size (by incrementing dt by the at each new iteration). This worked well at lesser number of iterations than the Constant time step, only if the incrementing time (t) is Small enough, which in my case, I choose t=1e-7 for dt=dt+t

Otherwise the solution will explode if t is not small enough.

```
!**********Begin Header***********************
!This program was written by Godswill Ezeorah, Student Number: 501012886 on July 20, 2021.
!This program solves an unsteady linear/non-linear equation using finite difference method
!and was written as a solution to AE8112 PS5 q1(b,c,d)
program unsteady_Dfinite
   implicit none
   !Variable declaration
   DOUBLE PRECISION, dimension(:), ALLOCATABLE :: omg_i, V_r, V_z, r_i, z_i
   DOUBLE PRECISION, PARAMETER :: rh=8.3D-4, RR=1, L=10, mu=1.4D-4, CLVZ=5
   DOUBLE PRECISION, PARAMETER :: Pi = 4*atan(1.0)
   DOUBLE PRECISION :: dr, dz, dt, tols, rrr, start, finish
   INTEGER :: nr, nz, jj, i1, j1, l1, nnr, nnz
   call cpu_time(start) !gets the start time, for timing purpose
   open(2, file = 'PS5_Q1w.txt', status = 'unknown')
   open(3, file = 'PS5_Q1vr.txt', status = 'unknown')
   open(4, file = 'PS5_Q1vz.txt', status = 'unknown')
   Nr=21
          !Number of grid points along radial direction
   Nz=201 !Number of grid points along axial direction
   nnr=nr-2; nnz=nz-2 !Number of grid points in the interior region
   dr=RR/(nr-1)
   dz=L/(nz-1)
   dt=1D-6
              !given timestep
   tols=1D-5
   ALLOCATE(omg_i(nr*nz), V_r(nr*nz), V_z(nr*nz), r_i(nr), z_i(nz))
   !Initialization
   omg_i=0; V_r=0; V_z=0; r_i=0; z_i=0 !first initial consideration
   !rrr=0.75; omg_i=2*CLVZ*(rrr/RR**2); V_r=0; V_z=CLVZ*(1-(rrr/1)**2) !second initial consideration
   call unsteady(omg_i, V_r, V_z, r_i, z_i) !solves the unsteady system
   !Printing and writing results
   j1=nr*nz-(nr-1)
   l1=nr*nz
   jj=nz
   do i1 = j1, 1, -nr
       write(2,*) z_i(jj), omg_i(i1:l1)
       write(3,*) z_i(jj), V_r(i1:l1)
       write(4,*) z_i(jj), V_z(i1:l1)
       11=11-nr
       jj=jj-1
   end do
   call cpu_time(finish) !gets the end time
   write(2,*) r_i(1:nr); write(3,*) r_i(1:nr); write(4,*) r_i(1:nr)
```

```
print *, "Computation Time = ", (finish-start)/(60*60), "hours"
   close(2);close(3);close(4)
   contains
subroutine unsteady(oma, Vr, Vz, ri, zi)
       DOUBLE PRECISION, dimension(nr*nz) :: oma, Vz, omo, b, Vr
       DOUBLE PRECISION, dimension(nnr*nnz):: d, e, f, g, h, bi, Vri
       DOUBLE PRECISION, dimension(nr) :: ri
       DOUBLE PRECISION, dimension(nz) :: zi, mdot, som
       DOUBLE PRECISION :: aa,aw,as,ap,ae,an, k1, k2, k3, t, rp, zp
       DOUBLE PRECISION :: Mmdot, Amdot
       integer :: i, j, k, nj, nk, ll, ij, iii
   !!This subroutine solves unsteady fluid flow through a pipe problem using vorticity method
       open(1, file = 'PS5_Q1.txt', status = 'unknown')
       !Variable initialization
       d=0; e=0; f=0; g=0; h=0
       t=1D-8; rp=0; zp=0
