



UNIVERSITAT POLITÈCNICA DE CATALUNYA  
BARCELONATECH

Escola Superior d'Enginyeries Industrial,  
Aeroespacial i Audiovisual de Terrassa

# Advanced study for the numerical resolution of the Navier-Stokes equations. Application to thermal energy storage systems

## Document:

Annexes

## Author:

Oriol Sanmarti Perona

## Director / Co-director:

Carlos David Perez Segarra - Jordi Vera Fernández

## Degree:

Master's degree in Space and Aeronautical Engineering

## Examination session:

Spring, 2022

TREBALL DE FI D'ESTUDIS



## Contents

<b>1</b>	<b>4 Materials</b>	<b>2</b>
<b>2</b>	<b>Smith Hutton CDS</b>	<b>12</b>
<b>3</b>	<b>Smith Hutton UDS</b>	<b>16</b>
<b>4</b>	<b>Lid Driven Cavity</b>	<b>20</b>
<b>5</b>	<b>Differentially Heated Cavity</b>	<b>26</b>
<b>6</b>	<b>Square Cylinder</b>	<b>36</b>
<b>7</b>	<b>Square Cylinder, coefficients</b>	<b>46</b>
<b>8</b>	<b>Burgers' equation</b>	<b>51</b>
<b>9</b>	<b>Burgers' equation LES</b>	<b>54</b>
<b>10</b>	<b>Lid Driven Cavity 3D</b>	<b>57</b>
<b>11</b>	<b>Parallelization Monte Carlo method</b>	<b>69</b>
<b>12</b>	<b>Parallelization 1D bar</b>	<b>71</b>

This document presents the annexes of the study carried out. These annexes consist of the C++ codes that have been used to carry out the cases presented. As for Thermocline's study, the code is not presented due to privacy reasons, since a code of the CTTC research group has been used.

## 1 4 Materials

```
#include <iostream>
#include <fstream>
#include <math.h> // To use pow
#include <vector> // To use vectors
using namespace std;

const double q = 60; //W/m
const double cp1 = 750; //J/(KgK)
const double cp2 = 770;
const double cp3 = 810;
const double cp4 = 930;
const double densitat1 = 1500; //Kg/m^3
const double densitat2 = 1600;
const double densitat3 = 1900;
const double densitat4 = 2500;
const double lambda1 = 170; //W/(mK)
const double lambda2 = 140;
const double lambda3 = 200;
const double lambda4 = 140;
const double lambda42 = 2/((1/lambda2)+(1/lambda4));
const double lambda13 = 2/((1/lambda1)+(1/lambda3));
const double lambda34 = 2/((1/lambda3)+(1/lambda4));
const double lambda32 = 2/((1/lambda3)+(1/lambda2));
const double lambda12 = 2/((1/lambda1)+(1/lambda2));
const double L = 1.1;
const double H = 0.8;
const int Nx = 55; //NODES x
const int Ny = 40; //NODES Y
const double epsilon = 0.000001; //ERROR
const double inct = 1; //CAMBIAR INCREMENT DE TEMPS
const double tempsmaxim = 5000; //CAMBIAR TEMPS
const double temps = tempsmaxim/inct; //NO CAMBIAR
const double incx = L/(Nx);
const double incy = H/(Ny);
const double gamma1 = (densitat1*cp1*incx*incy)/(inct); //450
const double gamma2 = (densitat2*cp2*incx*incy)/(inct); //492.8
const double gamma3 = (densitat3*cp3*incx*incy)/(inct); //615.6
const double gamma4 = (densitat4*cp4*incx*incy)/(inct); //930
const double alpha = 9; //W/m^2K
const double Tf = 33; //C
const double archius = 500;
const int iteprint = tempsmaxim/archius;
```

```

double CalculTiM2(double TE, double TW, double TN, double TS, double Tp){
    double r;
    double ae = (lambda2*incy)/(gamma2*incx);
    double aw = (lambda2*incy)/(gamma2*incx);
    double an = (lambda2*incx)/(gamma2*incy);
    double as = (lambda2*incx)/(gamma2*incy);
    double b = Tp;
    double ap = 1 + ae + aw + an + as;
    r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
    return r;
}

double CalculTiM1(double TE, double TW, double TN, double TS, double Tp){
    double r;
    double ae = (lambda1*incy)/(gamma1*incx);
    double aw = (lambda1*incy)/(gamma1*incx);
    double an = (lambda1*incx)/(gamma1*incy);
    double as = (lambda1*incx)/(gamma1*incy);
    double b = Tp;
    double ap = 1 + ae + aw + an + as;
    r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
    return r;
}

double CalculTiM3(double TE, double TW, double TN, double TS, double Tp){
    double r;
    double ae = (lambda3*incy)/(gamma3*incx);
    double aw = (lambda3*incy)/(gamma3*incx);
    double an = (lambda3*incx)/(gamma3*incy);
    double as = (lambda3*incx)/(gamma3*incy);
    double b = Tp;
    double ap = 1 + ae + aw + an + as;
    r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
    return r;
}

double CalculTiM4(double TE, double TW, double TN, double TS, double Tp){
    double r;
    double ae = (lambda4*incy)/(gamma4*incx);
    double aw = (lambda4*incy)/(gamma4*incx);
    double an = (lambda4*incx)/(gamma4*incy);
    double as = (lambda4*incx)/(gamma4*incy);
    double b = Tp;
    double ap = 1 + ae + aw + an + as;
    r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
    return r;
}

double CalculTiM24(double TE, double TW, double TN, double TS, double Tp){
    double r;
    double ae = (lambda2*incy)/(((gamma2))*incx);
    double aw = (lambda2*incy)/(((gamma2))*incx);

```

```

double an = (lambda42*incx)/(((gamma2))*incx);
double as = (lambda2*incx)/(((gamma2))*incx);
double b = Tp;
double ap = 1 + ae + aw + an + as;
r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
return r;
}
double CalculTiM42(double TE, double TW, double TN, double TS, double Tp){
double r;
double ae = (lambda4*incx)/(((gamma4))*incx);
double aw = (lambda4*incx)/(((gamma4))*incx);
double an = (lambda4*incx)/(((gamma4))*incx);
double as = (lambda42*incx)/(((gamma4))*incx);
double b = Tp;
double ap = 1 + ae + aw + an + as;
r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
return r;
}
double CalculTiM13(double TE, double TW, double TN, double TS, double Tp){
double r;
double ae = (lambda1*incx)/(((gamma1))*incx);
double aw = (lambda1*incx)/(((gamma1))*incx);
double an = (lambda13*incx)/(((gamma1))*incx);
double as = (lambda1*incx)/(((gamma1))*incx);
double b = Tp;
double ap = 1 + ae + aw + an + as;
r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
return r;
}
double CalculTiM31(double TE, double TW, double TN, double TS, double Tp){
double r;
double ae = (lambda3*incx)/(((gamma3))*incx);
double aw = (lambda3*incx)/(((gamma3))*incx);
double an = (lambda3*incx)/(((gamma3))*incx);
double as = (lambda13*incx)/(((gamma3))*incx);
double b = Tp;
double ap = 1 + ae + aw + an + as;
r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
return r;
}
double CalculTiM34(double TE, double TW, double TN, double TS, double Tp){
double r;
double ae = (lambda34*incx)/(((gamma3))*incx);
double aw = (lambda3*incx)/(((gamma3))*incx);
double an = (lambda3*incx)/(((gamma3))*incx);
double as = (lambda3*incx)/(((gamma3))*incx);
double b = Tp;
double ap = 1 + ae + aw + an + as;
r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);

```

```

    return r;
}
double CalculTiM43(double TE, double TW, double TN, double TS, double Tp){
    double r;
    double ae = (lambda4*incy)/(((gamma4)*incx);
    double aw = (lambda34*incy)/(((gamma4)*incx);
    double an = (lambda4*incx)/(((gamma4)*incy);
    double as = (lambda4*incx)/(((gamma4)*incy);
    double b = Tp;
    double ap = 1 + ae + aw + an + as;
    r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
    return r;
}
double CalculTiM32(double TE, double TW, double TN, double TS, double Tp){
    double r;
    double ae = (lambda32*incy)/(((gamma3)*incx);
    double aw = (lambda3*incy)/(((gamma3)*incx);
    double an = (lambda3*incx)/(((gamma3)*incy);
    double as = (lambda3*incx)/(((gamma3)*incy);
    double b = Tp;
    double ap = 1 + ae + aw + an + as;
    r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
    return r;
}
double CalculTiM23(double TE, double TW, double TN, double TS, double Tp){
    double r;
    double ae = (lambda2*incy)/(((gamma2)*incx);
    double aw = (lambda32*incy)/(((gamma2)*incx);
    double an = (lambda2*incx)/(((gamma2)*incy);
    double as = (lambda2*incx)/(((gamma2)*incy);
    double b = Tp;
    double ap = 1 + ae + aw + an + as;
    r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
    return r;
}
double CalculTiM12(double TE, double TW, double TN, double TS, double Tp){
    double r;
    double ae = (lambda12*incy)/(((gamma1)*incx);
    double aw = (lambda1*incy)/(((gamma1)*incx);
    double an = (lambda1*incx)/(((gamma1)*incy);
    double as = (lambda1*incx)/(((gamma1)*incy);
    double b = Tp;
    double ap = 1 + ae + aw + an + as;
    r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
    return r;
}
double CalculTiM21(double TE, double TW, double TN, double TS, double Tp){
    double r;
    double ae = (lambda2*incy)/(((gamma2)*incx);

```

```

double aw = (lambda12*incy)/(((gamma2))*incy);
double an = (lambda2*incy)/(((gamma2))*incy);
double as = (lambda2*incy)/(((gamma2))*incy);
double b = Tp;
double ap = 1 + ae + aw + an + as;
r = ((ae*TE+aw*TW+an*TN+as*TS)+b)/(ap);
return r;
}
double CalculNeumannesquerra(double TE, double TW, double TS, double Tp){
double r;
double ae = (lambda3*incy)/(gamma3*incy);
double aw = (lambda3*incy)/(gamma3*incy);
double as = (lambda3*incy)/(gamma3*incy);
double b = Tp;
double ap = 1 + ae + aw + as;
r = ((ae*TE+aw*TW+((q*incy)/gamma3)+as*TS)+b)/(ap);
return r;
}
double CalculNeumanndreta(double TE, double TW, double TS, double Tp){
double r;
double ae = (lambda4*incy)/(gamma4*incy);
double aw = (lambda4*incy)/(gamma4*incy);
double as = (lambda4*incy)/(gamma4*incy);
double b = Tp;
double ap = 1 + ae + aw + as;
r = ((ae*TE+aw*TW+((q*incy)/gamma4)+as*TS)+b)/(ap);
return r;
}
double CalculConvectionabaix(double TN, double TE, double TS, double Tp){
double r;
double an = (lambda1*incy)/(gamma1*incy);
double ae = (lambda1*incy)/(gamma1*incy);
double as = (lambda1*incy)/(gamma1*incy);
double b = Tp + ((Tf*alpha*incy)/gamma1);
double ap = 1 + ae + as + an + ((alpha*incy)/gamma1);
r = ((ae*TE+an*TN+as*TS)+b)/(ap);
return r;
}
double CalculConvectionadalt(double TN, double TE, double TS, double Tp){
double r;
double an = (lambda3*incy)/(gamma3*incy);
double ae = (lambda3*incy)/(gamma3*incy);
double as = (lambda3*incy)/(gamma3*incy);
double b = Tp + ((Tf*alpha*incy)/gamma3);
double ap = 1 + ae + as + an + ((alpha*incy)/gamma3);
r = ((ae*TE+an*TN+as*TS)+b)/(ap);
return r;
}
double CalculAdaltEsquerra(double TE, double TS, double Tp){

```

```

double r;
double ae = (lambda3*incx)/(gamma3*incx);
double as = (lambda3*incx)/(gamma3*incx);
double aw = (alpha*incx)/gamma3;
double b = Tp + Tf*aw + (q*incx)/gamma3;
double ap = 1 + ae + as + aw;
r = ((ae*TE+as*TS)+b)/(ap);
return r;
}
double CalculAbaixDreta(double TW, double TN, double Tp){
double r;
double aw = (lambda2*incx)/(gamma2*incx);
double ae = (lambda2*incx)/(gamma2*(incx/2));
double an = (lambda2*incx)/(gamma2*incx);
double as = (lambda2*(incx/2))/(gamma2*(incx/2));
double b = Tp;
double ap = 1 + aw + an + as +ae;
r = ((aw*TW+an*TN+as*23+ae*23)+b)/(ap);
return r;
}
double CalculAbaixEsquerra( double TN, double Tp){
double r;
double ae = (lambda1*incx)/(gamma1*incx);
double an = (lambda1*incx)/(gamma1*incx);
double as = (lambda1*(incx))/(gamma1*(incx/2));
double b = Tp + (alpha*incx*Tf)/gamma1;
double ap = 1 + ae + an + as + (alpha*incx)/gamma1;
r = ((ae*23 + an*TN + as*23 )+b)/(ap);
return r;
}
double CalculAdaltDreta(double TE, double TS, double TW, double Tp){
double r;
double ae = (lambda3*incx)/(gamma3*(incx/2));
double as = (lambda3*incx)/(gamma3*incx);
double aw = (lambda3*incx)/(gamma3*incx);
double b = Tp + (q*incx)/gamma3;
double ap = 1 + ae + as + aw + (alpha*(incx))/gamma3;
r = ((ae*TE+as*TS+aw*TW)+b)/(ap);
return r;
}
int main(){
double T[Nx][Ny];
double Tant[Nx][Ny];
double Tguarda[Nx][Ny];
double x = 0;
double y = 0;

```



```

for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    T[i][j] = 8;
    Tguarda[i][j]=20;
    Tant[i][j] = 8;
    if(j==Ny-1){
      T[i][j] = 23;
    }
  }
}

ofstream Archivo;
char nombre[500];
Archivo.open("Temp.0.txt");
Archivo << "t" << "t" << "T" << "t" << "x" << "t" << "y" << "t" << "z" << "\n";
for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    x = incx*i + incx/2;
    y = incy*j + incy/2;
    Archivo << 0 << "t" << T[i][j] << "t" << x << "t" << y << "t" << 0 << "\n";
  }
}

Archivo.close();
ofstream Archivo2;
Archivo2.open("Apartat.b.txt");
Archivo2 << "t" << "t" << "T(0.65,0.56)" << "t" << "T(0.74,0.72)" << "\n";
Archivo2 << 0 << "t" << 8 << "t" << 8 << "\n";

for(int k = 1; k<=temps; k++){
  bool itefin = false;
  while(!itefin){
    itefin = true;
    for(int j = 0; j<Ny; j++){
      for(int i = 0; i<Nx; i++){
        x = incx*i + incx/2;
        y = incy*j + incy/2;
        //PARET DRETA:
        if(x==L-(incx/2) and y < H - (incy/2) and y > incy/2){
          Tguarda[i][j] = T[i][j];
          T[i][j] = 8 + (0.005 * (inct * k));
        }
        //PARET ABAIX:
        if(y == H - (incy/2) and x < L-(incx/2) and x > incx/2){
          Tguarda[i][j] = T[i][j];
          T[i][j] = 23;
        }
      }
    }
  }
  //PARET ADALT Neumann

```

```

if(y == incy/2 and x > incx/2 and x <= 0.5){
  Tguarda[i][j] = T[i][j];
  T[i][j] = CalculNeumannesquerra(T[i+1][j], T[i-1][j], T[i][j+1], Tant[i][j]);
}
if(y == incy/2 and x > 0.5 and x < L-(incx/2)){
  Tguarda[i][j] = T[i][j];
  T[i][j] = CalculNeumanndreta(T[i+1][j], T[i-1][j], T[i][j+1], Tant[i][j]);
}
//PARET ESQUERRA CONVECCIO

if(x == incx/2 and y >= 0.4 and y < H-(incy/2)){
  Tguarda[i][j] = T[i][j];
  T[i][j] = CalculConvectionabaix(T[i][j-1], T[i+1][j], T[i][j+1], Tant[i][j]);
}

if(x == incx/2 and y > incy/2 and y < 0.4){
  Tguarda[i][j] = T[i][j];
  T[i][j] = CalculConvectionadalt(T[i][j-1], T[i+1][j], T[i][j+1], Tant[i][j]);
}

//M2
if(x >= 0.5 and x < L-(incx/2) and y >= 0.1 and y < H-(incy/2)){
  //M21
  if(x-incx < 0.5 and y > 0.4){
    Tguarda[i][j] = T[i][j];
    T[i][j] = CalculTiM21(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
  }
  //M24
  else if (y-incy < 0.1){
    Tguarda[i][j] = T[i][j];
    T[i][j] = CalculTiM24(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
  }
  //M23
  else if(x-incx < 0.5 and y < 0.4){
    Tguarda[i][j] = T[i][j];
    T[i][j] = CalculTiM23(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
  }
  else {
    Tguarda[i][j] = T[i][j];
    T[i][j] = CalculTiM2(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
  }
}

//M1
if(x > incx/2 and x < 0.5 and y >= 0.4 and y < H-(incy/2)){
  //M13
  if(y-incy < 0.4){