       b=0
       Vri=0
       Mmdot=1
       ij=1
       do i = 1,nz !Loop for creating the axial and radial gridpoint dimensions
           if ( i<=nr ) then</pre>
              ri(i)=rp
               zi(i)=zp
           else
               zi(i)=zp
           end if
           rp=rp+dr
           zp=zp+dz
       do while (Mmdot>tols) !Main loop for time step
           j=nr+1
           k=nr+nr
           11=1
           omo=oma
           !The below loop solves the vorticity at each gridpoint
           do i = 1, nr*nz
               aw=mu*dt/(rh*dr**2)-mu*dt/(2*rh*ri(11)*dr)+Vr(i)*dt/(2*dr)
               as=mu*dt/(rh*dz**2)+Vz(i)*dt/(2*dz)
               ap=2*mu*dt/(rh*dr**2)+2*mu*dt/(rh*dz**2)+mu*dt/(rh*ri(11)**2)-Vr(i)*dt/ri(11) -1
               ae=mu*dt/(rh*dr**2)+mu*dt/(2*rh*ri(11)*dr)-Vr(i)*dt/(2*dr)
               an=mu*dt/(rh*dz**2)-Vz(i)*dt/(2*dz)
```

```
k2=(Vr(i+nr)/dz)-((Vz(i+1)-Vz(i-1))/(2*dr))
    k3=Vz(i-1)/(2*dr)
    !The 9 if-statements below are for the 9 regions of our CV formulation
    if ( i==1 ) then !bottom left region
        oma(i)=0 !symmetry condition
        11=11+1
    else if ( i < nr ) then
                                  !bottom region
       oma(i)=k2
       11=11+1
   else if ( i==nr ) then !bottom right region
        oma(i)=k3
        11=1
    else if ( i==nr*nz-(nr-1) ) then !top left region
        oma(i)=0 !symmetry condition
        11=11+1
    else if ( i==nr*nz ) then
                                   !top right region
       oma(i)=k3 !wall condition
    else if ( i==j ) then
                                 !left region
        oma(i)=0 !symmetry condition
        j=j+nr
       11=11+1
    else if ( i==k ) then
                                 !right region
       oma(i)=k3 !wall condition
        k=k+nr
        11=1
   else if ( i < (nr*nz)-nr ) then !Interior region</pre>
        oma(i)=aw*omo(i-1)+as*omo(i-nr)-ap*omo(i)+ae*omo(i+1)+an*omo(i+nr)
        11=11+1
   else
                                 !top region
       oma(i)=oma(i-nr) !outflow condition
        11=11+1
   end if
end do
!print *,'Oma=', oma(1:nr*nz)
!The below Loop composes the radial velocity at the boundary
j=nr+1
k=nr+nr
11=1
do i = 1, nr*nz
    !The 9 if-statements below are for the 9 regions of our formulation
    if ( i==1 ) then !bottom left region
       Vr(i)=0 !inflow or symmetry condition
        b(i)=(ri(11)**2)*(0-0)/(2*dz)
        11=11+1
   else if ( i < nr ) then
                                 !bottom region
```

```
Vr(i)=0 !inflow condition
        b(i)=(ri(11)**2)*(oma(i+nr)+((Vz(i+1)-Vz(i-1))/(2*dr)))/(2*dz)
        11=11+1
    else if ( i==nr ) then
                                   !bottom right region
        Vr(i)=0 !wall conditions
        b(i)=(ri(11)**2)*(oma(i+nr)-(Vz(i-1)/(2*dr)))/(2*dz)
        11=1
    else if ( i==nr*nz-(nr-1) ) then !top left region
        Vr(i)=0 !symmetry condition
        b(i)=ri(11)**2*(0-0)/(2*dz)
        11=11+1
    else if ( i==nr*nz ) then !top right region
        Vr(i)=0
        b(i)=(ri(11)**2)*(oma(i)-oma(i-nr-nr))/(2*dz)
    else if ( i==j ) then
                                  !left region
        Vr(i)=0 !symmetry conditions
        b(i)=ri(11)**2*(0-0)/(2*dz)
        j=j+nr
        11=11+1
    else if ( i==k ) then
                                  !right region
        Vr(i)=0 !wall conditions
        b(i)=(ri(11)**2)*(oma(i+nr)-oma(i-nr))/(2*dz)
        k=k+nr
        11=1
    else if ( i < (nr*nz)-nr ) then !Interior region</pre>
        b(i)=(ri(11)**2)*(oma(i+nr)-oma(i-nr))/(2*dz)
        11=11+1
    else
                                  !top region
        b(i)=(ri(11)**2)*(oma(i)-oma(i-nr-nr))/(2*dz)
    end if
end do
nj=nnr+1; nk=nnr+nnr
j=nr+1; k=1
11=2
do i = 1,nnr*nnz !This loop composes the interior radial velocity pentadiagonal vectors
    aw=(ri(11)**2)/(dr**2) - ri(11)/(2*dr)
    aa=(ri(11)**2)/(dz**2)
    ap=(2*ri(11)**2)/(dr**2) + (2*ri(11)**2)/(dz**2) + 1
    ae=(ri(11)**2)/(dr**2) + ri(11)/(2*dr)
    !The 9 if-statements below are for the 9 regions of our formulation
    if ( i==1 ) then !bottom left region
        f(i)=-ap
        g(i+1)=ae
        h(i+nnr)=aa
        bi(i)=b(j+k)
```

```
k=k+1
    11=11+1
else if ( i < nnr ) then !bottom region</pre>
   e(i-1)=aw
   f(i)=-ap
   g(i+1)=ae
   h(i+nnr)=aa
   bi(i)=b(j+k)
   k=k+1
   11=11+1
else if ( i==nnr ) then !bottom right region
   e(i-1)=aw
   f(i)=-ap
   h(i+nnr)=aa
   bi(i)=b(j+k)
   j=j+nr
   k=1
   11=2
else if ( i==nnr*nnz-(nnr-1) ) then !top left region
   d(i-nnr)=aa
   f(i)=-ap+aa
   g(i+1)=ae
   bi(i)=b(j+k)
   k=k+1
   11=11+1
else if ( i==nnr*nnz ) then !top right region
   d(i-nnr)=aa
   e(i-1)=aw
   f(i)=-ap+aa
   bi(i)=b(j+k)
else if ( i==nj ) then    !left region
   d(i-nnr)=aa
   f(i)=-ap
   g(i+1)=ae
   h(i+nnr)=aa
   bi(i)=b(j+k)
   nj=nj+nnr
   k=k+1
   11=11+1
else if ( i==nk ) then
                             !right region
   d(i-nnr)=aa
   e(i-1)=aw
   f(i)=-ap
   h(i+nnr)=aa
   bi(i)=b(j+k)
   nk=nk+nnr
```

```
j=j+nr
         k=1
         11=2
     else if ( i < (nnr*nnz)-nnr ) then !Interior region</pre>
         d(i-nnr)=aa
         e(i-1)=aw
        f(i)=-ap
         g(i+1)=ae
         h(i+nnr)=aa
         bi(i)=b(j+k)
         k=k+1
        11=11+1
    else
                                   !top region
        d(i-nnr)=aa
         e(i-1)=aw
        f(i)=-ap+aa
         g(i+1)=ae
         bi(i)=b(j+k)
         k=k+1
        11=11+1
    end if
end do
!For computational efficiency, the, A penta-diagonal matrix is split to 5 vectors
call Bi_CGSTAB_P(d,e,f,g,h,bi,Vri) !Solves the linear system at each time-step
j=nr+1
k=nr+nr
11=1
do i = 1,nr*nz!This loop adds the Bi_CGSTAB solved interior to the boundary radial velocity
     if ( i>nr+1 .and. i/=j .and. i/=k) then
        Vr(i)=Vri(ll)
        11=11+1
    else if ( i==nr*nz-(nr-1) ) then !interior region