```

```

Tguarda[i][j] = T[i][j];
T[i][j] = CalculTiM13(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
}
//M12
else if(x+incx > 0.5){
Tguarda[i][j] = T[i][j];
T[i][j] = CalculTiM12(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);

}
else{
Tguarda[i][j] = T[i][j];
T[i][j] = CalculTiM1(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
}
}
//M3
if(x > incx/2 and x < 0.5 and y > incy/2 and y < 0.4){
//M32
if(x+incx > 0.5 and y > 0.1 and y <= 0.4){
Tguarda[i][j] = T[i][j];
T[i][j] = CalculTiM32(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
}
//M31
else if(x > incx/2 and x < 0.5 and y+incy > 0.4){
Tguarda[i][j] = T[i][j];
T[i][j] = CalculTiM31(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
}
//M34
else if(x+incx > 0.5 and y > incy/2 and y < 0.1){
Tguarda[i][j] = T[i][j];
T[i][j] = CalculTiM34(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
}
else{
Tguarda[i][j] = T[i][j];
T[i][j] = CalculTiM3(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
}
}
//M4
if(x >= 0.5 and x < L-(incx/2) and y > incy/2 and y < 0.1){
//M43
if(x-incx < 0.5){
Tguarda[i][j] = T[i][j];
T[i][j] = CalculTiM43(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
}
//M42
else if(y+incy > 0.1 ){
Tguarda[i][j] = T[i][j];
T[i][j] = CalculTiM42(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
}
}

```

```

    else {
        Tguarda[i][j] = T[i][j];
        T[i][j] = CalculTiM4(T[i+1][j], T[i-1][j], T[i][j-1], T[i][j+1], Tant[i][j]);
    }
}
//NODE ADALT ESQUERRA

if(x == incx/2 and y == incy/2){
    Tguarda[i][j] = T[i][j];
    T[i][j] = CalculAdaltEsquerra(T[i+1][j],T[i][j+1], Tant[i][j]);
}
//NODE ABAIX DRETA
if(x == L-(incx/2) and y == H-(incy/2)){
    Tguarda[i][j] = T[i][j];
    T[i][j] = CalculAbaixDreta(T[i-1][j], T[i][j-1], Tant[i][j]);
}
//NODE ABAIX ESQUERRA
if(x == incx/2 and y == H-(incy/2)){
    Tguarda[i][j] = T[i][j];
    T[i][j] = CalculAbaixEsquerra( T[i][j-1], Tant[i][j]);
}
//NODE ADALT DRETA
if(x == L-(incx/2) and y == incy/2){
    Tguarda[i][j] = T[i][j];
    T[i][j] = CalculAdaltDreta(T[i][j+1], T[i][j+1], T[i-1][j], Tant[i][j]);
}
}
}
for(int j = 1; j<Ny-1; j++){
for(int i = 1; i<Nx-1; i++){
double error = 0;
error = abs(Tguarda[i][j] - T[i][j]);
if(error > epsilon) {itefin = false;
    //cout<<"la i es: " << i << " la j es : " <<j<<" La t guarda es: "<<Tguarda[i][j]<<" La T es : " <<T[i][j]<<" El temps es : " <<k*
    inct<<endl;
    //cout<<error<<endl;
}

//cout<<"L'error es: " << itefin<<endl;
}
}
//cout<<endl;
}
//cout<<inct*k<<endl;
int num = k;
if(k%iteprint == 0){

```

```

sprintf(nombre, "Temp-%05d.txt", num);
Archivo.open(nombre, ios::trunc);
Archivo << "t" << "\t" << "T" << "\t" << "x" << "\t" << "y" << "\t" << "z" << "\n";
}

for(int j = 0; j<Ny; j++){
for(int i = 0; i<Nx; i++){
  Tant[i][j] = T[i][j];
  if(k%iteprint == 0){

    x = incx*i + incx/2;
    y = incy*j + incy/2;
    Archivo << k*inct << "\t" << T[i][j] << "\t" << x << "\t" << y << "\t" << 0 << "\n";

  }
}
}
x = 0;
y = 0;
Archivo.close();

//APARTAT B
int punt1 = 0.64/incx;
int punt1prima = 0.66/incx;
int punt2 = 0.44/incy;
int punt3 = 0.74/incx;
int punt4 = 0.28/incy;
double Tpunt1 = ((T[32][12])+(T[32][11]))/2;
double Tpunt2 = ((T[37][4])+(T[36][4])+(T[37][3])+(T[36][3]))/4;
Archivo2 << k*inct << "\t" << Tpunt1 << "\t" << Tpunt2 << "\n";
}
Archivo2.close();
for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    cout<<T[i][j]<<" ";
  }
  cout<<endl;
}
}
}

```

## 2 Smith Hutton CDS

```

#include <iostream>
#include <fstream>
#include <math.h> // To use pow
#include <vector> // To use vectors
#include <ctime>

```

```

using namespace std;

double const L = 2;
double const H = 1;
int const Nx = 100;
int const Ny = 50;
double const inct = 0.001;
double const densitat = 1000000;
double const gamma = 1;
double const incx = L/Nx;
double const incy = H/Ny;
//double const temps = 200;
double const alpha = 10;
double const epsilon = 0.0001;

double Calculvelu(double x, double y){
    double u = 2*y*(1-pow(x,2));
    return u;
}
double Calculvelv(double x, double y){
    double v = -2*x*(1-pow(y,2));
    return v;
}
double Calculinlet(double PHI, double x){
    double r = 1 + tanh((2*x+1)*alpha);
    return r;
}

int main(){

    unsigned t0, t1;

    t0=clock();

    double PHI[Nx][Ny];
    double PHiant[Nx][Ny];
    double u[Nx][Ny];
    double v[Nx][Ny];
    double xi = -(L/2) + (incx/2);
    double yi = H - (incy/2);
    double x;
    double y;
    ofstream Archivo1;
    ofstream Archivo2;
    ofstream Archivo3;

    Archivo1.open("Mapa de velocitats.txt");
    Archivo1<<"x"<<"\t"<<"y"<<"\t"<<"z"<<"\t"<<"u"<<"\t"<<"v"<<"\n";

```

```

Archivo2.open("Mapa Propietat.txt");
Archivo2<<"x"<<"\t"<<"y"<<"\t"<<"z"<<"\t"<<"PHI"<<"\n";

Archivo3.open("Outlet propietat.txt");
Archivo3<<"x"<<"\t"<<"PHI"<<"\n";

//DECLAREM MAPA DE VELOCITATS
for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    x = xi + i*incx;
    y = yi - j*incy;
    u[i][j] = Calculvelu(x,y);
    v[i][j] = Calculvelv(x,y);
    PHI[i][j] = 0;
    PHIant[i][j] = PHI[i][j];
    //cout<<" ("<<Vu[i][j]<<" "<<Vv[i][j]<<" ") ";
    cout<<PHI[i][j]<<" ";
    Archivo1<<x<<"\t"<<y<<"\t"<<0<<"\t"<<u[i][j]<<"\t"<<v[i][j]<<"\n";
  }
  cout<<endl;
}
Archivo1.close();

// double tempsmaxim = temps/inct;

bool conv = false;
int k = 0;
while(!conv){

  conv = true;

  for(int j = 0; j<Ny; j++){
    for(int i = 0; i<Nx; i++){
      x = xi + i*incx;
      y = yi - j*incy;
      //CALCUL NODES CENTRALS
      if(x > xi and y < yi and x < -xi and j<Ny-1){
        PHIant[i][j] = PHI[i][j];
        double uw = u[i-1][j];
        double ue = u[i+1][j];
        double vn = v[i][j-1];
        double vs = v[i][j+1];
        double ae = (gamma*incy)/incx;
        double aw = (gamma*incy)/incx;
        double an = (gamma*incx)/incy;
        double as = (gamma*incx)/incy;
        double PHIp = PHI[i][j];

```

```

double PHIE = (PHI[i+1][j] + PHIp)/2;
double PHIw = (PHI[i-1][j] + PHIp)/2;
double PHIn = (PHI[i][j-1] + PHIp)/2;
double PHIs = (PHI[i][j+1] + PHIp)/2;
double PHIE2 = PHI[i+1][j];
double PHIW2 = PHI[i-1][j];
double PHIN2 = PHI[i][j-1];
double PHIS2 = PHI[i][j+1];
double conveccio = densitat*(PHIE*incy*ue - PHIw*incy*uw + PHIn*incx*vn - PHIs*incx*vs);
double diffusio = ae*PHIE2 + aw*PHIW2 + an*PHIN2 + as*PHIS2 - PHIp*(an+aw+as+ae);
PHI[i][j] = (inct/(incx*incy*densitat)) * (-conveccio + diffusio) + PHIp;

if(abs(PHIant[i][j]-PHI[i][j]) > epsilon){
  conv = false;
  //cout<<PHIant[i][j]<<"--"<<PHI[i][j]<<endl;
}
}
//INLET
else if(j == Ny-1 and x < 0 and x > xi){
  PHIant[i][j] = PHI[i][j];
  PHI[i][j] = Calculinlet(PHI[i][j], x);
}
//OUTLET
else if(j == Ny-1 and x > 0 and x < -xi){
  PHIant[i][j] = PHI[i][j];
  PHI[i][j] = PHI[i][j-1];
  if(abs(PHIant[i][j]-PHI[i][j]) > epsilon){
    conv = false;
    //cout<<PHIant[i][j]<<"--"<<PHI[i][j]<<endl;
  }
}
}
//REST OF WALLS
else if(abs(x-xi)<0.0001 or abs(x+xi)<0.0001 or abs(y-yi)<0.0001){
  PHIant[i][j] = PHI[i][j];
  PHI[i][j] = 1-tanh(10);
}
}
}

k++;
if(k < 15000){
  conv = false;
}
}

```



```

cout<<"-----" <<endl;
for(int j = 0; j<Ny; j++){
    for(int i = 0; i<Nx; i++){
        x = xi + i*incx;
        y = yi - j*incy;
        cout<<PHI[i][j]<<" ";
        Archivo2<<x<<"\t"<<y<<"\t"<<0<<"\t"<<PHI[i][j]<<"\n";
        if(j==Ny-1){
            Archivo3<<x<<"\t"<<PHI[i][j]<<"\n";
        }
    }
}
cout<<endl;
}
Archivo2.close();
Archivo3.close();
t1 = clock();

double time = (double)(t1-t0)/CLOCKS_PER_SEC;
cout << "Execution Time: " << time << endl;
ofstream Archivo6;
Archivo6.open("Execution time.txt");
Archivo6<<time;
}

```

### 3 Smith Hutton UDS

```

#include <iostream>
#include <fstream>
#include <math.h> // To use pow
#include <vector> // To use vectors
#include <ctime>
using namespace std;

double const L = 2;
double const H = 1;
int const Nx = 200;
int const Ny = 100;
double const inct = 0.0001;
double const densitat = 1000000;
double const gamma = 1;
double const incx = L/Nx;
double const incy = H/Ny;
//double const temps = 200;
double const alpha = 10;

```

```

double const epsilon = 0.000001;

double Calculvelu(double x, double y){
    double u = 2*y*(1-pow(x,2));
    return u;
}

double Calculvelv(double x, double y){
    double v = -2*x*(1-pow(y,2));
    return v;
}

double Calculinlet(double PHI, double x){
    double r = 1 + tanh((2*x+1)*alpha);
    return r;
}

double CalculUDS(double PHLee, double PHIww, double PHInn, double PHIss, double PHI, double uw, double ue, double vn, double vs){
    double ae = (gamma*incy)/incx;
    double aw = (gamma*incy)/incx;
    double an = (gamma*incx)/incy;
    double as = (gamma*incx)/incy;
    double PHLe, PHLw, PHIn, PHIs;
    if(ue > 0){
        PHLe = PHI;
    }else if( ue <= 0 ){
        PHLe = PHLee;
    }
    if(uw > 0){
        PHLw = PHIww;
    }else if( uw <= 0 ){
        PHLw = PHI;
    }
    if(vn > 0){
        PHIn = PHI;
    }else if( vn <= 0 ){
        PHIn = PHInn;
    }
    if(vs > 0){
        PHIs = PHIss;
    }else if( vs <= 0 ){
        PHIs = PHI;
    }

    double conveccio = densitat*(PHLe*incy*ue - PHLw*incy*uw + PHIn*incx*vn - PHIs*incx*vs);
    double diffusio = ae*PHLee + aw*PHLww + an*PHInn + as*PHIss - PHI*(an+aw+as+ae);
    double r = (inct/(incx*incy*densitat)) * (-conveccio + diffusio) + PHI;
    return r;
}

```

```

int main(){
  unsigned t0, t1;
  t0=clock();

  double PHI[Nx][Ny];
  double PHiant[Nx][Ny];
  double Vu[Nx][Ny];
  double Vv[Nx][Ny];
  double xi = -(L/2) + (incx/2);
  double yi = H - (incy/2);
  double x;
  double y;
  double v;
  double u;
  ofstream Archivo1;
  ofstream Archivo2;
  ofstream Archivo3;
  Archivo1.open("Mapa de velocitats.txt");
  Archivo1<<"x"<<"\t"<<"y"<<"\t"<<"z"<<"\t"<<"u"<<"\t"<<"v"<<"\n";

  Archivo2.open("Mapa Propietat.txt");
  Archivo2<<"x"<<"\t"<<"y"<<"\t"<<"z"<<"\t"<<"PHI"<<"\n";

  Archivo3.open("Outlet propietat.txt");
  Archivo3<<"x"<<"\t"<<"PHI"<<"\n";

  //DECLAREM MAPA DE VELOCITATS
  for(int j = 0; j<Ny; j++){
    for(int i = 0; i<Nx; i++){
      x = xi + i*incx;
      y = yi - j*incy;
      u = Calculvelu(x,y);
      v = Calculvelv(x,y);
      Vu[i][j] = u;
      Vv[i][j] = v;
      PHI[i][j] = 0;
      PHiant[i][j] = PHI[i][j];
      //cout<<" ("<<Vu[i][j]<<","<<Vv[i][j]<<") ";
      cout<<PHI[i][j]<<" ";
      Archivo1<<x<<"\t"<<y<<"\t"<<0<<"\t"<<Vu[i][j]<<"\t"<<Vv[i][j]<<"\n";
    }
    cout<<endl;
  }
  Archivo1.close();

  // double tempsmaxim = temps/inct;

  bool conv = false;
  int k = 0;

```

```

while(!conv){

    conv = true;

    for(int j = 0; j<Ny; j++){
        for(int i = 0; i<Nx; i++){
            x = xi + i*incx;
            y = yi - j*incy;
            //CALCUL NODES CENTRALS
            if(x > xi and y < yi and x < -xi and j<Ny-1){
                PHIant[i][j] = PHI[i][j];
                PHI[i][j] = CalculUDS(PHI[i+1][j], PHI[i-1][j], PHI[i][j-1], PHI[i][j+1], PHI[i][j], Vu[i-1][j], Vu[i+1][j], Vv[i][j-1], Vv[i][j+1]);
                if(abs(PHIant[i][j]-PHI[i][j]) > epsilon){
                    conv = false;
                    //cout<<PHIant[i][j]<<"--"<<PHI[i][j]<<endl;
                }
            }
            //INLET
            else if(j == Ny-1 and x < 0 and x > xi){
                PHIant[i][j] = PHI[i][j];
                PHI[i][j] = Calculinlet(PHI[i][j], x);
            }
            //OUTLET
            else if(j == Ny-1 and x > 0 and x < -xi){
                PHIant[i][j] = PHI[i][j];
                PHI[i][j] = PHI[i][j-1];
                if(abs(PHIant[i][j]-PHI[i][j]) > epsilon){
                    conv = false;
                    //cout<<PHIant[i][j]<<"--"<<PHI[i][j]<<endl;
                }
            }
            //REST OF WALLS
            else if(abs(x-xi)<0.0001 or abs(x+xi)<0.0001 or abs(y-yi)<0.0001){
                PHIant[i][j] = PHI[i][j];
                PHI[i][j] = 1-tanh(10);
            }
        }
    }

    k++;
    if(k < 150000){
        conv = false;
    }
}