         11=nnr*nnz-(nnr-1)
    else if ( i==j ) then
                                !left region
         j=j+nr
    else if ( i==k ) then
                              !right region
         k=k+nr
    end if
 end do
 !print "(a5,25f10.7)", 'Vr=', Vr(1:nr*nz)
 !The below loop solves the axial velocity at each grid point
 j=nr+1
k=nr+nr
11=1
```

```
do i = 1, nr*nz
    aa=dz/(2*dr)
    ap=dz/ri(ll)
    k1=CLVZ-CLVZ*ri(ll)
    !The 9 if-statements below are for the 9 regions of our CV formulation
    if ( i==1 ) then !bottom left region
        Vz(i)=k1 !inflow condition
        11=11+1
    else if ( i < nr ) then</pre>
                                   !bottom region
        Vz(i)=k1
        11=11+1
    else if ( i==nr ) then
                                  !bottom right region
        Vz(i)=0
        11=1
    else if ( i==nr*nz-(nr-1) ) then !top left region
        Vz(i)=Vz(i-nr) !symmetry condition
        11=11+1
    else if ( i==nr*nz ) then
                                    !top right region
        Vz(i)=0 !wall conditions
    else if ( i==j ) then
                                 !left region
        Vz(i)=Vz(i-nr) !symmetry conditions
        j=j+nr
        11=11+1
    else if ( i==k ) then
                                  !right region
        Vz(i)=0 !wall conditions
        k=k+nr
        11=1
    else if ( i < nr*nz-nr ) then !Interior region</pre>
        Vz(i)=Vz(i-nr)-ap*Vr(i)-aa*Vr(i+1)+aa*Vr(i-1)
        11=11+1
                                  !top region
    else
        Vz(i)=Vz(i-nr) !outflow condition
        11=11+1
    end if
!print "(a5,25f8.5)", 'Vz=', Vz(1:nr*nz)
do j = 1, nz !this loop solves the mass flow-rate along each z-gridpoint
    mdot(j) = 0.d0
    do i = 1, nr
        k=k+1
        mdot(j) = mdot(j) + 2.d0*pi*rh*Vz(k)*ri(i)
    end do
end do
do i = 1,Nz
```

```
som(i)=abs(mdot(i)-mdot(1))
          end do
          !print "(a5,25f8.3)", 'som=', som(1:nz)
          !print "(a5,25f8.3)", 'mdot=', mdot(1:nz)
          Mmdot=maxval(som)/mdot(1)
          Amdot=(sum(som)/nz)/mdot(1)
          !for printing results
          print "(i9,E12.3,E12.3)",ij, Mmdot, Amdot
          if (ij == 1) then
             Mmdot=1
              iii=101
          elseif (ij <= 100) then
              write(1,*) ij, Mmdot, Amdot
          elseif (ij == iii) then
              print *, achar(27)//"[2]" !clears the console
          elseif (ij \geq 3D6) then !set the maximum number of iteration
             Mmdot=tols
          end if
          !dt=dt+t !Adaptive time-step
          ij=ij+1
      end do
      close(1)
   end subroutine unsteady
subroutine Bi_CGSTAB_P(d,e,f,g,h,b,x)
      implicit none
      DOUBLE PRECISION, DIMENSION(nnr*nnz), INTENT(OUT) :: x
      DOUBLE PRECISION, dimension(nnr*nnz), INTENT(IN):: d, e, f, g, h, b
      DOUBLE PRECISION, DIMENSION(nnr*nnz) ::d1,e1,g1,h1,d2,e2,g2,h2
      DOUBLE PRECISION, DIMENSION(nnr*nnz) :: p, r, r0, y, z, v, s, t, k, ks, kt
      DOUBLE PRECISION :: rho0, rho, w, alpha, beta, r_check, tol, nan
      INTEGER :: i
   !!This subroutine solves a penta-diagonal linear system using
   !!The Preconditioned Bi CGSTAB Algorithm by Van Der Vorst
      !Variable initialization
      d2=0; e2=0; g2=0; h2=0
      !since we are dealing with vectors, I have done the multiplication of two vectors,
      !using this pattern
      !this multiples each vectors with x, for A*x
      d1=d(1:nnr*nnz)*x(1:nnr*nnz);e1=e(1:nnr*nnz)*x(1:nnr*nnz);
```

```
g1=g(1:nnr*nnz)*x(1:nnr*nnz);h1=h(1:nnr*nnz)*x(1:nnr*nnz);
!This circularlly shifts the vectors to the appropriate position, before addition
d2(nnr+1:nnr*nnz)=d1(1:nnr*nnz-nnr);e2(2:nnr*nnz)=e1(1:nnr*nnz-1);
g2(1:nnr*nnz-1)=g1(2:nnr*nnz)
h2(1:nnr*nnz-nnr)=h1(nnr+1:nnr*nnz)
!Hence A*x = d2+e2+f(1:nnr*nnz)*x(1:nnr*nnz)+g2+h2
r0=b-(d2+e2+f(1:nnr*nnz)*x(1:nnr*nnz)+g2+h2)
r=r0
w=1; alpha=1; rho0=1; r_check=1
v=0; p=0; k=0
tol=10E-8
!For the inverse of K
do i = 1, nnr*nnz
    k(i)=1/f(i)
end do
i=0
!main algorithm loop
do while (r_check>tol)
    i=i+1
    rho=dot_product(r0,r)
    beta=(rho/rho0)*(alpha/w)
    rho0=dot_product(r0,r)
    p=r+beta*(p-w*v)
    y=k(1:nnr*nnz)*p(1:nnr*nnz)
    d1=d(1:nnr*nnz)*y(1:nnr*nnz);e1=e(1:nnr*nnz)*y(1:nnr*nnz);
     g1=g(1:nnr*nnz)*y(1:nnr*nnz);h1=h(1:nnr*nnz)*y(1:nnr*nnz);
    d2(nnr+1:nnr*nnz)=d1(1:nnr*nnz-nnr);e2(2:nnr*nnz)=e1(1:nnr*nnz-1);
     g2(1:nnr*nnz-1)=g1(2:nnr*nnz)
    h2(1:nnr*nnz-nnr)=h1(nnr+1:nnr*nnz)
    v=(d2+e2+f(1:nnr*nnz)*y(1:nnr*nnz)+g2+h2)
    alpha=rho/dot_product(r0,v)
    if (rho==0) then
        alpha=0
    end if
    s=r-alpha*v
    z=k(1:nnr*nnz)*s(1:nnr*nnz)
    d1=d(1:nnr*nnz)*z(1:nnr*nnz);e1=e(1:nnr*nnz)*z(1:nnr*nnz);
     g1=g(1:nnr*nnz)*z(1:nnr*nnz);h1=h(1:nnr*nnz)*z(1:nnr*nnz);
    d2(nnr+1:nnr*nnz)=d1(1:nnr*nnz-nnr);e2(2:nnr*nnz)=e1(1:nnr*nnz-1);
     g2(1:nnr*nnz-1)=g1(2:nnr*nnz)
    h2(1:nnr*nnz-nnr)=h1(nnr+1:nnr*nnz)
    t=(d2+e2+f(1:nnr*nnz)*z(1:nnr*nnz)+g2+h2)
    d1=d(1:nnr*nnz)*s(1:nnr*nnz);e1=e(1:nnr*nnz)*s(1:nnr*nnz);
     g1=g(1:nnr*nnz)*s(1:nnr*nnz);h1=h(1:nnr*nnz)*s(1:nnr*nnz);
    d2(nnr+1:nnr*nnz)=d1(1:nnr*nnz-nnr);e2(2:nnr*nnz)=e1(1:nnr*nnz-1);
     g2(1:nnr*nnz-1)=g1(2:nnr*nnz)
```

```
h2(1:nnr*nnz-nnr)=h1(nnr+1:nnr*nnz)
          ks=(d2+e2+f(1:nnr*nnz)*s(1:nnr*nnz)+g2+h2)
          d1=d(1:nnr*nnz)*t(1:nnr*nnz);e1=e(1:nnr*nnz)*t(1:nnr*nnz);
           g1=g(1:nnr*nnz)*t(1:nnr*nnz);h1=h(1:nnr*nnz)*t(1:nnr*nnz);
          d2(nnr+1:nnr*nnz)=d1(1:nnr*nnz-nnr);e2(2:nnr*nnz)=e1(1:nnr*nnz-1);
           g2(1:nnr*nnz-1)=g1(2:nnr*nnz)
          h2(1:nnr*nnz-nnr)=h1(nnr+1:nnr*nnz)
          kt=(d2+e2+f(1:nnr*nnz)*t(1:nnr*nnz)+g2+h2)
          nan=dot_product(kt,ks)
          w=nan/dot_product(kt,kt)
          if (nan==0) then
             w=0
          end if
          x=x+alpha*y+w*z
          r=s-w*t
          r_check=norm2(r)
   end subroutine Bi_CGSTAB_P
```