```

```

cout<<"-----" << endl;
for(int j = 0; j<Ny; j++){ for(int i = 0; i<Nx; i++){
  x = xi + i*incx;
  y = yi - j*incy;
  cout<<PHI[i][j]<<" ";
  Archivo2<<x<<"\t"<<y<<"\t"<<0<<"\t"<<PHI[i][j]<<"\n";
  if(j==Ny-1){
    Archivo3<<x<<"\t"<<PHI[i][j]<<"\n";
  }
}
}
cout<<endl;
}
Archivo2.close();
Archivo3.close();
t1 = clock();

double time = (double(t1-t0)/CLOCKS_PER_SEC);
cout << "Execution Time: " << time << endl;
ofstream Archivo6;
Archivo6.open("Execution time.txt");
Archivo6<<time;
}

```

## 4 Lid Driven Cavity

```

#include <iostream>
#include <fstream>
#include <math.h> // To use pow
#include <vector> // To use vectors
#include <ctime>
using namespace std;

int main(){
  unsigned t0, t1;

  t0=clock();
  double L = 1;
  int Ny = 128;
  int Nx = 128;
  double incx = L/Nx;
  double incy = L/Ny;

```

```

double Uref = 1;
double temps = 45;
double inct = 0.001;
double tempsmaxim = temps/inct;
double densitat = 100;
double nyu = 1;
double Re = (densitat*Uref*L)/(nyu);
double Ru = 0;
double Rv = 0;
double itefin = false;
double epsilon = 0.00001;

double Vpu[Nx+1][Ny];
double Vu[Nx+1][Ny];
double Vpv[Nx][Ny+1];
double Vv[Nx][Ny+1];
double P[Nx][Ny];
double Ruant[Nx+1][Ny];
double Rvant[Nx][Ny+1];

//INTRODUIM MALLA Vpu i Vu
for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx+1; i++){
    if(j == Ny-1){
      Vpu[i][j] = Uref;
      Vu[i][j] = Uref;
    }
    else{
      Vpu[i][j] = 0;
      Vu[i][j] = 0;
    }
    Ruant[i][j] = 0;
  }
}

//INTRODUIM MALLA Vpv i Vv
for(int j = 0; j<Ny+1; j++){
  for(int i = 0; i<Nx; i++){
    Vpv[i][j] = 0;
    Vv[i][j] = 0;
    Rvant[i][j] = 0;
  }
}

//INTRODUIM MALLA P
for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    P[i][j] = 0;
  }
}

```

```

for(int k = 0; k<tempsmaxim; k++){

if(k%1000==0){
  cout<<k/1000<<endl;
}

//*****
//          STEP 1
//*****
//Calcul Velocitat predictoria u

for(int j = 0; j<Ny; j++){
for(int i = 0; i<Nx+1; i++){
  if(i>0 and i <Nx and j>0 and j<Ny-1){
    double Vup = Vu[i][j];
    double Vvn = 0.5*(Vv[i-1][j+1]+Vv[i][j+1]);
    double Vvs = 0.5*(Vv[i-1][j]+Vv[i][j]);
    double Vue = 0.5*(Vu[i+1][j]+Vu[i][j]);
    double Vuw = 0.5*(Vu[i-1][j]+Vu[i][j]);
    double Vun = 0.5*(Vu[i][j+1]+Vu[i][j]);
    double Vus = 0.5*(Vu[i][j-1]+Vu[i][j]);
    double Un = Vu[i][j+1];
    double Us = Vu[i][j-1];
    double Ue = Vu[i+1][j];
    double Uw = Vu[i-1][j];
    //Calcul up
    Ru = -((densitat*incx)*(Vvn*Vun-Vvs*Vus+Vue*Vue-Vuw*Vuw)) + nyu*(Un+Us+Ue+Uw-4*Vup);
    //Vpu[i][j] = Vu[i][j] + ((inct/(densitat*incx*incx))*(1.5*Ru-0.5*Ruant[i][j]));
    Vpu[i][j] = Vu[i][j] + ((inct/(densitat*incx*incx))*(1*Ru));
    Ruant[i][j] = Ru;
  }
  else if(j == Ny-1){
    Vpu[i][j] = Uref;
  }
  else {
    Vpu[i][j] = 0;
  }
}
}

//Calcul velocitat predictoria v

for(int j = 0; j<Ny+1; j++){
for(int i = 0; i<Nx; i++){
  if(i>0 and i <Nx-1 and j>0 and j<Ny){
    double Vvp = Vv[i][j];
    double Vvn = 0.5*(Vv[i][j+1]+Vv[i][j]);
    double Vvs = 0.5*(Vv[i][j-1]+Vv[i][j]);
    double Vve = 0.5*(Vv[i+1][j]+Vv[i][j]);
    double Vvw = 0.5*(Vv[i-1][j]+Vv[i][j]);
  }
}
}

```

```

double Vue = 0.5*(Vu[i+1][j]+Vu[i+1][j-1]);
double Vuw = 0.5*(Vu[i][j]+Vu[i][j-1]);
double Vn = Vv[i][j+1];
double Vs = Vv[i][j-1];
double Ve = Vv[i+1][j];
double Vw = Vv[i-1][j];
//Calcul vp
Rv = -(densitat*incx)*(Vvn*Vvn-Vvs*Vvs+Vue*Vve-Vuw*Vvw) + nyu*(Vn+Vs+Ve+Vw-4*Vvp);
//Vpv[i][j] = Vv[i][j] + (inct/(densitat*incx*incx))*(1.5*Rv - 0.5*Rvant[i][j]);
Vpv[i][j] = Vv[i][j] + ((inct/(densitat*incx*incx))*(1*Rv));
Rvant[i][j] = Rv;
}
else{
  Vpv[i][j] = 0;
}
}
}

//*****
//                STEP 2
//*****

//Calcul camp de Pressio
itefin = false;
while(!itefin){
  itefin = true;
  for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    double Pant = P[i][j];
    double Pn = 0;
    double Ps = 0;
    double Pe = 0;
    double Pw = 0;
    double vpn = Vpv[i][j+1];
    double vps = Vpv[i][j];
    double upe = Vpu[i+1][j];
    double upw = Vpu[i][j];
    double an = 0;
    double as = 0;
    double ae = 0;
    double aw = 0;

    if(i != 0){
      aw = 1;
      Pw = P[i-1][j];
    }
    if(i != Nx-1){
      ae = 1;
      Pe = P[i+1][j];
    }
  }
}

```



```

}
if(j != 0){
  as = 1;
  Ps = P[i][j-1];
}
if(j != Ny-1){
  an = 1;
  Pn = P[i][j+1];
}
double ap = an+as+ae+aw;
P[i][j] = (1/ap)*(an*Pn+as*Ps+ae*Pe+aw*Pw - ((densitat*incx)/(inct))*(vpn-vps+upe-upw));
double error = fabs(Pant - P[i][j]);
if(error > epsilon) itefin = false;
}
}
}

//*****
//          STEP 3
//*****

//Calcul velocitat u

for(int j = 0; j<Ny; j++){
for(int i = 0; i<Nx+1; i++){
  if(i>0 and i <Nx and j>0 and j<Ny-1){
    Vu[i][j] = Vpu[i][j] - (inct/densitat)*(P[i][j]-P[i-1][j])/incx);
  }
  else if(j == Ny-1){
    Vu[i][j] = Uref;
  }
  else {
    Vu[i][j] = 0;
  }
}
}

//Calcul velocitat v

for(int j = 0; j<Ny+1; j++){
for(int i = 0; i<Nx; i++){
  if(i>0 and i <Nx-1 and j>0 and j<Ny){
    Vv[i][j] = Vpv[i][j] - (inct/densitat)*(P[i][j]-P[i][j-1])/incy);
  }
  else{
    Vv[i][j] = 0;
  }
}
}
}

```

```

}
ofstream Archivo1;
Archivo1.open("Velocitat.u.txt");
Archivo1<<"Y"<<"\t"<<"u"<<"\n";
ofstream Archivo2;
Archivo2.open("Velocitat.v.txt");
Archivo2<<"X"<<"\t"<<"v"<<"\n";

cout<<"-----MAPA Vpu-----"<<endl;

for(int j = Ny-1; j>=0; j--){
  for(int i = 0; i<=Nx; i++){
    cout<<Vpu[i][j]<<" ";
  }
  cout<<endl;
}
cout<<endl;
cout<<"-----MAPA Vu-----"<<endl;
cout<<endl;
for(int j = Ny-1; j>=0; j--){
  for(int i = 0; i<=Nx; i++){
    cout<<Vu[i][j]<<" ";
    if(i == Nx/2){
      Archivo1<<(incy/2) + j*incy<<"\t"<<Vu[i][j]<<"\n";
    }
  }
  cout<<endl;
}
cout<<endl;
cout<<"-----MAPA Vpv-----"<<endl;
cout<<endl;
for(int j = Ny; j>=0; j--){
  for(int i = 0; i<Nx; i++){
    cout<<Vpv[i][j]<<" ";
  }
  cout<<endl;
}
cout<<endl;
cout<<"-----MAPA Vv-----"<<endl;
cout<<endl;
for(int j = Ny; j>=0; j--){
  for(int i = 0; i<Nx; i++){
    cout<<Vv[i][j]<<" ";
    if(j == Nx/2){
      Archivo2<<(incx/2) + i*incx<<"\t"<<Vv[i][j]<<"\n";
    }
  }
  cout<<endl;
}
}

```

```

cout<<endl;
cout<<"-----MAPA P-----"<<endl;
cout<<endl;
for(int j = Ny-1; j>=0; j--){
    for(int i = 0; i<Nx; i++){
        cout<<P[i][j]<<" ";
    }
    cout<<endl;
}

ofstream Archivo3;
Archivo3.open("Mapa_velocitat.txt");
Archivo3<<"X"<<"\t"<<"Y"<<"\t"<<"Z"<<"\t"<<"u"<<"\t"<<"v"<<"\t"<<"P"<<"\n";
double u = 0;
double v = 0;
for(int j = 0; j<Ny; j++){
    for(int i = 0; i<Nx; i++){
        u = 0.5*(Vu[i][j]+Vu[i+1][j]);
        v = 0.5*(Vv[i][j]+Vv[i+1][j]);
        Archivo3<<(incx/2)+i*incx<<"\t"<<(incy/2)+j*incy<<"\t"<<0<<"\t"<<u<<"\t"<<v<<"\t"<<P[i][j]<<"\n";
    }
}
t1 = clock();

double time = (double)(t1-t0)/CLOCKS_PER_SEC;
cout << "Execution Time: " << time << endl;
ofstream Archivo6;
Archivo6.open("Execution time.txt");
Archivo6<<time;
}

```

## 5 Differentially Heated Cavity

```

#include <iostream>
#include <fstream>
#include <math.h> // To use pow
#include <vector> // To use vectors
using namespace std;

int main(){

double L = 1;
int Ny = 81;
int Nx = 81;
double incx = L/Nx;
double incy = L/Ny;

```

```

double Uref = 0;
double temps = 30;
double inct = 0.00001;
double tempsmaxim = temps/inct;
double densitat = 1;
double nyu = 1;
double Re = (densitat*Uref*L)/(nyu);
double Ru = 0;
double Rv = 0;
double itefin = false;
double epsilon = 0.00001;

double Thot = 1;
double Tcold = 0;
double cp = 1;
double beta = 1;
double lambda = 1;
double gravetat = 1E6;
double gamma = lambda/cp;
double Ra = (cp*gravetat*beta*pow(densitat,2)*(Thot-Tcold)*pow(L,3))/(nyu*lambda);
double Tfluid = 0.5;
cout<<"El nombre Ra es: "<<Ra<<endl;

string esquema = "CDS"; //UDS or CDS
cout<<"L'esquema seleccionat es: "<<esquema<<endl;
double up[Nx+1][Ny];
double u[Nx+1][Ny];
double vp[Nx][Ny+1];
double v[Nx][Ny+1];
double P[Nx][Ny];
double T[Nx][Ny];

//INTRODUIM MALLA Vpu i Vu
for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx+1; i++){
    if(j == Ny-1){
      up[i][j] = Uref;
      u[i][j] = Uref;
    }
    else{
      up[i][j] = 0;
      u[i][j] = 0;
    }
  }
}

//INTRODUIM MALLA Vpv i Vv
for(int j = 0; j<Ny+1; j++){
  for(int i = 0; i<Nx; i++){
    vp[i][j] = 0;
  }
}

```

```

    v[i][j] = 0;
  }
}
//INTRODUIM MALLA P i T
for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    if(i==0 and j<Ny-1 and j>0){
      T[i][j] = Thot;
    }
    else if(i==Nx-1 and j<Ny-1 and j>0){
      T[i][j] = Tcold;
    }
    else{
      T[i][j] = 0;
    }
    P[i][j] = 0;
  }
}

for(int k = 0; k<tempsmaxim; k++){

int iteprint = 10000;
if(k%iteprint==0){
  cout<<k<<endl;
}

//*****
//                               STEP 1
//*****
//Calcul Velocitat predictorua u

for(int j = 0; j<Ny; j++){
for(int i = 0; i<Nx+1; i++){
  if(i>0 and i <Nx and j>0 and j<Ny-1){
    double uP = u[i][j];
    double vn = 0.5*(v[i-1][j+1]+v[i][j+1]);
    double vs = 0.5*(v[i-1][j]+v[i][j]);
    double ue = 0.5*(u[i+1][j]+u[i][j]);
    double uw = 0.5*(u[i-1][j]+u[i][j]);
    double un = 0.5*(u[i][j+1]+u[i][j]);
    double us = 0.5*(u[i][j-1]+u[i][j]);
    double Un = u[i][j+1];
    double Us = u[i][j-1];
    double Ue = u[i+1][j];
    double Uw = u[i-1][j];
    //Calcul up
    Ru = -((densitat*incx)*(vn*un-vs*us+ue*ue-uw*uw)) + nyu*(Un+Us+Ue+Uw-4*uP); // + incx*beta*(T[i][j]-Tfluid)*gravetat;
    up[i][j] = u[i][j] + ((inct/(densitat * incx * incx)) * Ru);
  }
}
}

```

```

}
else if(j == Ny-1){
  up[i][j] = Uref;
}
else {
  up[i][j] = 0;
}
}
}
}

//Calcul velocitat predictoria v

for(int j = 0; j<Ny+1; j++){
for(int i = 0; i<Nx; i++){
  if(i>0 and i <Nx-1 and j>0 and j<Ny){
    double vP = v[i][j];
    double vn = 0.5*(v[i][j+1]+v[i][j]);
    double vs = 0.5*(v[i][j-1]+v[i][j]);
    double ve = 0.5*(v[i+1][j]+v[i][j]);
    double vw = 0.5*(v[i-1][j]+v[i][j]);
    double ue = 0.5*(u[i+1][j]+u[i+1][j-1]);
    double uw = 0.5*(u[i][j]+u[i][j-1]);
    double Vn = v[i][j+1];
    double Vs = v[i][j-1];
    double Ve = v[i+1][j];
    double Vw = v[i-1][j];
    //Calcul vp
    Rv = -(densitat*incx)*(vn*vn-vs*vs+ue*ve-uw*vw) + nyu*(Vn+Vs+Ve+Vw-4*vP) ;
    vp[i][j] = v[i][j] + ((inct/(densitat*incx*incx))*(1*Rv) + inct*beta*(T[i][j]-Tfluid)*gravetat);
  }
  else{
    vp[i][j] = 0;
  }
}
}

//*****
//                STEP 2
//*****

//Calcul camp de Pressi
itefin = false;
while(!itefin){
  itefin = true;
  for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    double Pant = P[i][j];
    double Pn = 0;
    double Ps = 0;
    double Pe = 0;

```

```

double Pw = 0;
double vpn = vp[i][j+1];
double vps = vp[i][j];
double upe = up[i+1][j];
double upw = up[i][j];
double an = 0;
double as = 0;
double ae = 0;
double aw = 0;

if(i != 0){
    aw = 1;
    Pw = P[i-1][j];
}
if(i != Nx-1){
    ae = 1;
    Pe = P[i+1][j];
}
if(j != 0){
    as = 1;
    Ps = P[i][j-1];
}
if(j != Ny-1){
    an = 1;
    Pn = P[i][j+1];
}
double ap = an+as+ae+aw;
P[i][j] = (1/ap)*(an*Pn+as*Ps+ae*Pe+aw*Pw - ((densitat*incx)/(inct))*(vpn-vps+upe-upw));
double error = fabs(Pant - P[i][j]);
if(error > epsilon) itefin = false;

}
}

//*****
//                               STEP 3
//*****

//Calcul velocitat u

for(int j = 0; j<Ny; j++){
for(int i = 0; i<Nx+1; i++){
    if(i>0 and i <Nx and j>0 and j<Ny-1){
        u[i][j] = up[i][j] - (inct/densitat)*((P[i][j]-P[i-1][j])/incx);
    }
    else if(j == Ny-1){

```

```

    u[i][j] = Uref;
  }
  else {
    u[i][j] = 0;
  }
}
}
}

//Calcul velocitat v

for(int j = 0; j<Ny+1; j++){
  for(int i = 0; i<Nx; i++){
    if(i>0 and i <Nx-1 and j>0 and j<Ny){
      v[i][j] = vp[i][j] - (inct/densitat)*((P[i][j]-P[i][j-1])/incy);
    }
    else{
      v[i][j] = 0;
    }
  }
}

//*****
//
//          STEP 4
//*****

if(esquema == "CDS"){
  //Calcul camp de Temperatura CDS

  for(int j = 0; j<Ny; j++){
    for(int i = 0; i<Nx; i++){

      if(i>0 and j>0 and i<Nx-1 and j<Ny-1){
        double uw = u[i][j];
        double ue = u[i+1][j];
        double vn = v[i][j+1];
        double vs = v[i][j];
        double ae = (gamma*incy)/incx;
        double aw = (gamma*incy)/incx;
        double an = (gamma*incx)/incy;
        double as = (gamma*incx)/incy;
        double Tp = T[i][j];
        double Te = (T[i+1][j] + Tp)/2.0;
        double Tw = (T[i-1][j] + Tp)/2.0;
        double Tn = (T[i][j+1] + Tp)/2.0;
        double Ts = (T[i][j-1] + Tp)/2.0;
        double TE2 = T[i+1][j];
        double TW2 = T[i-1][j];
        double TN2 = T[i][j+1];
        double TS2 = T[i][j-1];
        double conveccio = densitat*(Te*incy*ue - Tw*incy*uw + Tn*incx*vn - Ts*incx*vs);

```



```

    double diffusio = ae*TE2 + aw*TW2 + an*TN2 + as*TS2 - Tp*(an+aw+as+ae);
    T[i][j] = (inct/(incx*incy*densitat)) * (-conveccio + diffusio) + Tp;
  } else if(i==0 and j<Ny-1 and j>0){
    T[i][j] = Thot;
  }
  else if(i==Nx-1 and j<Ny-1 and j>0){
    T[i][j] = Tcold;
  }
  else if(j == 0){
    T[i][j] = T[i][j+1];
  }
  else if(j == Ny-1){
    T[i][j] = T[i][j-1];
  }
}
} else if(esquema == "UDS"){

//Calcul camp de Temperatura UDS

for(int j = 0; j<Ny; j++){
for(int i = 0; i<Nx; i++){

if(i>0 and j>0 and i<Nx-1 and j<Ny-1){
  double uw = u[i][j];
  double ue = u[i+1][j];
  double vn = v[i][j+1];
  double vs = v[i][j];
  double ae = (gamma*incy)/incx;
  double aw = (gamma*incy)/incx;
  double an = (gamma*incx)/incy;
  double as = (gamma*incx)/incy;
  double Tp = T[i][j];
  double TE2 = T[i+1][j];
  double TW2 = T[i-1][j];
  double TN2 = T[i][j+1];
  double TS2 = T[i][j-1];
  double Te, Tw, Tn, Ts;
if(ue > 0){
  Te = Tp;
} else if( ue <= 0 ){
  Te = TE2;
}
if(uw > 0){
  Tw = TW2;
} else if( uw <= 0 ){
  Tw = Tp;
}
}
}
}

```

```

if(vn > 0){
    Tn = Tp;
}else if( vn <= 0){
    Tn = TN2;
}
if(vs > 0){
    Ts = TS2;
}else if( vs <= 0 ){
    Ts = Tp;
}
double conveccio = densitat*(Te*incy*ue - Tw*incy*uw + Tn*incx*vn - Ts*incx*vs);
double diffusio = ae*TE2 + aw*TW2 + an*TN2 + as*TS2 - Tp*(an+aw+as+ae);
T[i][j] = (inct/(incx*incy*densitat)) * (-conveccio + diffusio) + Tp;
}
else if(i==0 and j<Ny-1 and j>0){
    T[i][j] = Thot;
}
else if(i==Nx-1 and j<Ny-1 and j>0){
    T[i][j] = Tcold;
}
else if(j == 0){
    T[i][j] = T[i][j+1];
}
else if(j == Ny-1){
    T[i][j] = T[i][j-1];
}
}
}
}
//*****
//          CALCUL NUSSELT
//*****
//Es parteix dels punts següents:
double alpha = lambda/(densitat*cp);
double Nu = 0;
for (int i = 1; i<Nx; i++){
    double Nux = 0;
    for(int j = 0; j<Ny; j++){
        double uprima = (u[i][j] * L)/(alpha);
        double Tprima = (T[i][j] - Tcold)/(Thot-Tcold);
        double Tprimaesquerra = (T[i-1][j] - Tcold)/(Thot-Tcold);
        Nux = Nux + (uprima*Tprima - ((Tprima-Tprimaesquerra)/incx)) * incy;
    }
    Nu = Nu + (Nux) * incx;
}
cout<<i<<" : "<<Nux<<endl;

```

```

}

ofstream Archivo1;
Archivo1.open("Velocitat_u.txt");
Archivo1<<"Y"<<"\t"<<"u"<<"\n";
ofstream Archivo2;
Archivo2.open("Velocitat_v.txt");
Archivo2<<"X"<<"\t"<<"v"<<"\n";

cout<<"-----MAPA Vpu-----"<<endl;

for(int j = Ny-1; j>=0; j--){
  for(int i = 0; i<=Nx; i++){
    cout<<up[i][j]<<" ";
  }
  cout<<endl;
}

double Umax = 0;
double yymax = 0;

cout<<endl;
cout<<"-----MAPA Vu-----"<<endl;
cout<<endl;
for(int j = Ny-1; j>=0; j--){
  for(int i = 0; i<=Nx; i++){
    cout<<u[i][j]<<" ";
    if(i == Nx/2){
      if(u[i][j] > Umax){
        Umax = u[i][j];
        yymax = (incy/2) + j*incy;
      }
      Archivo1<<(incy/2) + j*incy<<"\t"<<u[i][j]<<"\n";
    }
  }
  cout<<endl;
}
cout<<endl;
cout<<"-----MAPA Vpv-----"<<endl;
cout<<endl;
for(int j = Ny; j>=0; j--){
  for(int i = 0; i<Nx; i++){
    cout<<vp[i][j]<<" ";
  }
  cout<<endl;
}
double Vmax = 0;
double xvmax = 0;
cout<<endl;

```

```

cout<<"-----MAPA Vv-----"<<endl;
cout<<endl;
for(int j = Ny; j>=0; j--){
  for(int i = 0; i<Nx; i++){
    cout<<v[i][j]<<" ";
    if(j == Nx/2){
      if(v[i][j] > Vmax){
        Vmax = v[i][j];
        xvmax = (incx/2) + i*incx;
      }
      Archivo2<<(incx/2) + i*incx<<"\t"<<v[i][j]<<"\n";
    }
  }
  cout<<endl;
}
cout<<endl;
cout<<"-----MAPA P-----"<<endl;
cout<<endl;
for(int j = Ny-1; j>=0; j--){
  for(int i = 0; i<Nx; i++){
    cout<<P[i][j]<<" ";
  }
  cout<<endl;
}
cout<<"-----MAPA T-----"<<endl;
cout<<endl;
for(int j = Ny-1; j>=0; j--){
  for(int i = 0; i<Nx; i++){
    cout<<T[i][j]<<" ";
  }
  cout<<endl;
}
cout<<"-----Nusselt-----"<<endl;
ofstream Archivo5;
Archivo5.open("Nusselt.txt");
Archivo5<<"El nombre de Nusselt es:"<<"\t"<<Nu<<"\n";
Archivo5<<"La velocitat m gran u : "<<"\t"<<Umax<<"\t"<<"En la posici y: "<<"\t"<<yumax<<"\n";
Archivo5<<"La velocitat m gran v : "<<"\t"<<Vmax<<"\t"<<"En la posici x: "<<"\t"<<xvmax<<"\n";
cout<<Nu<<endl;

ofstream Archivo3;
Archivo3.open("Mapa_velocitat.txt");
Archivo3<<"X"<<"\t"<<"Y"<<"\t"<<"Z"<<"\t"<<"u"<<"\t"<<"v"<<"\t"<<"P"<<"\n";
double U = 0;
double V = 0;
for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    U = 0.5*(u[i][j]+u[i+1][j]);
    V = 0.5*(v[i][j]+v[i][j+1]);
  }
}

```

```

    Archivo3<<(incx/2)+i*incx<<"\t"<<(incy/2)+j*incy<<"\t"<<0<<"\t"<<U<<"\t"<<V<<"\t"<<T[i][j]<<"\n";
  }
}

ofstream Archivo4;
Archivo4.open("Mapa_Temperatura.txt");
Archivo4<<"X"<<"\t"<<"Y"<<"\t"<<"Z"<<"\t"<<"T"<<"\n";
for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    Archivo4<<(incx/2)+i*incx<<"\t"<<(incy/2)+j*incy<<"\t"<<0<<"\t"<<T[i][j]<<"\n";
  }
}
}

```

## 6 Square Cylinder

```

#include <iostream>
#include <fstream>
#include <math.h> // To use pow
#include <vector> // To use vectors
#include <ctime>
using namespace std;

int main(){
  unsigned t0, t1;
  t0=clock();
  double D = 1;
  double Lxx = 26;
  double Lyy = 20;
  int Nx = 150;
  int Ny = 120;
  double Uref2= 1;
  double temps = 200;
  double inct = 0.0001;
  double tempmaxim = temps/inct;
  double nyu = 1;
  double Reynolds = 100;
  double densitat = Reynolds*nyu/(Uref2*D);
  double Re = (densitat*Uref2*D)/(nyu);
  double Ru = 0;
  double Rv = 0;
  double itefin = false;
  double epsilon = 0.00001;

  cout<<"El Reynolds es: "<<Re<<endl;
  cout<<"La densitat es "<<densitat<<endl;

```

```

vector< vector<double> > up(Nx+1, vector<double> (Ny, 0));
vector< vector<double> > Ruant(Nx+1, vector<double> (Ny, 0));
vector< vector<double> > u(Nx+1, vector<double> (Ny, 0));
vector< vector<double> > vp(Nx, vector<double> (Ny+1, 0));
vector< vector<double> > v(Nx, vector<double> (Ny+1, 0));
vector< vector<double> > Rvant(Nx, vector<double> (Ny+1, 0));
vector< vector<double> > P(Nx, vector<double> (Ny, 0));

vector< vector<double> > x(Nx, vector<double> (Ny, 0));
vector< vector<double> > y(Nx, vector<double> (Ny, 0));
vector< vector<double> > incx(Nx, vector<double> (Ny, 0));
vector< vector<double> > incy(Nx, vector<double> (Ny, 0));

//INTRODUIM MALLA

double La = 8;
double Lb = 2;
double Lc = 16;
double Lu = 9;
double Lv = 2;
double Lw = 9;

int N1 = Nx/5;
int N2 = Nx/5;
int N3 = 3.0*Nx/5.0;
int N4 = 2*Ny/5.0;
int N5 = Ny/5.0;
int N6 = 2*Ny/5.0;

int Na = N1 + 0.5/(Lb/N2);
int Nb = Na + 1/(Lb/N2);
int Nu = N4 + 0.5/(Lv/N5);
int Nv = Nu + 1/(Lv/N5);

double Lx, Ly, Nxx, Nyy, A, B, c1, c2, ii, jj, xo, yo, incxx, incyy;

for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    if(j<N4){

      if(i<N1){
        incx[i][j] = La/N1;
        x[i][j] = (La/N1)/2 + i*La/N1;
      }
      if(i>=N1 and i<=N2+N1){
        incx[i][j] = Lb/N2;
        x[i][j] = x[i-1][j] + incx[i-1][j]/2 + incx[i][j]/2;
      }
    }
  }
}

```

```

if(i>N2+N1){
    incx[i][j] = Lc/N3;
    x[i][j] = x[i-1][j] + incx[i-1][j]/2 + incx[i][j]/2;
}

incy[i][j] = Lu/N4;
y[i][j] = (Lu/N4)/2 + j*Lu/N4;
}
if(j>=N4 and j<=N5+N4){

if(i<N1){
    incx[i][j] = La/N1;
    x[i][j] = (La/N1)/2 + i*La/N1;
}
if(i>=N1 and i<=N2+N1){
    incx[i][j] = Lb/N2;
    x[i][j] = x[i-1][j] + incx[i-1][j]/2 + incx[i][j]/2;
}
if(i>N2+N1){
    incx[i][j] = Lc/N3;
    x[i][j] = x[i-1][j] + incx[i-1][j]/2 + incx[i][j]/2;
}

incy[i][j] = Lv/N5;
y[i][j] = y[i][j-1] + incy[i][j-1]/2 + incy[i][j]/2;
}
if(j>N5+N4){

if(i<N1){
    incx[i][j] = La/N1;
    x[i][j] = (La/N1)/2 + i*La/N1;
}
if(i>=N1 and i<=N2+N1){
    incx[i][j] = Lb/N2;
    x[i][j] = x[i-1][j] + incx[i-1][j]/2 + incx[i][j]/2;
}
if(i>N2+N1){
    incx[i][j] = Lc/N3;
    x[i][j] = x[i-1][j] + incx[i-1][j]/2 + incx[i][j]/2;
}

incy[i][j] = Lw/N6;
y[i][j] = y[i][j-1] + incy[i][j-1]/2 + incy[i][j]/2;
}
}
}

//INTRODUIM MALLA Vpu i Vu
for(int j = 0; j<Ny; j++){

```

```

for(int i = 0; i<Nx+1; i++){
  if(j == Ny-1){
    up[i][j] = 0;
    u[i][j] = 0;
  }else if(i == 0){
    up[i][j] = Uref2;
    u[i][j] = Uref2;
  }else if(i == Nx){
    up[i][j] = 0;
    u[i][j] = 0;
  }else if(j == 0){
    up[i][j] = 0;
    u[i][j] = 0;
  }else{
    up[i][j] = 0;
    u[i][j] = 0;
  }
  Ruant[i][j] = 0;
}
}
//INTRODUIM MALLA Vpv i Vv
for(int j = 0; j<Ny+1; j++){
  for(int i = 0; i<Nx; i++){
    vp[i][j] = 0;
    v[i][j] = 0;
    Rvant[i][j] = 0;
  }
}
//INTRODUIM MALLA P
for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    P[i][j] = 0;
  }
}

for(int k = 0; k<tempsmaxim; k++){

  if(k%1000==0){
    cout<<k/1000<<endl;
  }

  //*****
  //                STEP 1
  //*****
  //Calcul Velocitat predictoria u

  for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx+1; i++){

```



```

if(i>0 and i <Nx and j>0 and j<Ny-1){
  if(i>=Na and i<=Nb and j>=Nu and j<=Nv-1){
    up[i][j] = 0;
  }else{
    double uP = u[i][j];
    double vn = 0.5*(v[i-1][j+1]+v[i][j+1]);
    double vs = 0.5*(v[i-1][j]+v[i][j]);
    double ue = 0.5*(u[i+1][j]+u[i][j]);
    double uw = 0.5*(u[i-1][j]+u[i][j]);
    double un = 0.5*(u[i][j+1]+u[i][j]);
    double us = 0.5*(u[i][j-1]+u[i][j]);
    double Un = u[i][j+1];
    double Us = u[i][j-1];
    double Ue = u[i+1][j];
    double Uw = u[i-1][j];

    double dN = y[i][j+1]-y[i][j];
    double dS = y[i][j]-y[i][j-1];
    double dE = incx[i-1][j];
    double dW = incx[i][j];
    double Sn = x[i][j]-x[i-1][j];
    double Ss = x[i][j]-x[i-1][j];
    double Se = incy[i][j];
    double Sw = incy[i][j];

    //Calcul up
    Ru = -((densitat)*(vn*un*Sn-vs*us*Ss+ue*ue*Se-uw*uw*Sw)) + nyu*(((Un-uP)*Sn/dN)+((Us-uP)*Ss/dS)+((Ue-uP)*Se/dE)+((Uw-uP)*Sw/
    dW));
    //up[i][j] = u[i][j] + ((inct/(densitat * incx[i][j] * incy[i][j])) * (1.5 * Ru - 0.5 * Ruant[i][j]));
    up[i][j] = u[i][j] + ((inct/(densitat * Sn * Se)) * Ru);
    Ruant[i][j] = Ru;
  }
}
else if(j == Ny-1){
  up[i][j] = up[i][j-1];
}
else if(i == 0){
  up[i][j] = Uref2;
}
else if(i == Nx){
  up[i][j] = up[i-1][j];
}
else if(j == 0){
  up[i][j] = up[i][j+1];
}
}
}

//Calcul velocitat predictoria v

for(int j = 0; j<Ny+1; j++){
for(int i = 0; i<Nx; i++){

```

```

if(i>0 and i <Nx-1 and j>0 and j<Ny){
  if(i>=Na and i<=Nb-1 and j>=Nu and j<=Nv){
    vp[i][j] = 0;
  }else{
    double vP = v[i][j];
    double vn = 0.5*(v[i][j+1]+v[i][j]);
    double vs = 0.5*(v[i][j-1]+v[i][j]);
    double ve = 0.5*(v[i+1][j]+v[i][j]);
    double vw = 0.5*(v[i-1][j]+v[i][j]);
    double ue = 0.5*(u[i+1][j]+u[i+1][j-1]);
    double uw = 0.5*(u[i][j]+u[i][j-1]);
    double Vn = v[i][j+1];
    double Vs = v[i][j-1];
    double Ve = v[i+1][j];
    double Vw = v[i-1][j];

    double dN = incy[i][j];
    double dS = incy[i][j-1];
    double dE = x[i+1][j]-x[i][j];
    double dW = x[i][j]-x[i-1][j];
    double Sn = incx[i][j];
    double Ss = incx[i][j];
    double Se = y[i][j]-y[i][j-1];
    double Sw = y[i][j]-y[i][j-1];
    //Calcul vp
    Rv = -(densitat)*(vn*vn*Sn-vs*vs*Ss+ue*ve*Se-uw*vw*Sw) + nyu*(((Vn-vP)*Sn/dN)+((Vs-vP)*Ss/dS)+((Ve-vP)*Se/dE)+((Vw-vP)*Sw/
    dW));
    //vp[i][j] = v[i][j] + (inct/(densitat*incx[i][j]*incy[i][j]))*(1.5*Rv - 0.5*Rvant[i][j]);
    vp[i][j] = v[i][j] + (inct/(densitat*Sn*Se))*(1*Rv);
    Rvant[i][j] = Rv;
  }
}
else if(j == Ny){
  vp[i][j] = 0;
}else if(i == 0){
  vp[i][j] = 0;
}else if(i == Nx-1){
  vp[i][j] = vp[i-1][j];
}else if(j == 0){
  vp[i][j] = 0;
}
}
}

//*****
//                               STEP 2
//*****

//Calcul camp de Pressio

```

```

itefin = false;
double contpres = 0;
while(!itefin){
itefin = true;
for(int j = 0; j<Ny; j++){
for(int i = 0; i<Nx; i++){
double Pant = P[i][j];
double Pn = 0;
double Ps = 0;
double Pe = 0;
double Pw = 0;
double vpn = vp[i][j+1];
double vps = vp[i][j];
double upe = up[i+1][j];
double upw = up[i][j];
double an = 0;
double as = 0;
double ae = 0;
double aw = 0;

if(i != 0){
aw = incy[i][j]/(x[i][j]-x[i-1][j]);
Pw = P[i-1][j];
}
if(i != Nx-1){
ae = incy[i][j]/(x[i+1][j]-x[i][j]);
Pe = P[i+1][j];
}
if(j != 0){
as = incx[i][j]/(y[i][j]-y[i][j-1]);
Ps = P[i][j-1];
}
if(j != Ny-1){
an = incx[i][j]/(y[i][j+1]-y[i][j]);
Pn = P[i][j+1];
}

if(i==Na-1 and j>=Nu and j<=Nv-1){ //Paret esquerra quadrat
ae = 0;
}

if(i==Nb and j>=Nu and j<=Nv-1){ //Paret dreta quadrat
aw = 0;
}

if(j==Nv and i>=Na and i<=Nb-1){ //Paret adalt quadrat
as = 0;
}
}
}

```

```

if(j==Nu-1 and i>=Na and i<=Nb-1){ //Paret abaix quadrat
  an = 0;
}

double ap = an + as + ae + aw;
if(i == Nx-1){ //Sortida
  P[i][j] = 0;
}else if(i>=Na and i<=Nb-1 and j>=Nu and j<=Nv-1){ //Cuadrat
  P[i][j] = 0;
}
else{
  P[i][j] = (1/ap)*(an*Pn+as*Ps+ae*Pe+aw*Pw - ((densitat*incx[i][j])/(inct))*(vpn-vps) - ((densitat*incy[i][j])/(inct))*(upe-upw) );
}
double error = fabs(Pant - P[i][j]);
if(error > epsilon) itefin = false;
//cout<<error<<endl;
//cout<<P[i][j]<<endl;
}
} if(contpres >= 1000000){
  cout<<"Error calculant Pressio, No convergencia"<<endl;
  return 0;
}
contpres = contpres + 1;

if(isnan(P[Nx/4][Ny/4])){
  cout<<"Error NaN pressio"<<endl;
  ofstream Archivo7;
  Archivo7.open("Informe_Errors.txt");
  Archivo7<<"NaN Pressio";
  return 0;
}
}

//*****
// STEP 3
//*****

//Calcul velocitat u

for(int j = 0; j<Ny; j++){
for(int i = 0; i<Nx+1; i++){
  if(i>0 and i <Nx and j>0 and j<Ny-1){
    if(i>=Na and i<=Nb-1 and j>=Nu and j<=Nv-1){
      u[i][j] = 0;
    }else{
      u[i][j] = up[i][j] - (inct/densitat)*((P[i][j]-P[i-1][j])/(x[i][j]-x[i-1][j]));
    }
  }
}
}

```

```

}
else if(j == Ny-1){
    u[i][j] = u[i][j-1];
}else if(i == 0){
    u[i][j] = Uref2;
}else if(i == Nx){
    u[i][j] = u[i-1][j];
}else if(j == 0){
    u[i][j] = u[i][j+1];
}
}
}
}

//Calcul velocitat v

for(int j = 0; j<Ny+1; j++){
for(int i = 0; i<Nx; i++){
    if(i>0 and i <Nx-1 and j>0 and j<Ny){
        if(i>=Na and i<=Nb-1 and j>=Nu and j<=Nv-1){
            v[i][j] = 0;
        }else{
            v[i][j] = vp[i][j] - (inct/densitat)*((P[i][j]-P[i][j-1])/(y[i][j]-y[i][j-1]));
        }
    }
}
else if(j == Ny){
    v[i][j] = 0;
}else if(i == 0){
    v[i][j] = 0;
}else if(i == Nx-1){
    v[i][j] = v[i-1][j];
}else if(j == 0){
    v[i][j] = 0;
}
}
}

int num = k;
ofstream Archivo;
char nombre[500];
double U = 0;
double V = 0;
if(k%1000 == 0){
    sprintf(nombre, "Vel_%05d.txt", num);
    Archivo.open(nombre, ios::trunc);
    Archivo<<"X"<<"\t"<<"Y"<<"\t"<<"Z"<<"\t"<<"u"<<"\t"<<"v"<<"\t"<<"P"<<"\t"<<"incx"<<"\t"<<"incy"<<"\n";
    for(int j = 0; j<Ny; j++){
        for(int i = 0; i<Nx; i++){
            U = 0.5*(u[i][j]+u[i+1][j]);
            V = 0.5*(v[i][j]+v[i][j+1]);
            Archivo<<x[i][j]<<"\t"<<y[i][j]<<"\t"<<0<<"\t"<<U<<"\t"<<V<<"\t"<<P[i][j]<<"\t"<<incx[i][j]<<"\t"<<incy[i][j]<<"\n";

```

```

    }
  }
}

ofstream Archivo1;
Archivo1.open("Velocitat_u.txt");
Archivo1<<"Y"<<"\t"<<"u"<<"\n";
ofstream Archivo2;
Archivo2.open("Velocitat_v.txt");
Archivo2<<"X"<<"\t"<<"v"<<"\n";

cout<<"-----MAPA Vpu-----"<<endl;

for(int j = Ny-1; j>=0; j--){
  for(int i = 0; i<=Nx; i++){
    cout<<up[i][j]<<" ";
  }
  cout<<endl;
}
cout<<endl;
cout<<"-----MAPA Vu-----"<<endl;
cout<<endl;
for(int j = Ny-1; j>=0; j--){
  for(int i = 0; i<=Nx; i++){
    cout<<u[i][j]<<" ";
    if(i == Nx/2){
      Archivo1<<y[i][j]<<"\t"<<u[i][j]<<"\n";
    }
  }
  cout<<endl;
}
cout<<endl;
cout<<"-----MAPA Vpv-----"<<endl;
cout<<endl;
for(int j = Ny; j>=0; j--){
  for(int i = 0; i<Nx; i++){
    cout<<vp[i][j]<<" ";
  }
  cout<<endl;
}
cout<<endl;
cout<<"-----MAPA Vv-----"<<endl;
cout<<endl;
for(int j = Ny; j>=0; j--){
  for(int i = 0; i<Nx; i++){
    cout<<v[i][j]<<" ";
    if(j == Nx/2){
      Archivo2<<x[i][j]<<"\t"<<v[i][j]<<"\n";
    }
  }
}

```

```

    }
  }
  cout<<endl;
}
cout<<endl;
cout<<"-----MAPA P-----" <<endl;
cout<<endl;
for(int j = Ny-1; j>=0; j--){
  for(int i = 0; i<Nx; i++){
    cout<<P[i][j]<<" ";
  }
  cout<<endl;
}

ofstream Archivo3;
Archivo3.open("Mapa_velocitat.txt");
Archivo3<<"X" <<"\t" <<"Y" <<"\t" <<"Z" <<"\t" <<"u" <<"\t" <<"v" <<"\t" <<"P" <<"\t" <<"incx" <<"\t" <<"incy" <<"\n";
double U = 0;
double V = 0;
for(int j = 0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    U = 0.5*(u[i][j]+u[i+1][j]);
    V = 0.5*(v[i][j]+v[i][j+1]);
    if(i>=Na and i<=Nb-1 and j>=Nu and j<=Nv-1){ //Cuadrat
      U = 0;
      V = 0;
    }
    Archivo3<<x[i][j]<<"\t" <<y[i][j]<<"\t" <<0 <<"\t" <<U <<"\t" <<V <<"\t" <<P[i][j]<<"\t" <<incx[i][j]<<"\t" <<incy[i][j]
    ]<<"\n";
  }
}

// Code to execute
t1 = clock();

double time = (double)(t1-t0)/CLOCKS_PER_SEC;
cout << "Execution Time: " << time << endl;
ofstream Archivo6;
Archivo6.open("Execution time.txt");
Archivo6<<time;
}

```

## 7 Square Cylinder, coefficients

```

#include <iostream>
#include <fstream>
#include <math.h> // To use pow
#include <vector> // To use vectors

```

```
#include <ctime>
using namespace std;

//i>=Na and i<=Nb-1 and j>=Nu and j<=Nv-1
int main(){
    double densitat = 100;
    double temps = 200;
    double inct = 0.0001;
    double tempsmaxim = temps/inct;
    int Nx = 150;
    int Ny = 120;
    vector< vector<double> > u(Nx, vector<double> (Ny, 0));
    vector< vector<double> > v(Nx, vector<double> (Ny, 0));
    vector< vector<double> > P(Nx, vector<double> (Ny, 0));
    vector< vector<double> > incx(Nx, vector<double> (Ny, 0));
    vector< vector<double> > incy(Nx, vector<double> (Ny, 0));
    vector< vector<double> > x(Nx, vector<double> (Ny, 0));
    vector< vector<double> > y(Nx, vector<double> (Ny, 0));
    int N1 = Nx/3;
    int N2 = Nx/3;
    int N3 = Nx/3;
    int N4 = Ny/3;
    int N5 = Ny/3;
    int N6 = Ny/3;
    double La = 8;
    double Lb = 2;
    double Lc = 16;
    double Lu = 9;
    double Lv = 2;
    double Lw = 9;int Na = Nx/3 + 0.5/(Lb/N2);
    int Nb = Na + 1/(Lb/N2);
    int Nu = Ny/3 + 0.5/(Lv/N5);
    int Nv = Nu + 1/(Lv/N5);

    double cdpaverage = 0;
    double cdaverage = 0;
    double claverage = 0;
    double total = 0;

    ofstream Archivo4;
    Archivo4.open("Coeficientsglobals.txt");
    Archivo4<<"t"<<"\t"<<"Cdp"<<"\t"<<"Cd"<<"\t"<<"Cl"<<"\n";

for(int k = 0; k<1999; k++){
    ifstream Archivo;
    char nombre[500000];
    int num = (k+1)*1000;

    sprintf(nombre, "Vel_%05d.txt", num);
```



```

Archivo.open(nombre);

vector<double> col1V;
vector<double> col2V;
vector<double> col3V;
vector<double> col4V;
vector<double> col5V;
vector<double> col6V;
vector<double> col7V;
vector<double> col8V;

cout<<nombre<<endl;
Archivo.seekg(0);

while(!Archivo.eof()) {

    string col1,col2,col3,col4,col5,col6,col7,col8;
    getline(Archivo,col1,'\t'); //Separation of a tabulation
    getline(Archivo,col2,'\t'); //Separation of a tabulation
    getline(Archivo,col3,'\t'); //Separation of a tabulation
    getline(Archivo,col4,'\t'); //Separation of a tabulation
    getline(Archivo,col5,'\t'); //Separation of a tabulation
    getline(Archivo,col6,'\t'); //Separation of a tabulation
    getline(Archivo,col7,'\t'); //Separation of a tabulation
    getline(Archivo,col8,'\n'); //Separation of a new line
    if (Archivo.eof()) break;
    double a = atof(col1.c_str()); // Transforms the string to a floating point number
    double b = atof(col2.c_str());
    double c = atof(col3.c_str());
    double d = atof(col4.c_str());
    double e = atof(col5.c_str());
    double f = atof(col6.c_str());
    double g = atof(col7.c_str());
    double h = atof(col8.c_str());
    col1V.push_back(a);
    col2V.push_back(b);
    col3V.push_back(c);
    col4V.push_back(d);
    col5V.push_back(e);
    col6V.push_back(f);
    col7V.push_back(g);
    col8V.push_back(h);
    //cout << a << " " << b << endl;
}

double cont = 1;

```

```

for(int j=0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){

    x[i][j] = col1V[cont];
    y[i][j] = col2V[cont];
    u[i][j] = col4V[cont];
    v[i][j] = col5V[cont];
    P[i][j] = col6V[cont];
    incx[i][j] = col7V[cont];
    incy[i][j] = col8V[cont];
    cont++;
  }
}
Archivo.close();
//cout<<"hola"<<endl;
/*
ofstream Archivo3;
Archivo3.open("Mapa_velocitat.txt");
Archivo3<<"X"<<"Y"<<"Z"<<"u"<<"v"<<"P"<<"incx"<<"incy"<<"\n";
for(int j=0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){

    Archivo3<<x[i][j]<<"\t"<<y[i][j]<<"\t"<<0<<"\t"<<u[i][j]<<"\t"<<v[i][j]<<"\t"<<P[i][j]<<"\t"<<incx[i][j]<<"\t"
    <<incy[i][j]<<"\n";

  }
}
*/

double dragvn = 0;
double liftpn = 0;
double dragvs = 0;
double liftps = 0;
double dragpw = 0;
double liftvw = 0;
double dragpe = 0;
double liftve = 0;
for(int j=0; j<Ny; j++){
  for(int i = 0; i<Nx; i++){
    if(i>=Na and i<=Nb-1 and j == Nv){
      dragvn = dragvn + ((u[i][j]-u[i][j-1])/incy[i][j])*incx[i][j];
      liftpn = liftpn + P[i][j]*incx[i][j];
    }
    if(i>=Na and i<=Nb-1 and j == Nu-1){
      dragvs = dragvs + ((-u[i][j+1]+u[i][j])/incy[i][j])*incx[i][j];
      liftps = liftps + P[i][j]*incx[i][j];
    }
    if(j>=Nu and j<=Nv-1 and i == Na-1){

```

```

    dragpw = dragpw + P[i][j]*incy[i][j];
    liftvw = liftvw + ((v[i][j]-v[i+1][j])/incx[i][j])*incy[i][j];
  }
  if(j>=Nu and j<=Nv-1 and i == Nb){
    dragpe = dragpe + P[i][j]*incy[i][j];
    liftve = liftve + ((v[i][j]-v[i-1][j])/incx[i][j])*incy[i][j];
  }
}
}

double dragp = dragpw - dragpe;
double dragv = dragvn + dragvs;

double drag = dragvn + dragvs + dragpw - dragpe;
double lift = liftpn - liftps + liftvw + liftve;

Archivo4<<1000*inct*(k+1)<<"\t"<<fabs(dragp/(0.5*densitat))<<"\t"<<fabs(drag/(0.5*densitat))<<"\t"<<fabs(lift/(0.5*densitat))<<"\n";

cdpaverage = cdpaverage + fabs(dragp/(0.5*densitat));
cdaverage = cdaverage + fabs(drag/(0.5*densitat));
claverage = claverage + fabs(lift/(0.5*densitat));
total = k;

cout<<inct*k*1000<<endl;
}

ofstream Archivo7;
Archivo7.open("CoeficientsAverage.txt");
Archivo7<<"Cdp average : "<<cdpaverage/total<<"\n";
Archivo7<<"Cd average : "<<cdaverage/total<<"\n";
Archivo7<<"Cl average : "<<claverage/total<<"\n";

/*
ofstream Archivo2;
Archivo2.open("Coeficients.txt");
Archivo2<<"dragvn = "<<dragvn<<"\n";
Archivo2<<"liftpn = "<<liftpn<<"\n";
Archivo2<<"dragvs = "<<dragvs<<"\n";
Archivo2<<"liftps = "<<liftps<<"\n";
Archivo2<<"dragpw = "<<dragpw<<"\n";
Archivo2<<"liftvw = "<<liftvw<<"\n";
Archivo2<<"dragpe = "<<dragpe<<"\n";
Archivo2<<"liftve = "<<liftve<<"\n";
Archivo2<<"El coeficient de drag de Pressio es : "<<dragp/(0.5*densitat)<<"\n";
Archivo2<<"El coeficient de drag de Viscositat es : "<<dragv/(0.5*densitat)<<"\n";
Archivo2<<"El coeficient de drag es : "<<drag/(0.5*densitat)<<"\n";
Archivo2<<"El coeficient de lift es : "<<lift/(0.5*densitat)<<"\n";

```

```

Archivo2.close();
*/
}

```

## 8 Burgers' equation

```

#include <iostream>
#include <fstream>
#include <math.h> // To use pow
#include <vector> // To use vectors
#include <ctime>
using namespace std;

class ComplexNum{
public:
    double Re;
    double Im;

    ComplexNum(){
    }

    ComplexNum(double Re_, double Im_){
        Re = Re_;
        Im = Im_;
    }
};

static ComplexNum operator * (const ComplexNum& l, const ComplexNum& r){
    ComplexNum result;
    result.Re = (r.Re * l.Re) - (l.Im * r.Im);
    result.Im = (l.Re * r.Im) + (l.Im * r.Re);
    return result;
};

static ComplexNum operator * (const double& l, const ComplexNum& r){
    ComplexNum result;
    result.Re = r.Re * l;
    result.Im = r.Im * l;
    return result;
};

static ComplexNum operator + (const ComplexNum& l, const ComplexNum& r){
    ComplexNum result;
    result.Re = l.Re + r.Re;
    result.Im = l.Im + r.Im;
    return result;
};

```

```

};
static ComplexNum operator - (const ComplexNum& l, const ComplexNum& r){
    ComplexNum result;
    result.Re = l.Re - r.Re;
    result.Im = l.Im - r.Im;
    return result;
};

int main(){
    int N = 21;
    double Re = 40;
    double inct = 0.0001;
    double temps = 10;
    double instant = temps/inct;
    ComplexNum i(0,1);

    ComplexNum uri(1,1);
    ComplexNum uriol(3,3);
    uri = operator+(operator*(3,uriol) , uri);
    uri.Re = uri.Re;
    uri.Im = -uri.Im;
    cout<<uri.Re<<" "<<uri.Im<<endl;

    vector<ComplexNum> u(N);

    //Introduim vector velocitat

    for(int k = 0; k<N; k++){
        if(k==0){
            u[k].Re = 0;
            u[k].Im= 0;
        }
        else{
            u[k].Re = double(1)/k;
            u[k].Im= 0;
        }
    }

    //bucle temps

    for(int t = 0; t<instant; t++){
        //Calcul velocitat
        for(int k = 0; k<N; k++){
            if(k>1){

                //Terme Reynolds
                ComplexNum termeRe = operator*(-(pow(k,2)/Re),u[k]);

                //Terme sumatori
            }
        }
    }
}

```

```

ComplexNum termeSum(0,0);
ComplexNum up(0,0);
ComplexNum uq(0,0);
for(int p=-N; p<=N; p++){
  for(int q=-N; q<=N; q++){
    if(q+p==k){
      if(p<0){
        up.Re = u[-p].Re;
        up.Im = -(u[-p].Im);
      }
      if(p>=0){
        up = u[p];
      }
      if(q<0){
        uq.Re = u[-q].Re;
        uq.Im = -(u[-q].Im);
      }
      if(q>=0){
        uq = u[q];
      }
      termeSum = operator+(termeSum , (operator*(operator*(up,i),operator*(q,uq)) ));
    }
  }
}

//cout<<"("<<termeRe.Re<<" "<<termeRe.Im<<endl;
ComplexNum claudator = operator-(termeRe, termeSum);
u[k] = u[k] + operator*(inct, claudator);
}

else{
  u[k] = u[k];
}
}

//Mostrar per pantalla els vectors
ofstream Archivo1;
Archivo1.open("Energia20.txt");
Archivo1<<<"k"<<"\t"<<"E"<<"\n";
//velocitat
for(int k = 1; k<N; k++){
  ComplexNum uk(u[k].Re, -(u[k].Im));
  ComplexNum Energia = operator*(u[k],uk);
  cout<<"("<<u[k].Re<<" + "<<u[k].Im<<")"<<" Energia:"<<Energia.Re<<endl;
  Archivo1<<k<<"\t"<<Energia.Re<<"\n";
}

```

```
}  
  
}
```

## 9 Burgers' equation LES

```
#include <iostream>
#include <fstream>
#include <math.h> // To use pow
#include <vector> // To use vectors
#include <ctime>
using namespace std;

class ComplexNum{
public:
    double Re;
    double Im;

    ComplexNum(){
    }

    ComplexNum(double Re_, double Im_){
        Re = Re_;
        Im = Im_;
    }
};

static ComplexNum operator * (const ComplexNum& l, const ComplexNum& r){
    ComplexNum result;
    result.Re = (r.Re * l.Re) - (l.Im * r.Im);
    result.Im = (l.Re * r.Im) + (l.Im * r.Re);
    return result;
};

static ComplexNum operator * (const double& l, const ComplexNum& r){
    ComplexNum result;
    result.Re = r.Re * l;
    result.Im = r.Im * l;
    return result;
};

static ComplexNum operator + (const ComplexNum& l, const ComplexNum& r){
    ComplexNum result;
    result.Re = l.Re + r.Re;
    result.Im = l.Im + r.Im;
    return result;
};

static ComplexNum operator - (const ComplexNum& l, const ComplexNum& r){
```

```

ComplexNum result;
result.Re = l.Re - r.Re;
result.Im = l.Im - r.Im;
return result;
};

int main(){

int N = 21;
double Re = 40;
double inct = 0.001;
double temps = 10;
double instant = temps/inct;
ComplexNum i(0,1);

ComplexNum uri(1,1);
ComplexNum uriol(3,3);
uri = operator+(operator*(3,uriol) , uri);
uri.Re = uri.Re;
uri.Im = -uri.Im;
cout<<uri.Re<<" "<<uri.Im<<endl;

vector<ComplexNum> u(N);

//Introduim vector velocitat

for(int k = 0; k<N; k++){
  if(k==0){
    u[k].Re = 0;
    u[k].Im= 0;
  }
  else{
    u[k].Re = double(1)/k;
    u[k].Im= 0;
  }
}

//bucle temps
ComplexNum Energiaultima(0,0);
for(int t = 0; t<instant; t++){
  //Calcul velocitat
  for(int k = 0; k<N; k++){

    if(k>1){

      //Terme Reynolds
      double m = 2.0;
      double ck = 0.4523;

```



```

double vtinfnf = 0.31 * (5.0-m)/(m+1.0) * pow((3.0-m),0.5) * pow(ck,(-3.0/2.0));

if(k==N-1){
    ComplexNum uk(u[k].Re, -(u[k].Im));
    Energiaultima = operator*(u[k],uk);
}

double vtinf = (vtinfnf) * pow((Energiaultima.Re/double(N)),0.5);
double vtasterisc = 1 + 34.5*exp(-3.03*(double(N)/double(k)));
double vtk = vtinf * vtasterisc;
double viscositat = 1.0/Re + vtk;
ComplexNum termeRe = operator*(viscositat, operator*(-(pow(k,2.0)),u[k] ));
//Terme sumatori
ComplexNum termeSum(0,0);
ComplexNum up(0,0);
ComplexNum uq(0,0);
for(int p=-N; p<=N; p++){
    for(int q=-N; q<=N; q++){
        if(q+p==k){
            if(p<0){
                up.Re = u[-p].Re;
                up.Im = -(u[-p].Im);
            }
            if(p>=0){
                up = u[p];
            }
            if(q<0){
                uq.Re = u[-q].Re;
                uq.Im = -(u[-q].Im);
            }
            if(q>=0){
                uq = u[q];
            }
            termeSum = operator+(termeSum , (operator*(operator*(up,i),operator*(q,uq) )));
        }
    }
}
//cout<<" "<<termeRe.Re<<" "<<termeRe.Im<<endl;
ComplexNum claudator = operator-(termeRe, termeSum);
u[k] = u[k] + operator*(inct, claudator);
}

else{
    u[k] = u[k];
}
}
}

```

```
//Mostrar per pantalla els vectors
ofstream Archivo1;
Archivo1.open("Energia20LES.txt");
Archivo1<<"k"<<"\t"<<"E"<<"\n";
//velocitat
for(int k = 1; k<N; k++){
    ComplexNum uk(u[k].Re, -(u[k].Im));
    ComplexNum Energia = operator*(u[k],uk);
    cout<<"("<<u[k].Re<<" + "<<u[k].Im<<")"<<" Energia:"<<Energia.Re<<endl;
    Archivo1<<k<<"\t"<<Energia.Re<<"\n";
}
}
```

## 10 Lid Driven Cavity 3D

```
#include <iostream>
#include <fstream>
#include <vector>
#include <math.h>
#include <cstdlib>
#include <random>
#include "mpi.h"
using namespace std;

int main(int argc, char *argv[]){
MPI_Init(&argc, &argv); //Inicia del MPI

int rang, tamany;
    MPI_Comm_rank(MPI_COMM_WORLD, &rang); //Asigna a la variable "rang" el n de procesador
    MPI_Comm_size(MPI_COMM_WORLD, &tamany); //Asigna a la variable "tamany" la cantidad total de procesadores
    MPI_Barrier(MPI_COMM_WORLD);

int Nx = 25;
int Ny = 25;
int Nz = 25;
int Lx = 1;
int Ly = 1;
int Lz = 1;
double incx = (double)Lx/(double)Nx;
double incy = (double)Ly/(double)Ny;
double incz = (double)Lz/(double)Nz;
vector<vector<vector<double>>> P(Nx, vector<vector<double>> (Ny, vector<double>(Nz, 0)));
vector<vector<vector<double>>> up(Nx+1, vector<vector<double>> (Ny, vector<double>(Nz, 0)));
vector<vector<vector<double>>> vp(Nx, vector<vector<double>> (Ny+1, vector<double>(Nz, 0)));
vector<vector<vector<double>>> wp(Nx, vector<vector<double>> (Ny, vector<double>(Nz+1, 0)));
vector<vector<vector<double>>> u(Nx+1, vector<vector<double>> (Ny, vector<double>(Nz, 0)));
```

```

vector<vector<vector<double>>> v(Nx, vector<vector<double>> (Ny+1, vector<double>(Nz, 0)));
vector<vector<vector<double>>> w(Nx, vector<vector<double>> (Ny, vector<double>(Nz+1, 0)));
vector<vector<vector<double>>> Cdyn(Nx, vector<vector<double>> (Ny, vector<double>(Nz, 0.11)));

bool temps = true;
double inct = 0.001; //Increment de temps en segons
double tempsmaxim = 200; //Temps maxim en Segons
double tempstotal = tempsmaxim/inct;
int cont = 0; //Contador iteracions temps
double Uref = 1; //Velocitat de referencia
double mu = 1;
vector<vector<vector<double>>> muu(Nx, vector<vector<double>> (Ny, vector<double>(Nz, mu)));
double Cs = 0.11;
double lo = pow(incx*incy*incz,1.0/3.0);
double epsilon = 0.00001;
double Re = 10000;
double densitat = Re*mu/(Lx*Uref);
char nombre[500];
cout<<"El Reynolds es: " << densitat << endl;
//Condicions de frontera
for(int i = 0; i<Nx+1; i++){
  for(int j = 0; j<Ny; j++){
    for(int k = 0; k<Nz; k++){
      if(j == Ny-1){
        u[i][j][k] = Uref;
        up[i][j][k] = Uref;
      }
    }
  }
}

ofstream Archivo5;
Archivo5.open("Temps", ios::trunc);
Archivo5<<"u" <<" " <<"temps" <<"\n";
//Comena bucle temps

ofstream Arxiu;
ofstream Arxiu2;

while(temps){

//*****
//          STEP 1
//*****
//Calcul velocitat predictor u

for(int i = 1; i<Nx; i++){
  for(int j = 1; j<Ny-1; j++){

```

```

for(int k = 1; k<Nz-1; k++){
    double vn = 0.5*(v[i-1][j+1][k] + v[i][j+1][k]);
    double vs = 0.5*(v[i-1][j][k] + v[i][j][k]);
    double un = 0.5*(u[i][j+1][k] + u[i][j][k]);
    double us = 0.5*(u[i][j-1][k] + u[i][j][k]);
    double ue = 0.5*(u[i+1][j][k] + u[i][j][k]);
    double uw = 0.5*(u[i-1][j][k] + u[i][j][k]);
    double uf = 0.5*(u[i][j][k+1] + u[i][j][k]);
    double ub = 0.5*(u[i][j][k-1] + u[i][j][k]);
    double wf = 0.5*(w[i-1][j][k+1] + w[i][j][k+1]);
    double wb = 0.5*(w[i-1][j][k] + w[i][j][k]);
    double uN = u[i][j+1][k];
    double uS = u[i][j-1][k];
    double uE = u[i+1][j][k];
    double uW = u[i-1][j][k];
    double uF = u[i][j][k+1];
    double uB = u[i][j][k-1];
    double uP = u[i][j][k];
    double conv = densitat * (incx*incz*(vn*un-vs*us) + incz*incy*(ue*ue-uw*uw) + incx*incy*(wf*uf-wb*ub));
    double diff = muu[i][j][k] * (incx*incz*(uN-uP)/incz + incx*incz*(uS-uP)/incz + incy*incz*(uE-uP)/incx + incy*incz*(uW-uP)/incx + incx
*incy*(uF-uP)/incz + incx*incy*(uB-uP)/incz);
    double Ru = -conv + diff;
    up[i][j][k] =u[i][j][k] + inct/(densitat*incx*incy*incz) * Ru;
}
}
}

//Calcul velocitat predictor a v

for(int i = 1; i<Nx-1; i++){
    for(int j = 1; j<Ny; j++){
        for(int k = 1; k<Nz-1; k++){
            double vn = 0.5*(v[i][j+1][k] + v[i][j][k]);
            double vs = 0.5*(v[i][j-1][k] + v[i][j][k]);
            double ve = 0.5*(v[i+1][j][k] + v[i][j][k]);
            double vw = 0.5*(v[i-1][j][k] + v[i][j][k]);
            double uw = 0.5*(u[i][j-1][k] + u[i][j][k]);
            double ue = 0.5*(u[i+1][j-1][k] + u[i+1][j][k]);
            double vf = 0.5*(v[i][j][k+1] + v[i][j][k]);
            double vb = 0.5*(v[i][j][k-1] + v[i][j][k]);
            double wf = 0.5*(w[i][j-1][k+1] + w[i][j][k+1]);
            double wb = 0.5*(w[i][j-1][k] + w[i][j][k]);
            double vN = v[i][j+1][k];
            double vS = v[i][j-1][k];
            double vE = v[i+1][j][k];
            double vW = v[i-1][j][k];
            double vF = v[i][j][k+1];
            double vB = v[i][j][k-1];
            double vP = v[i][j][k];
        }
    }
}

```

```

double conv = densitat * (incx*incz*(vn*vn-vs*vs) + incz*incy*(ue*ve-uw*vw) + incx*incy*(wf*vf-wb*wb));
double diff = muu[i][j][k] * (incx*incz*(vN-vP)/incy + incx*incz*(vS-vP)/incy + incy*incz*(vE-vP)/incx + incy*incz*(vW-vP)/incx + incx*
incy*(vF-vP)/incz + incx*incy*(vB-vP)/incz);
double Rv = -conv + diff;
vp[i][j][k] =v[i][j][k] + inct/(densitat*incx*incy*incz) * Rv;
}
}
}

//Calcul velocitat predictor w

for(int i = 1; i<Nx-1; i++){
for(int j = 1; j<Ny-1; j++){
for(int k = 1; k<Nz; k++){
double vn = 0.5*(v[i][j+1][k] + v[i][j+1][k+1]);
double vs = 0.5*(v[i][j][k] + v[i][j][k+1]);
double uw = 0.5*(u[i][j][k] + u[i][j][k+1]);
double ue = 0.5*(u[i+1][j][k] + u[i+1][j][k+1]);
double wn = 0.5*(w[i][j][k] + w[i][j+1][k]);
double ws = 0.5*(w[i][j][k] + w[i][j-1][k]);
double we = 0.5*(w[i][j][k] + w[i+1][j][k]);
double ww = 0.5*(w[i][j][k] + w[i-1][j][k]);
double wf = 0.5*(w[i][j][k] + w[i][j][k+1]);
double wb = 0.5*(w[i][j][k] + w[i][j][k-1]);
double wN = w[i][j+1][k];
double wS = w[i][j-1][k];
double wE = w[i+1][j][k];
double wW = w[i-1][j][k];
double wF = w[i][j][k+1];
double wB = w[i][j][k-1];
double wP = w[i][j][k];
double conv = densitat * (incx*incz*(vn*wn-vs*ws) + incz*incy*(ue*we-uw*ww) + incx*incy*(wf*wf-wb*wb));
double diff = muu[i][j][k] * (incx*incz*(wN-wP)/incy + incx*incz*(wS-wP)/incy + incy*incz*(wE-wP)/incx + incy*incz*(wW-wP)/incx +
incx*incy*(wF-wP)/incz + incx*incy*(wB-wP)/incz);
double Rw = -conv + diff;
wp[i][j][k] =w[i][j][k] + inct/(densitat*incx*incy*incz) * Rw;
}
}
}

//*****
//          STEP 2
//*****

//Calcul gradient de Pressio

bool itefin = false;
while(!itefin){
itefin = true;

```

```

for(int i = 0; i<Nx; i++){
  for(int j = 0; j<Ny; j++){
    for(int k = 0; k<Nz; k++){
      double Pant = P[i][j][k];
      double Pn = 0;
      double Ps = 0;
      double Pe = 0;
      double Pw = 0;
      double Pf = 0;
      double Pb = 0;
      double vpn = vp[i][j+1][k];
      double vps = vp[i][j][k];
      double upe = up[i+1][j][k];
      double upw = up[i][j][k];
      double wpf = wp[i][j][k+1];
      double wpb = wp[i][j][k];
      double an = 0;
      double as = 0;
      double ae = 0;
      double aw = 0;
      double af = 0;
      double ab = 0;

      if(i != 0){
        aw = incy*incz/incx;
        Pw = P[i-1][j][k];
      }
      if(i != Nx-1){
        ae = incy*incz/incx;
        Pe = P[i+1][j][k];
      }
      if(j != 0){
        as = incx*incz/incy;
        Ps = P[i][j-1][k];
      }
      if(j != Ny-1){
        an = incx*incz/incy;
        Pn = P[i][j+1][k];
      }
      if(k != 0){
        ab = incy*incx/incz;
        Pb = P[i][j][k-1];
      }
      if(k != Nz-1){
        af = incy*incx/incz;
        Pf = P[i][j][k+1];
      }

      double ap = an + as + ae + aw + af + ab;
    }
  }
}

```

```

P[i][j][k] = (1.0/ap) * ((an*Pn + as*Ps + ae*Pe + aw*Pw + af*Pf + ab*Pb) - ((densitat/inct) * (incx*incz*(vpn-vps) + incy*incz*(upe-upw) +
incx*incy*(wfp-wpb) ) ) );
double error = fabs(Pant - P[i][j][k]);
if(error>epsilon) {itefin = false;}
if(P[i][j][k]>10000.0){
  cout<<"ERROR NAN"<<endl;
  return 1;
}
}
}
}

//*****
//          STEP 3
//*****

if(cont%1000==0 and cont/1000>60){
  int num = cont;
  sprintf(nombre, "Veluy_.%05d.txt", num);
  Arxiu.open(nombre,ios::trunc);}

//Calcul velocitat u
for(int i = 1; i<Nx; i++){
  for(int j = 1; j<Ny-1; j++){
    for(int k = 1; k<Nz-1; k++){
      u[i][j][k] = up[i][j][k] - inct/densitat * (P[i][j][k] - P[i-1][j][k])/incx;
      if(k == Nz/2 and j == Ny/3 and i == Nx/2 and cont%1000 == 0){
        Archivo5<<cont/1000<<" " <<u[i][j][k]<<"\n";
      }
      if(cont%1000==0 and cont/1000>60){
        if(i==Nx/2 and k==Nz/2){
          Arxiu<<incy*(j+0.5)<<" " <<0.5*(u[i][j][k]+u[i+1][j][k])<<"\n";}
        }
      }
    }
  }
}
Arxiu.close();

//Calcul velocitat v

if(cont%1000==0 and cont/1000>190){
  int num = cont;
  sprintf(nombre, "Velvx_.%05d.txt", num);
  Arxiu2.open(nombre,ios::trunc);}

for(int i = 1; i<Nx-1; i++){
  for(int j = 1; j<Ny; j++){
    for(int k = 1; k<Nz-1; k++){

```

```

v[i][j][k] = vp[i][j][k] - inct/densitat * (P[i][j][k] - P[i][j-1][k])/incy;
if(cont%1000==0 and cont/1000>60){
if(j==Ny/2 and k==Nz/2){
Arxiu2<<incx*(i+0.5)<<" " <<0.5*(v[i][j][k]+v[i][j+1][k])<<"\n";}
}
}
}
}
Arxiu2.close();

//Calcul velocitat w

for(int i = 1; i<Nx-1; i++){
for(int j = 1; j<Ny-1; j++){
for(int k = 1; k<Nz; k++){
w[i][j][k] = wp[i][j][k] - inct/densitat * (P[i][j][k] - P[i][j][k-1])/incz;
}
}
}

//*****
//          STEP 4
//*****

//Calcul viscositat sub grid scale
for(int i = 1; i<Nx-1; i++){
for(int j = 1; j<Ny-1; j++){
for(int k = 1; k<Nz-1; k++){
double dwdx = (v[i][j][k]-v[i-1][j][k])/incx;
double dudy = (u[i][j][k]-u[i][j-1][k])/incy;
double dvdz = (v[i][j][k]-v[i][j][k-1])/incz;
double dwdy = (w[i][j][k]-w[i][j-1][k])/incy;
double dudz = (u[i][j][k]-u[i][j][k-1])/incz;
double dwdx = (w[i][j][k]-w[i-1][j][k])/incx;
double dudx = (u[i][j][k]-u[i-1][j][k])/incx;
double dvdy = (v[i][j][k]-v[i][j-1][k])/incy;
double dwdz = (w[i][j][k]-w[i][j][k-1])/incz;
double Sij =pow((2*( 0.5*pow(dwdx+dudy,2) +0.5*pow(dudz+dwdx,2) +0.5*pow(dvdz+dwdy,2) + pow(dudx,2) + pow(dvdy,2) + pow(dwdz,2)
)), 0.5);
//double vis = pow((Cs*lo),2) * Sij;
double vis = pow((Cdyn[i][j][k]*lo),2) * Sij;
muu[i][j][k] = mu + vis*densitat;
}
}
}

//*****
//          STEP 5

```



```

//*****
//Calcul coeficient dynamic
for(int i = 1; i<Nx-1; i++){
  for(int j = 1; j<Ny-1; j++){
    for(int k = 1; k<Nz-1; k++){
      //-----Prova L-----
      double L[3][3];
      double L1[3][3];
      double L2[3][3];
      //L1
      L1[0][0] = (2*pow(u[i][j][k],2) + pow(u[i-1][j][k],2) + pow(u[i+1][j][k],2) ) /4;
      L1[1][1] = (2*pow(v[i][j][k],2) + pow(v[i][j-1][k],2) + pow(v[i][j+1][k],2) ) /4;
      L1[2][2] = (2*pow(w[i][j][k],2) + pow(w[i][j][k-1],2) + pow(w[i][j][k+1],2) ) /4;
      L1[0][1] = (2*(u[i][j][k]*v[i][j][k]) + u[i-1][j][k]*v[i][j-1][k] + u[i+1][j][k]*v[i][j+1][k]) / 4;
      L1[1][0] = (2*(u[i][j][k]*v[i][j][k]) + u[i-1][j][k]*v[i][j-1][k] + u[i+1][j][k]*v[i][j+1][k]) / 4;
      L1[0][2] = (2*(u[i][j][k]*w[i][j][k]) + u[i-1][j][k]*w[i][j][k-1] + u[i+1][j][k]*w[i][j][k+1]) / 4;
      L1[2][0] = (2*(u[i][j][k]*w[i][j][k]) + u[i-1][j][k]*w[i][j][k-1] + u[i+1][j][k]*w[i][j][k+1]) / 4;
      L1[2][1] = (2*(v[i][j][k]*w[i][j][k]) + v[i][j-1][k]*w[i][j][k-1] + v[i][j+1][k]*w[i][j][k+1]) / 4;
      L1[1][2] = (2*(v[i][j][k]*w[i][j][k]) + v[i][j-1][k]*w[i][j][k-1] + v[i][j+1][k]*w[i][j][k+1]) / 4;
      //L2
      L2[0][0] =pow( ( (2*u[i][j][k] + u[i+1][j][k] + u[i-1][j][k]) / 4 ) ,2);
      L2[1][1] =pow( ( (2*v[i][j][k] + v[i][j+1][k] + v[i][j-1][k]) / 4 ) ,2);
      L2[2][2] =pow( ( (2*w[i][j][k] + w[i][j][k+1] + w[i][j][k-1]) / 4 ) ,2);
      L2[0][1] = ( (2*v[i][j][k] + v[i][j+1][k] + v[i][j-1][k]) / 4 ) * ( (2*u[i][j][k] + u[i+1][j][k] + u[i-1][j][k]) / 4 );
      L2[1][0] = ( (2*v[i][j][k] + v[i][j+1][k] + v[i][j-1][k]) / 4 ) * ( (2*u[i][j][k] + u[i+1][j][k] + u[i-1][j][k]) / 4 );
      L2[0][2] = ( (2*w[i][j][k] + w[i][j][k+1] + w[i][j][k-1]) / 4 ) * ( (2*u[i][j][k] + u[i+1][j][k] + u[i-1][j][k]) / 4 );
      L2[2][0] = ( (2*w[i][j][k] + w[i][j][k+1] + w[i][j][k-1]) / 4 ) * ( (2*u[i][j][k] + u[i+1][j][k] + u[i-1][j][k]) / 4 );
      L2[1][2] = ( (2*w[i][j][k] + w[i][j][k+1] + w[i][j][k-1]) / 4 ) * ( (2*v[i][j][k] + v[i][j+1][k] + v[i][j-1][k]) / 4 );
      L2[2][1] = ( (2*w[i][j][k] + w[i][j][k+1] + w[i][j][k-1]) / 4 ) * ( (2*v[i][j][k] + v[i][j+1][k] + v[i][j-1][k]) / 4 );
      //Lfinal
      L[0][0] = L1[0][0] - L2[0][0];
      L[0][1] = L1[0][1] - L2[0][1];
      L[0][2] = L1[0][2] - L2[0][2];
      L[1][0] = L1[1][0] - L2[1][0];
      L[1][1] = L1[1][1] - L2[1][1];
      L[1][2] = L1[1][2] - L2[1][2];
      L[2][0] = L1[2][0] - L2[2][0];
      L[2][1] = L1[2][1] - L2[2][1];
      L[2][2] = L1[2][2] - L2[2][2];
      for(int i = 0; i<3; i++){
        for(int j = 0; j<3; j++){
          if(isnan(L[i][j])){
            cout<<"ERROR NaN L"<<endl;
            return 0;
          }
        }
      }
    }
  }
}
//-----Prova M-----

```

```

double M[3][3];
double Sij[3][3];
double Sijdreta[3][3];
double Sijesquerra[3][3];
    //Sij i modul S
double dvdx = (v[i][j][k]-v[i-1][j][k])/incx;
double dudy = (u[i][j][k]-u[i][j-1][k])/incy;
double dvdz = (v[i][j][k]-v[i][j][k-1])/incz;
double dwdy = (w[i][j][k]-w[i][j-1][k])/incy;
double dudz = (u[i][j][k]-u[i][j][k-1])/incz;
double dwdx = (w[i][j][k]-w[i-1][j][k])/incx;
double dudx = (u[i][j][k]-u[i-1][j][k])/incx;
double dvdy = (v[i][j][k]-v[i][j-1][k])/incy;
double dwdz = (w[i][j][k]-w[i][j][k-1])/incz;
Sij[0][0] = dudx;
Sij[1][1] = dvdy;
Sij[2][2] = dwdz;
Sij[0][1] = 0.5*(dvdx + dudy);
Sij[1][0] = 0.5*(dvdx + dudy);
Sij[0][2] = 0.5*(dwdx + dudz);
Sij[2][0] = 0.5*(dwdx + dudz);
Sij[2][1] = 0.5*(dwdy + dvdz);
Sij[1][2] = 0.5*(dwdy + dvdz);
double modulS = pow((2*( 0.5*pow(dvdx+dudy,2) +0.5*pow(dudz+dwdx,2) +0.5*pow(dvdz+dwdy,2) + pow(dudx,2) + pow(dvdy,2) + pow(
dwdz,2) )), 0.5);
    //Sij dreta i modul S dreta
dvdx = (v[i][j+1][k]-v[i-1][j+1][k])/incx;
dudy = (u[i+1][j][k]-u[i+1][j-1][k])/incy;
dvdz = (v[i][j+1][k]-v[i][j+1][k-1])/incz;
dwdy = (w[i][j][k+1]-w[i][j-1][k+1])/incy;
dudz = (u[i+1][j][k]-u[i+1][j][k-1])/incz;
dwdx = (w[i][j][k+1]-w[i-1][j][k+1])/incx;
dudx = (u[i+1][j][k]-u[i][j][k])/incx;
dvdy = (v[i][j+1][k]-v[i][j][k])/incy;
dwdz = (w[i][j][k+1]-w[i][j][k])/incz;
Sijdreta[0][0] = dudx;
Sijdreta[1][1] = dvdy;
Sijdreta[2][2] = dwdz;
Sijdreta[0][1] = 0.5*(dvdx + dudy);
Sijdreta[1][0] = 0.5*(dvdx + dudy);
Sijdreta[0][2] = 0.5*(dwdx + dudz);
Sijdreta[2][0] = 0.5*(dwdx + dudz);
Sijdreta[2][1] = 0.5*(dwdy + dvdz);
Sijdreta[1][2] = 0.5*(dwdy + dvdz);
double modulSdreta = pow((2*( 0.5*pow(dvdx+dudy,2) +0.5*pow(dudz+dwdx,2) +0.5*pow(dvdz+dwdy,2) + pow(dudx,2) + pow(dvdy,2) +
pow(dwdz,2) )), 0.5);
    //Sij esquerra i modul S esquerra
dvdx = (v[i][j-1][k]-v[i-1][j-1][k])/incx;
dudy = (u[i-1][j][k]-u[i-1][j-1][k])/incy;

```

```

dvdz = (v[i][j-1][k]-v[i][j-1][k-1])/incz;
dwdy = (w[i][j][k-1]-w[i][j-1][k-1])/incy;
dudz = (u[i-1][j][k]-u[i-1][j][k-1])/incz;
dwdx = (w[i][j][k-1]-w[i-1][j][k-1])/incx;
dudx = (u[i-1][j][k]-u[i][j][k])/incx;
dvdy = (v[i][j-1][k]-v[i][j][k])/incy;
dwdz = (w[i][j][k-1]-w[i][j][k])/incz;
Sijesquerra[0][0] = dudx;
Sijesquerra[1][1] = dvdy;
Sijesquerra[2][2] = dwdz;
Sijesquerra[0][1] = 0.5*(dvdz + dudy);
Sijesquerra[1][0] = 0.5*(dvdz + dudy);
Sijesquerra[0][2] = 0.5*(dwdx + dudz);
Sijesquerra[2][0] = 0.5*(dwdx + dudz);
Sijesquerra[2][1] = 0.5*(dwdy + dvdz);
Sijesquerra[1][2] = 0.5*(dwdy + dvdz);
double modulSesquerra = pow((2*( 0.5*pow(dvdz+dudy,2) +0.5*pow(dudz+dwdx,2) +0.5*pow(dvdz+dwdy,2) + pow(dudx,2) + pow(dvdy
,2) + pow(dwdz,2) ) ) , 0.5);
double tamanymalla = pow(incx, 2);
double Sij1[3][3];
double Sij2[3][3];
//Mij 1
for(int a = 0; a<3; a++){
  for(int b = 0; b<3;b++){
    Sij1[a][b] = ( (2*Sij[a][b] + Sijesquerra[a][b] + Sijdreta[a][b] ) /4) * ((2*modulS + modulSdreta + modulSesquerra) / 4) * 2 *
tamanymalla;
    if(isnan(Sij1[a][b])){
      cout<<"ERROR NaN Sij1"<<endl;
      return 0;
    }
  }
}
//Mij 2
for(int a = 0; a<3; a++){
  for(int b = 0; b<3;b++){
    Sij2[a][b] = ( (2*Sij[a][b]*modulS + Sijesquerra[a][b]*modulSesquerra + Sijdreta[a][b]*modulSdreta ) /4) * tamanymalla;
    if(isnan(Sij2[a][b])){
      cout<<"ERROR NaN Sij2"<<endl;
      return 0;
    }
  }
}
//Mij final
for(int a = 0; a<3; a++){
  for(int b = 0; b<3;b++){
    M[a][b] = Sij1[a][b] - Sij2[a][b];
    if(isnan(M[a][b])){
      cout<<"ERROR NaN M"<<endl;
      return 0;
    }
  }
}

```

```

    }
  }
}
//Calcul Coeficient Dinamic
double LM = L[0][0]*M[0][0] + L[0][1]*M[1][0] + L[0][2]*M[2][0] + L[1][0]*M[0][1] + L[1][1]*M[1][1] + L[1][2]*M[2][1] + L[2][0]*M
[0][2] + L[2][1]*M[1][2] + L[2][2]*M[2][2];
if(isnan(LM)){
  cout<<"ERROR NaN LM"<<endl;
  return 0;
}
//cout<<"LM es: "<<LM<<endl;
double MM = M[0][0]*M[0][0] + M[0][1]*M[1][0] + M[0][2]*M[2][0] + M[1][0]*M[0][1] + M[1][1]*M[1][1] + M[1][2]*M[2][1] + M
[2][0]*M[0][2] + M[2][1]*M[1][2] + M[2][2]*M[2][2];
if(isnan(MM)){
  cout<<"ERROR NaN MM"<<endl;
  return 0;
}
//cout<<"MM es: "<<MM<<endl;

double coefdyn =sqrt(-0.5*(LM/MM));
if(coefdyn < 0){
  coefdyn = 0;
}
if(isnan(coefdyn)){
  coefdyn = 0.11;
}
if(coefdyn > 0.11){
  coefdyn = 0.11;
}
Cdyn[i][j][k] = coefdyn;
//cout<<Cdyn[i][j][k]<<endl;
if(isnan(Cdyn[i][j][k])){
  cout<<"ERROR NaN Cdyn"<<endl;
  return 0;
}
//cout<<Cdyn[i][j][k]<<endl;
}
}
}
}

```

```

//Acaben steps
cont++;
if(cont%1000 == 0){
cout<<cont/1000<<endl;
}
if((double)cont >= tempstotal) temps = false;
} //Acaba bucle temps
Archivo5.close();

//*****
//Introduim datta al fitxer

ofstream Archivo;
Archivo.open("Datta", ios::trunc);
Archivo<<"X"<<" "<<"Y"<<" "<<"Z"<<" "<<"Pressio"<<" "<<"viscositat"<<" "<<"u"<<" "<<"v"<<" "<<"w"<<"\n";
for(int i = 0; i<Nx; i++){
  for(int j = 0; j<Ny; j++){
    for(int k = 0; k<Nz; k++){
      Archivo<<incx*(i+0.5)<<" "<<incy*(j+0.5)<<" "<<incz*(k+0.5)<<" "<<P[i][j][k]<<" "<<muu[i][j][k]<<" "<<0.5*(u[i][j][k]+
      u[i+1][j][k])<<" "<<0.5*(v[i][j][k]+v[i][j+1][k])<<" "<<0.5*(w[i][j][k]+w[i][j][k+1])<<"\n";
    }
  }
}
Archivo.close();

//Introduim velocitat u en el llarg d'y en el pla mitj
ofstream Archivo2;
Archivo2.open("UY", ios::trunc);
//Archivo2<<"y"<<" "<<"u"<<"\n";
for(int i = 0; i<Nx; i++){
  for(int j = 0; j<Ny; j++){
    for(int k = 0; k<Nz; k++){
      if(i==Nx/2 and k==Nz/2){
        Archivo2<<incx*(j+0.5)<<" "<<0.5*(u[i][j][k]+u[i+1][j][k])<<"\n";
      }
    }
  }
}
Archivo2.close();

//Introduim velocitat v en el llarg d'x en el pla mitj
ofstream Archivo3;
Archivo3.open("VX", ios::trunc);

```

```

//Archivo3<<"x"<<" "<<"v"<<"\n";
for(int i = 0; i<Nx; i++){
  for(int j = 0; j<Ny; j++){
    for(int k = 0; k<Nz; k++){
      if(j==Ny/2 and k==Nz/2){
        Archivo3<<incx*(i+0.5)<<" "<<0.5*(v[i][j][k]+v[i][j+1][k])<<"\n";
      }
    }
  }
}
Archivo3.close();

//Introduim viscositat sub grid scale al llarg del pla mitj

ofstream Archivo4;
Archivo4.open("Viscositat", ios::trunc);
Archivo4<<"x"<<" "<<"y"<<" "<<"z"<<" "<<"viscositat"<<"\n";
for(int i = 0; i<Nx; i++){
  for(int j = 0; j<Ny; j++){
    for(int k = 0; k<Nz; k++){
      if(k==Nz/2){
        Archivo4<<incx*(i+0.5)<<" "<<incy*(j+0.5)<<" "<<0<<" "<<muu[i][j][k]-mu<<"\n";
      }
    }
  }
}
Archivo4.close();

ofstream Archivo6;
Archivo6.open("CoeficientDynamic", ios::trunc);
Archivo6<<"x"<<" "<<"y"<<" "<<"z"<<" "<<"viscositat"<<"\n";
for(int i = 0; i<Nx; i++){
  for(int j = 0; j<Ny; j++){
    for(int k = 0; k<Nz; k++){
      if(k==Nz/2){
        Archivo6<<incx*(i+0.5)<<" "<<incy*(j+0.5)<<" "<<0<<" "<<Cdyn[i][j][k]<<"\n";
      }
    }
  }
}
Archivo6.close();

MPI_Finalize(); // Finaliza el MPI
}

```

## 11 Parallelization Monte Carlo method

```
#include <iostream>
```

```

#include <fstream>
#include <vector>
#include <math.h>
#include <cstdlib>
#include <random>
#include "mpi.h"
using namespace std;

int main(int argc, char *argv[]) {
MPI_Init(&argc, &argv); //Inicia el MPI

//cout<<"Holav3"<<endl;
int rang, tamany;
  MPI_Comm_rank(MPI_COMM_WORLD, &rang); //Asigna a la variable "rang" el n de procesador
  MPI_Comm_size(MPI_COMM_WORLD, &tamany); //Asigna a la variable "tamany" la cantidad total de procesadores
  MPI_Barrier(MPI_COMM_WORLD);
//  cout<<"Holav3! Soy el core : "<<rang<<" de "<<tamany<<endl;

int n = 1000000;
double randomx;
double randomy;
double cont = 0;

std::random_device rd; //Will be used to obtain a seed for the random number engine
std::mt19937 gen(rd()); //Standard mersenne_twister_engine seeded with rd()
std::uniform_real_distribution<double> unif(-1, 1);

  srand(time(NULL) + rang);
for(int i=0; i<n; i++){
//  randomx = -1+2*((double)rand())/((double)RAND_MAX);
  randomx = unif(gen);
//  cout<<randomx<<endl;
//  randomy = -1+2*((double)rand())/((double)RAND_MAX);
  randomy = unif(gen);
//  cout<<randomy<<endl;
  if((pow(randomx,2)+pow(randomy,2)) <= 1) cont++;
}
double pi = (cont/double(n)) * 4;
cout<<"Pel core "<<rang<<" Pi es: "<<pi<<endl;

double piaverage;
MPI_Allreduce(&pi, &piaverage, 1, MPI::DOUBLE, MPI_SUM, MPI_COMM_WORLD);
piaverage = piaverage/tamany;

cout<<"El valor de pi average es: "<<piaverage<<endl;

MPI_Finalize(); //Finaliza el MPI

```

```
}

```

## 12 Parallelization 1D bar

```
#include <iostream>
#include <fstream>
#include <vector>
#include <math.h>
#include <cstdlib>
#include <random>
#include "mpi.h"
using namespace std;

int main(int argc, char *argv[]) {
MPI_Init(&argc, &argv); //Inicia el MPI

int rang, tamany;
    MPI_Comm_rank(MPI_COMM_WORLD, &rang); //Asigna a la variable "rang" el n de procesador
    MPI_Comm_size(MPI_COMM_WORLD, &tamany); //Asigna a la variable "tamany" la cantidad total de procesadores
    MPI_Barrier(MPI_COMM_WORLD);

//int n = 5*tamany;
int n = 2400;
if(n%tamany != 0) {
    cout<<"Error numero de cores"<<endl;
    return 1;
}
vector<double> T(n/tamany, 0.0);
vector<double> Tant(n/tamany, 0.0);
double L = 1;
double incx = L/double(n);
double epsilon = 1E-8;
bool minim = false;
int iter = 0;
double error = 0;
double qv = -10;
if(rang == 0) T[0] = 1;

//ITERACIONS
double dreta = 0;
double esquerra = 0;
double number;

while(!minim){
    minim = true;
if(tamany > 1) {
    if(rang == 0){
        MPI_Send(&T[n/tamany-1], 1, MPI_DOUBLE, 1, 0, MPI_COMM_WORLD);
```



```

    MPI_Recv(&dreta, 1, MPI.DOUBLE, 1, 0, MPI.COMM_WORLD, MPI.STATUS_IGNORE);
} if(rang>0 and rang<tamany-1){
    MPI_Send(&T[0], 1, MPI.DOUBLE, (rang-1), 0, MPI.COMM_WORLD);
    MPI_Recv(&esquerra, 1, MPI.DOUBLE, (rang-1), 0, MPI.COMM_WORLD, MPI.STATUS_IGNORE);
    MPI_Send(&T[n/tamany-1], 1, MPI.DOUBLE, (rang+1), 0, MPI.COMM_WORLD);
    MPI_Recv(&dreta, 1, MPI.DOUBLE, (rang+1), 0, MPI.COMM_WORLD, MPI.STATUS_IGNORE);
} if(rang == tamany-1){
    MPI_Send(&T[0], 1, MPI.DOUBLE, rang-1, 0, MPI.COMM_WORLD);
    MPI_Recv(&esquerra, 1, MPI.DOUBLE, rang-1, 0, MPI.COMM_WORLD, MPI.STATUS_IGNORE);
}
}
double suma = 0;
for(int i = 0; i<n/tamany; i++){
    if(i>0 and i<n/tamany-1){
        T[i] = (T[i+1]+T[i-1]-qv*pow(incx,2))/2;
    } else{
        //CORE 0
        if(rang == 0){
            if(i == 0){
                T[i] = 1;
            }
            else if(i == (n/tamany)-1){
                if(tamany == 1) dreta = 0;
                T[i] = (dreta+T[i-1]-qv*pow(incx,2))/2;
            }
            //CORES CENTRE
        } else if (rang > 0 and rang < tamany-1){
            if(i == 0){
                T[i] = (T[i+1]+esquerra-qv*pow(incx,2))/2;
            }
            else if(i == (n/tamany)-1){
                T[i] = (dreta+T[i-1]-qv*pow(incx,2))/2;
            }
            //CORE FINAL
        } else if (rang == tamany-1){
            if(i == 0){
                T[i] = (T[i+1]+esquerra-qv*pow(incx,2))/2;
            }
            else if(i == (n/tamany)-1){
                T[i] = 0;
            }
        }
    }
}
error = fabs(T[i]-Tant[i]);
//MPI_Barrier(MPI.COMM_WORLD);
if(error > epsilon){ suma = 1;}
Tant[i] = T[i];
}
iter++;

```

```

double sumatotal = 0;
MPI_Allreduce(&suma, &sumatotal, 1, MPI::DOUBLE, MPI_SUM, MPI_COMM_WORLD);
if(sumatotal > 0) {minim = false;}
}

//VECTOR FINAL
vector<double> Tfinal(n, 0.0);
MPI_Gather(&T[0], T.size(), MPI_DOUBLE, &Tfinal[0], T.size(), MPI_DOUBLE, 0, MPI_COMM_WORLD);

cout<<"El vector de temperatura del core: "<<rang<<" es: ";
for(int i = 0; i<n/tamany; i++){
    cout<<T[i]<<" ";
}
cout<<endl;

MPI_Barrier(MPI_COMM_WORLD);
if (rang == 0){
ofstream Archivo;
Archivo.open("Temperatura", ios::trunc);
for(int i = 0; i<n; i++){
    cout<<Tfinal[i]<<" ";
    Archivo << i*incx + 0.5*incx <<" " << Tfinal[i] << endl;
}cout<<endl;
cout<<"S'han fet: "<<iter<<" iteracions"<<endl;
Archivo.close();
}

MPI_Finalize(); //Finaliza el MPI
}

```