end program unsteady_Dfinite

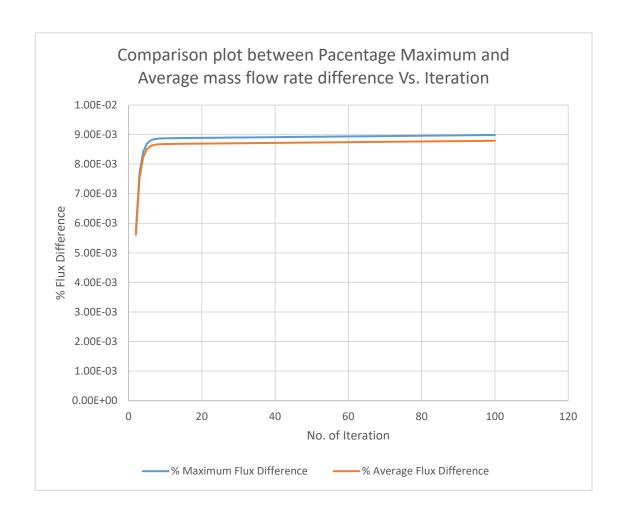
Q1b] contid!

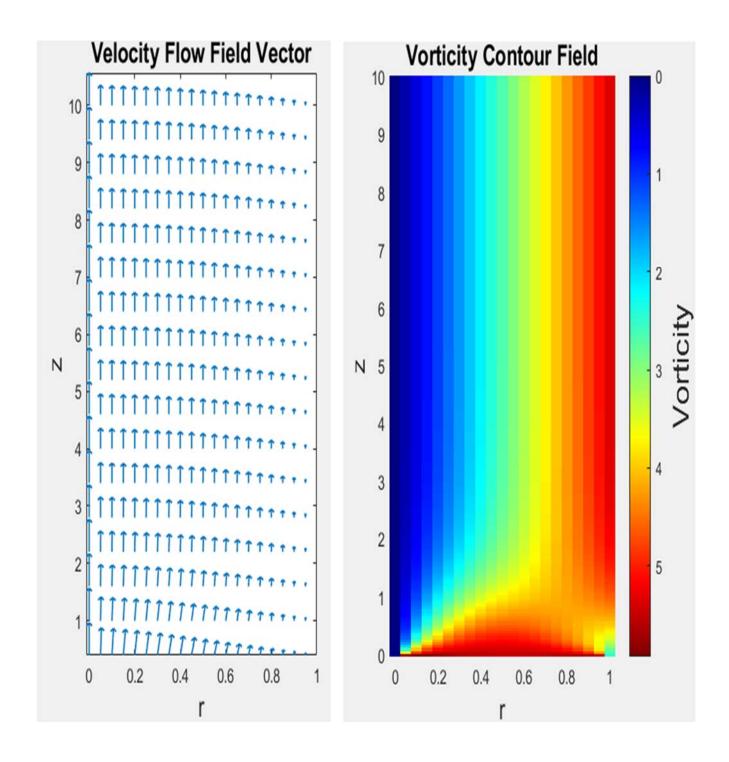
- >>> For computation time, I observed that slight adjustment to the toterance value greatly affects the computation time, For instance if adaptive timestep is not used. Even after 7 hours of timestep is not used. Even after 9 hours of simulation on a constant timstep (dt), the tot=1e-5 simulation on a constant timstep (dt), the tot=1e-5 is not still reached. But I was able to simulate for 3 million Iterations at just 3.088 hours using vector only Bi-CastaB instead of the conventional matrix type. This improved my computational time.

 Greatly.
- >> The accuracy which is a direct function of our Percenteage maximum/awerage flow rate difference (i.e based on the Conservation law), will be as accurate as we want it to be, in the expense of our computational time.

But what I've been able to deduce from our simulation, is that the solution gets more slightly inaccurate as we increase at new iterations up to about 50,000 iterations. And then it starts to get accurate agains and on-ward in expense of computation time.

- Q1C)>From below plot, we can see that both percentage Maximum and average flow rate difference nce have similar behavour against the number of iteration. And thus eather one of them can be used to theck for the accuracy of our solution. (This was taken from samples of just 100 iterations).
 - >> And setting our tolerance to tol = 8e-3 is small enough for us to understand the behavour of the solution (ie. after it begins to converge again, from about 50,000 iterations). With this we got about 3 million iterations. And the vector plot for the flow field, as while as the contour plot for the vorticity is shown below;





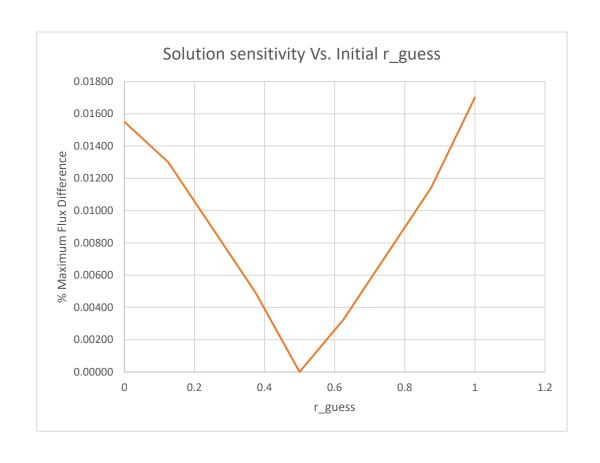
Q1d If we assume an initial guess of; $V_r = 0 \quad , \quad V_3 = CLV 2 \left(1 - \left(\frac{r_3 u_0^2}{R^2} \right) \right) \quad w = 2(CLV 2) \left(\frac{r_3 u_0^2}{R^2} \right)$ lets assume R to be our overall radius R=1cm and let's play with $V = 0 \le r \le 1$ cm (we can call this r-guess).

- >> The sensitivity of varying this r-guess is shown in the table below. Which at r-guess = 0.5 we can see the solution converging very at just 9 iteration.
- >> But the strange part is with the vector and contour ur plot which is shown below for reguess = 0.5 and it cooks much different from the solution if $V_r=0$, $V_3=0$ & $\omega=0$.

I also observed that rguess=0 is thields similar results to our previous guess of $v_r = v_z = \omega = 0$.

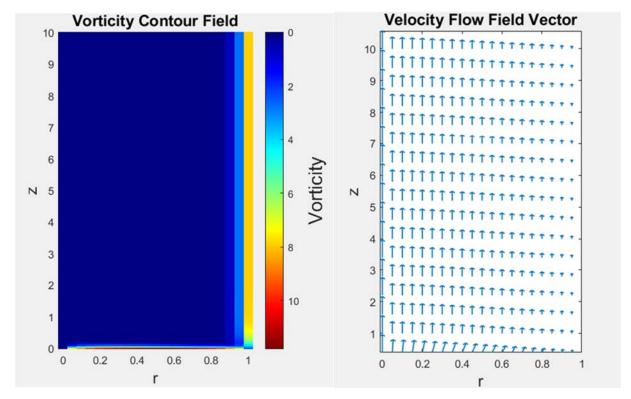
Q1d] cont'd.]

| No. of Iteration | r_guess | % Maximum Flux Difference |
|------------------|---------|---------------------------|
| 10000 | 1 | 0.01550 |
| 10000 | 0.125 | 0.01300 |
| 10000 | 0.25 | 0.00895 |
| 10000 | 0.375 | 0.00489 |
| 9 | 0.5 | 0.00001 |
| 10000 | 0.625 | 0.00324 |
| 10000 | 0.75 | 0.00732 |
| 10000 | 0.875 | 0.01140 |
| 10000 | 1 | 0.01700 |

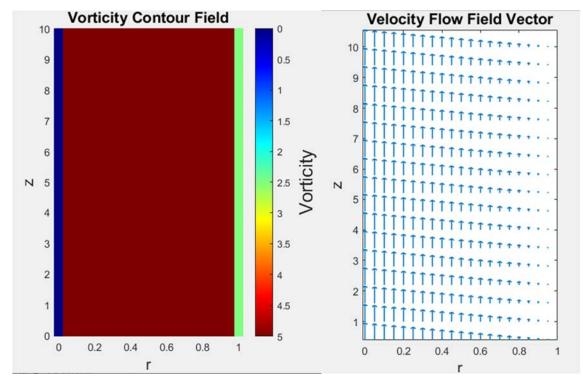


Q1d] cont'd.]

Plot when $r_{guess} = 0$:



Plot when $r_{guess} = 0.5$:



Q1d] cont'd.]

Plot when $r_{guess} = 1$:

