

MEEN 644 - Homework 3

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Problem statement

Consider a thin copper square plate of dimensions $0.5 \text{ m} \times 0.5 \text{ m}$. The temperature of the west and south edges are maintained at 50°C and the north edge is maintained at 100°C . The east edge is insulated. Using finite volume method, write a program to predict the steady-state temperature solution.

- (a) **(35 points)** Set the over relaxation factor α from 1.00 to 1.40 in steps of 0.05 to identify α_{opt} . Plot the number of iterations required for convergence for each α .
- (b) **(15 points)** Solve the same problem using $21^2, 25^2, 31^2$, and 41^2 CVs, respectively. Plot the temperature at the center of the plate (0.25 m, 0.25 m) vs CVs.
- (c) **(10 points)** Plot the steady state temperature contour in the 2D domain with the 41^2 CV solution.

Preliminaries

Two-dimensional heat conduction

With two-dimensional heat conduction with constant material properties, insulation on the right and prescribed temperatures on all other sides, we have the PDE

$$\begin{cases} k \frac{\partial^2 T}{\partial x^2} + k \frac{\partial^2 T}{\partial y^2} = 0, \\ T(x, 0) = T_B, \\ T(0, y) = T_L, \\ T(0, L) = T_T, \\ -k \frac{\partial T}{\partial x} \Big|_{x=L} = 0, \end{cases} \quad (1)$$

where

$$\begin{aligned} T_B &\equiv 50^\circ\text{C}, & T_L &\equiv 50^\circ\text{C}, & T_T &\equiv 100^\circ\text{C}, \\ k &\equiv 386 \text{ W/m }^\circ\text{C}, & L &\equiv 0.5 \text{ m}. \end{aligned}$$

We discretize the region on $x \times y = [0, L]^2$ by N^2 internal nodes with $\Delta x = x/N, \Delta y = y/N$.

Control volume equations

Integrate over an internal control volume (i, j) to obtain

$$k \iint_{CV_{i,j}} \left[\frac{\partial^2 T}{\partial x^2} + \frac{\partial^2 T}{\partial y^2} \right] dx dy = 0, \quad (i, j) \in [2, 3, \dots, N-1]^2.$$

$$k \Delta y \left[\frac{\partial T}{\partial x} \Big|_{w_{i,j}} - \frac{\partial T}{\partial x} \Big|_{e_{i,j}} \right] + k \Delta x \left[\frac{\partial T}{\partial y} \Big|_{n_{i,j}} - \frac{\partial T}{\partial y} \Big|_{s_{i,j}} \right] = 0, \quad (i, j) \in [2, 3, \dots, N-1]^2.$$

Now use the two node formulation for the derivative terms to obtain

$$k \Delta y \left[\frac{T_{E_{ij}} - T_{P_{ij}}}{\Delta x} - \frac{T_{P_{ij}} - T_{W_{ij}}}{\Delta x} \right] + k \Delta x \left[\frac{T_{N_{ij}} - T_{P_{ij}}}{\Delta y} - \frac{T_{P_{ij}} - T_{S_{ij}}}{\Delta y} \right] = 0, \quad (i, j) \in [2, 3, \dots, N-1]^2.$$

Collect like terms and modify the index to obtain

$$T_{i,j} a_p - T_{i,j+1} a_n - T_{i+1,j} a_e - T_{i,j-1} a_s - T_{i-1,j} a_w = b, \quad (i, j) \in [2, 3, \dots, N-1]^2, \quad (2)$$

where (note that the below applies only to *internal* control volumes)

$$a_n \equiv \frac{k \Delta y}{\Delta x}, \quad a_e \equiv \frac{k \Delta x}{\Delta y}, \quad a_s \equiv \frac{k \Delta y}{\Delta x}, \quad a_w \equiv \frac{k \Delta x}{\Delta y}, \quad a_p \equiv a_n + a_e + a_s + a_w, \quad \text{and} \quad b = 0.$$

The remaining equations are solved similarly but with slight differences depending on which boundary the CV is on. For the boundary control volumes, we first assume the above are the coefficients with the equation given in Equation (2). Next, the following changes are made; if the control volume is on the:

- North surface, then $a_n = 2k \Delta y / \Delta x$ and $T_{i,N+1} = T_T$ (known, moves to RHS),
- East surface, then, $a_e = 0$ and $T_{N+1,j} = 0$,
- South surface, then, $a_s = 2k \Delta y / \Delta x$ and $T_{i,0} = T_B$ (known, moves to RHS),
- West surface, then, $a_w = 2k \Delta x / \Delta y$ and $T_{0,j} = T_L$ (known, moves to RHS).

Solving method

The problem is to be solved by the line-by-line method. In specific, the sweeping arrangement is: **south to north, west to east, north to south, east to west**. In this method, the contribution from one direction in a given control volume is lagged and moved to the right hand side in order to solve a tri-diagonal system. For example, in a south to north sweep, a single row in the physical mesh is solved (integrated) at a time. The terms that are on the left hand side that come from CVs above and below the given row are lagged (not solved for, the most up-to-date solution is known) in order to move said terms to the right hand side. The resulting system of equations for a single row is a diagonally dominant tri-diagonal matrix, which allows for a quick solve. In the iteration process, the initial guess is the average of the Dirichlet boundary values, $(T_B + T_L + T_T)/3$.

Convergence is declared when

$$R = \sum_{CV} \left| a_p T_p - \sum_{nb} a_{nb} T_{nb} b_p \right| \leq 10^{-5}. \quad (3)$$

Upon solving an individual system $Ax^{\ell+1} = b$ with relaxation (where ℓ is the iteration index, i.e., x^ℓ is the solution from the previous iteration), the system is relaxed with the coefficient α by modifying it after construction by

$$\begin{cases} a_{ii} = a_{ii}\alpha^{-1}, \\ b_i = b_i + (\alpha^{-1} - 1)a_{ii}x_i^\ell, \end{cases} \quad i = 1, \dots, N,$$

and it is then solved using the standard TDMA algorithm.

Results

Part a

With the given range of α , it was determined for this specific problem with 15^2 CVs that $\alpha_{\text{opt}} \approx 1.3$ (given that it required the fewest iterations). The solve did not converge (with an attempted 2,000 iterations) for $\alpha = 1.4$. The requested figure showing the residuals follows in Figure 1.

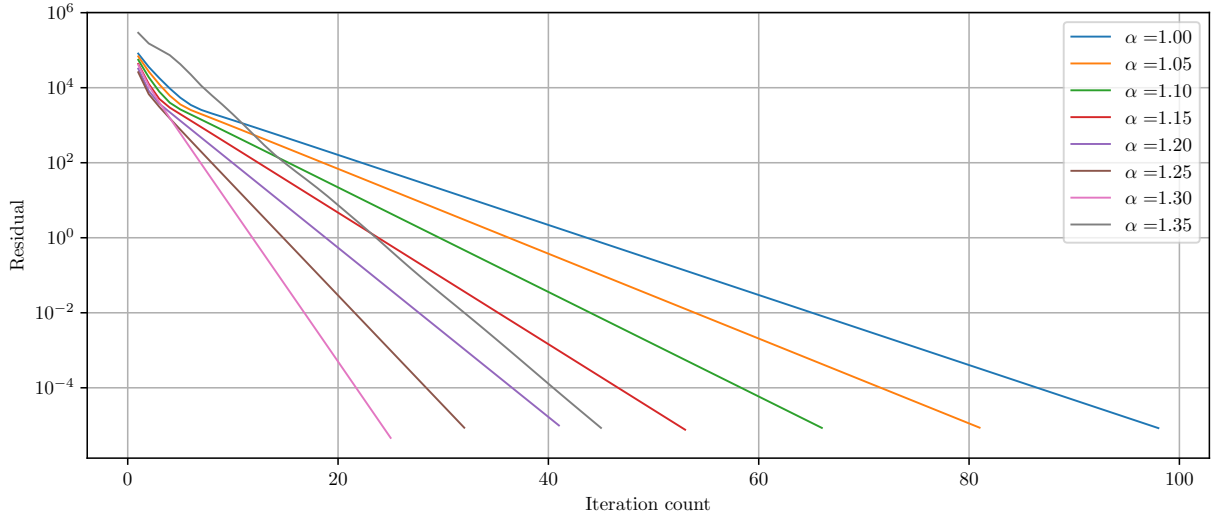


Figure 1: Plot of the residual vs. iteration count for each relaxation parameter, α .

Part b

With a mesh refinement of 21^2 , 25^2 , 31^2 , and 41^2 CVs, the center temperature for each refinement with varying relaxation parameter α is plotted below in Figure 2. The same result is tabulated in Table 1. Note the fact that to 6 digits the solution does not change for a given refinement with a change in α . This is due to the fact that (providing the scheme converges to within the desired tolerance) relaxation will not change the solution significantly (to within the tolerance).

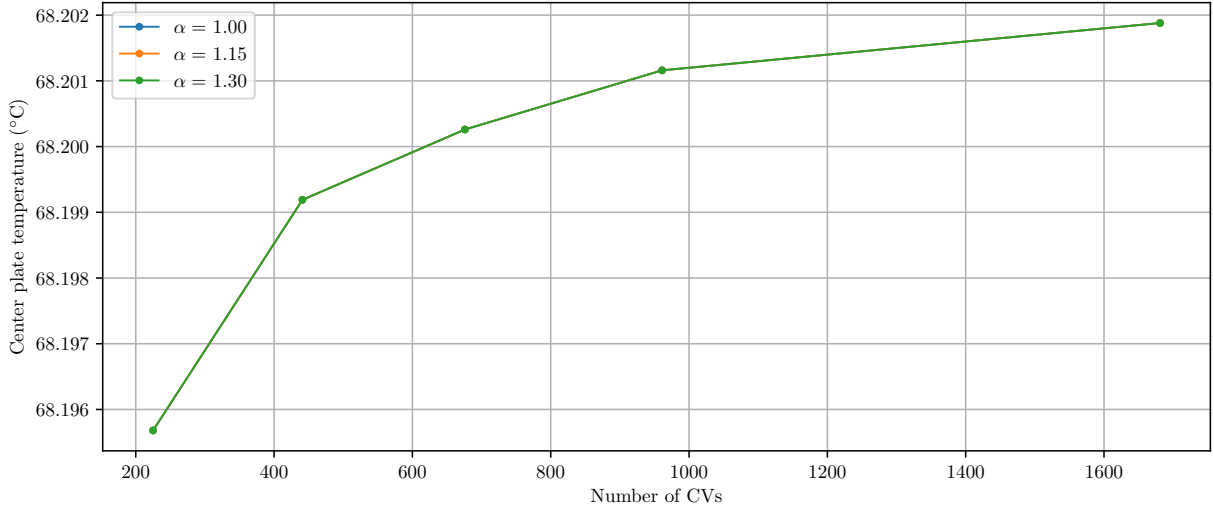


Figure 2: Plot of the center temperature with mesh refinement.

Table 1: The center temperature (°C) with varying refinements and relaxation parameters.

CVs	$\alpha = 1.0$	$\alpha = 1.15$	$\alpha = 1.30$
225	68.19568	68.19568	68.19568
441	68.19919	68.19919	68.19919
625	68.20026	68.20026	68.20026
961	68.20116	68.20116	68.20116
1681	68.20188	68.20188	68.20188

With refinement, the residuals were plotted below in Figure 3. No refinement levels for 21^2 CVs and above converged for $\alpha = 1.35, 1.4$ with 2,000 maximum iterations.

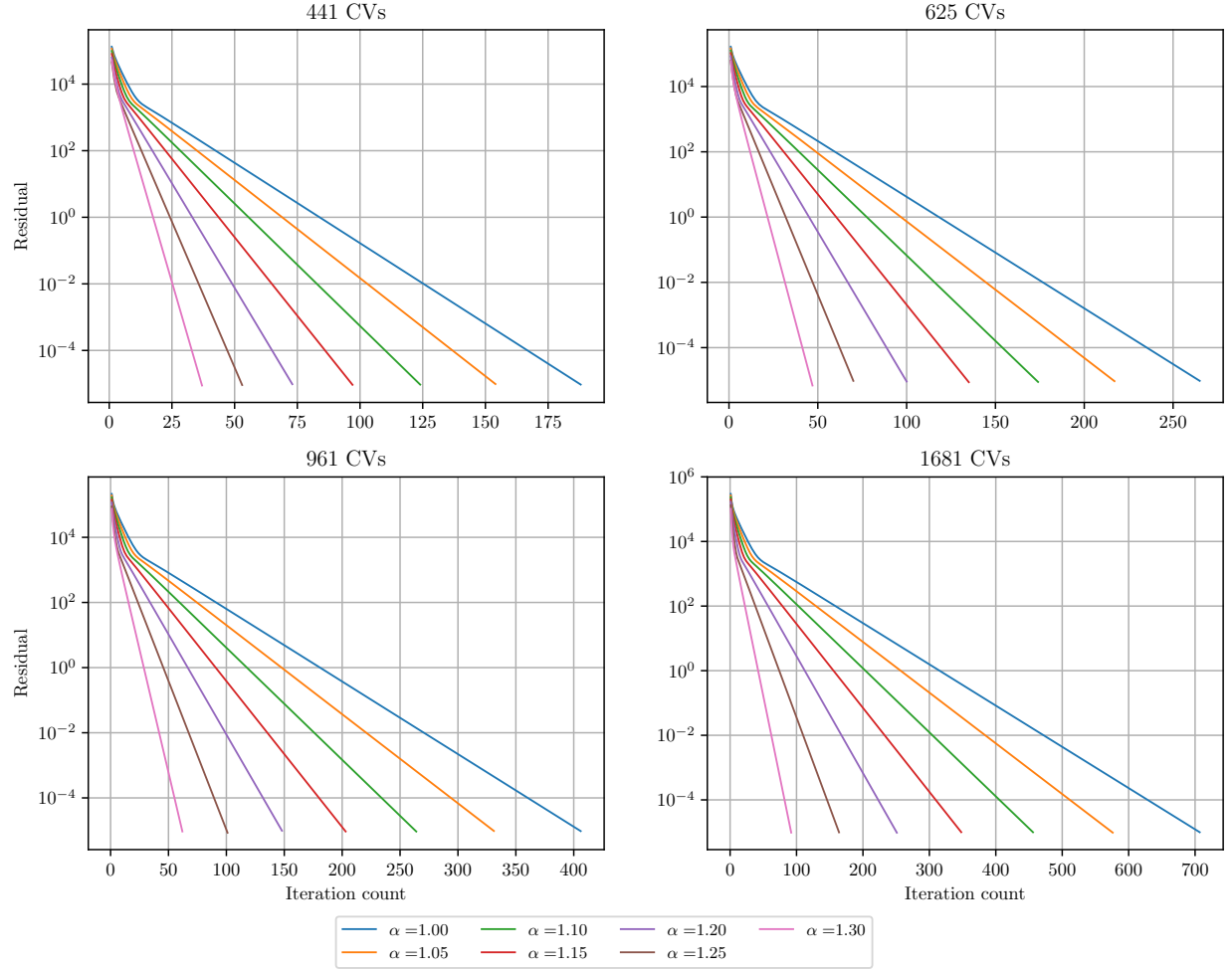


Figure 3: Plot of the residual vs. iteration count for the more refined problems.

Part c

With the final mesh refinement of 41^2 CVs and $\alpha = 1$, a colored contour plot of the temperature solution follows in Figure 4.

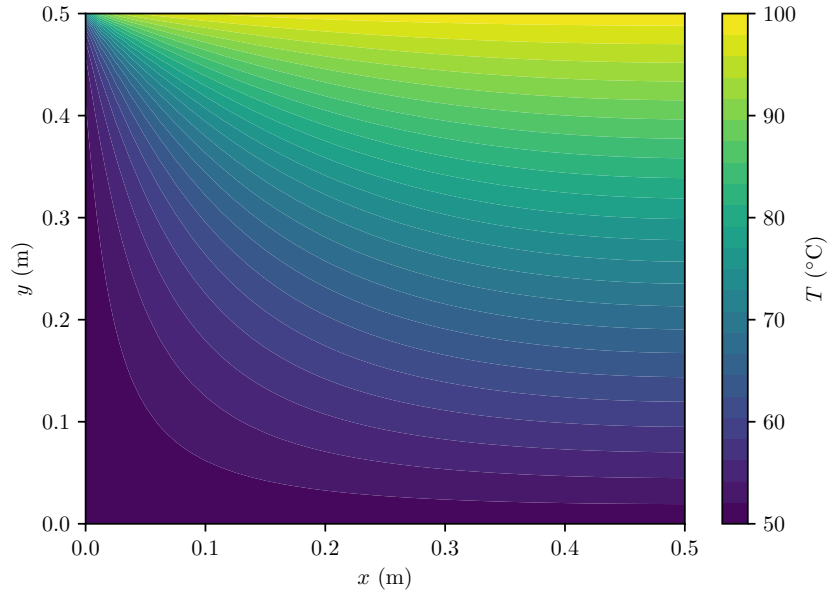


Figure 4: Plot of the solution with 41^2 CVs.

Code listing

For the implementation, we have the following files:

- **Makefile** – Allows for compiling the c++ project with **make**.
- **hwk3.cpp** – Contains the **main()** function that is required by C that runs the cases requested in this problem set.
- **Conduction2D.h / Conduction2D.cpp** – Contains the **Conduction2D** class which is the solver for the 2D heat conduction problem required in this homework.
- **Matrix.h** – Contains the **Matrix** class which provides storage for a matrix with various standard matrix operations.
- **TriDiagonal.h** – Contains the **TriDiagonal** class which provides storage for a tri-diagonal matrix including the TDMA solver found in the member function **solveTDMA()**.
- **plots.py** - Produces the plots in this report.

Makefile

```
src = $(wildcard *.cpp)
obj = $(src:.cpp=.o)
CXXFLAGS = -std=c++14
CCFLAGS = $(CXXFLAGS)

hwk-opt: $(obj)
    clang++ -o $@ $^
```

```
.PHONY: clean
clean:
    rm -f $(obj) hwk-opt
```

hwk3.cpp

```
#include "Conduction2D.h"
#include <boost/format.hpp>
#include <map>
#include <sstream>

template <typename T> void save(const std::vector<T> &v, std::string filename) {
    std::ofstream f;
    f.open(filename);
    for (unsigned int i = 0; i < v.size(); ++i)
        f << std::scientific << v[i] << std::endl;
    f.close();
}

int main() {
    // Part a: change relaxation factor
    std::cout << "Part a" << std::endl;
    for (double w : {1.0, 1.05, 1.1, 1.15, 1.2, 1.25, 1.3, 1.35, 1.4}) {
        std::cout << "    Running with w = " << w << ": ";
        Conduction2D problem(15, 15, w);
        problem.solve();

        // Save residuals if converged
        if (problem.converged()) {
            std::stringstream filename;
            filename << boost::format("results/a/w%1%.csv") % w;
            save(problem.getResiduals(), filename.str());
        }
    }
    std::cout << std::endl;

    // Part b: refine mesh and retrieve center value
    std::cout << "Part b" << std::endl;
    std::map<double, std::vector<double>> temps;
    for (unsigned int N : {15, 21, 25, 31, 41}) {
        std::cout << "    Running with " << N * N << " CVs: " << std::endl;
        for (double w : {1.0, 1.05, 1.1, 1.15, 1.2, 1.25, 1.3}) {
            std::cout << "        Running with w = " << w << ": ";

            Conduction2D problem(N, N, w);
            problem.solve();

            // Store iteration count and center solution if it converged
            if (problem.converged()) {
                std::stringstream filename;
                filename << boost::format("results/b-iterations/N%1%_w%2%.csv") % N % w;
                save(problem.getResiduals(), filename.str());
                temps[w].push_back(problem.getT((N - 1) / 2, (N - 1) / 2));
            }
        }

        // Part c: save solution for 41x41
    }
}
```

```

        if (N == 41 && w == 1.0)
            problem.saveT("results/c/solution.csv");
    }
}
// Save temperatures for each alpha
for (double w : {1.0, 1.15, 1.3})
    save(temps[w], "results/b-temps/" + std::to_string(w) + ".csv");

return 0;
}

```

Conduction2D.h

```

#ifndef CONDUCTION2D_H
#define CONDUCTION2D_H

#include <cmath>
#include <fstream>
#include <iomanip>
#include <iostream>

#include "Matrix.h"
#include "TriDiagonal.h"

/**
 * Solves a 2D heat conduction problem with dirichlet conditions on the top,
 * left, bottom and with a zero-flux condition on the right with Nx x Ny
 * internal control volumes.
 */
class Conduction2D {
public:
    Conduction2D(unsigned int Nx, unsigned int Ny, double alpha, double Lx = 0.5,
                 double Ly = 0.5, double k = 386.0, double T_L = 50.0,
                 double T_T = 100.0, double T_B = 50.0, double tol = 1e-5,
                 unsigned int max_its = 2000);

    void solve();

    // See if this is solved/converged
    bool converged() {
        return (residuals.size() != 0 && residuals.size() != max_its);
    }

    // Get the solution at the (i, j) internal node
    const double getT(unsigned int i, unsigned int j) const { return T(i, j); }

    // Get the residuals and number of iterations
    const std::vector<double> &getResiduals() const { return residuals; }
    unsigned int getNumIterations() { return residuals.size(); }

    // Save the solution
    void saveT(std::string filename) const { T.save(filename); }

private:
    double computeResidual() const;

    // Precompute operations

```



```

void precomputeProperties();
void precomputeColumn(unsigned int col);
void precomputeRow(unsigned int row);

// Solve and sweep operations
void solveColumn(unsigned int col);
void solveRow(unsigned int row);

protected:
// Number of interior nodal points in the x and y-dimensions
const unsigned int Nx, Ny;

// Geometry [m]
const double Lx, Ly, dx, dy;
// Heat conduction coefficient [W / m k]
const double k;
// Dirichlet boundary conditions (left, top, bottom) [C]
const double T_L, T_T, T_B;
// Properties stored in matrix form
Matrix a_p, a_n, a_e, a_s, a_w;

// Inverse of the relaxation coefficient
const double w_inv;
// Iteration tolerance
const double tol;
// Maximum iterations
const unsigned int max_its;

// Temperature solution
Matrix T;

// Precomputed matrices for the TDMA solves
std::vector<TriDiagonal> pre_A_x, pre_A_y;
// Precomputed RHS for the TDMA solves
std::vector<std::vector<double>> pre_b_x, pre_b_y;
// Matrices for the TDMA solves
TriDiagonal A_x, A_y;
// RHS/solution vector for the TDMA solves
std::vector<double> b_x, b_y;
// Residual for each iteration
std::vector<double> residuals;
};

#endif /* CONDUCTION2D_H */

```

Conduction2D.cpp

```

#include "Conduction2D.h"

Conduction2D::Conduction2D(unsigned int Nx, unsigned int Ny, double w,
                           double Lx, double Ly, double k, double T_L,
                           double T_T, double T_B, double tol,
                           unsigned int max_its)
: // Interior nodal points
  Nx(Nx), Ny(Ny), Lx(Lx), Ly(Ly), dx(Lx / Nx), dy(Ly / Ny),
  // Material properties
  k(k),

```

```

    // Boundary conditions
    T_L(T_L), T_T(T_T), T_B(T_B),
    // Material properties in matrix form
    a_p(Nx, Ny), a_n(Nx, Ny), a_e(Nx, Ny), a_s(Nx, Ny), a_w(Nx, Ny),
    // Solver properties
    w_inv(1.0 / w), tol(tol), max_its(max_its),
    // Initialize matrices and vectors
    T(Nx, Ny, (T_L + T_T + T_B) / 3.0), pre_A_x(Ny, Nx), pre_A_y(Nx, Ny),
    pre_b_x(Ny, std::vector<double>(Nx)),
    pre_b_y(Nx, std::vector<double>(Ny)), A_x(Nx), A_y(Ny), b_x(Nx), b_y(Ny) {
}

void Conduction2D::solve() {
    // Compute the a coefficients for each CV
    precomputeProperties();
    // Compute the unchanging LHS and RHS for each column
    for (unsigned int i = 0; i < Nx; ++i)
        precomputeColumn(i);
    // Compute the unchanging LHS and RHS for each row
    for (unsigned int j = 0; j < Ny; ++j)
        precomputeRow(j);

    // Iterate and exit when complete
    for (unsigned int l = 1; l <= max_its; ++l) {
        // Sweep south to north
        for (int j = 0; j < Ny; ++j)
            solveRow(j);
        // Sweep west to east
        for (int i = 0; i < Nx; ++i)
            solveColumn(i);
        // Sweep north to south
        for (int j = Ny - 1; j >= 0; --j)
            solveRow(j);
        // Sweep east to west
        for (int i = Nx - 1; i >= 0; --i)
            solveColumn(i);

        // Check for convergence and store residual
        double R = computeResidual();
        residuals.push_back(R);
        if (R < tol) {
            std::cout << "Converged with " << l << " iterations" << std::endl;
            return;
        }
    }

    std::cout << "Failed to converge in " << max_its << " iterations"
        << std::endl;
}

void Conduction2D::precomputeProperties() {
    // Set all neighbors to the default at first
    a_n = k * dx / dy;
    a_e = k * dy / dx;
    a_s = k * dx / dy;
    a_w = k * dy / dx;

    // Top Dirichlet
    a_n.setRow(Ny - 1, 2 * k * dx / dy);

```

```

// Right Neumann
a_e.setColumn(Nx - 1, 0);
// Bottom Dirichlet
a_s.setRow(0, 2 * k * dx / dy);
// Left dirichlet
a_w.setColumn(0, 2 * k * dy / dx);

// Center point
for (unsigned int i = 0; i < Nx; ++i)
    for (unsigned int j = 0; j < Ny; ++j)
        a_p(i, j) = a_n(i, j) + a_e(i, j) + a_s(i, j) + a_w(i, j);
}

void Conduction2D::precomputeRow(unsigned int j) {
    TriDiagonal &A = pre_A_x[j];
    std::vector<double> &b = pre_b_x[j];

    // First treat all as an internal volume
    A.addTopRow(a_p(0, j) * w_inv, -a_e(0, j));
    A.addBottomRow(-a_w(Nx - 1, j), a_p(Nx - 1, j) * w_inv);
    for (unsigned int i = 1; i < Nx - 1; ++i)
        A.addMiddleRow(i, -a_w(i, j), a_p(i, j) * w_inv, -a_e(i, j));

    // Left Dirichlet
    b[0] += T_L * a_w(0, j);
    // Top dirichlet
    if (j == Ny - 1)
        for (unsigned int i = 0; i < Nx; ++i)
            b[i] += T_T * a_n(i, j);
    // Bottom dirichlet
    else if (j == 0)
        for (unsigned int i = 0; i < Nx; ++i)
            b[i] += T_B * a_s(i, j);
}

void Conduction2D::precomputeColumn(unsigned int i) {
    TriDiagonal &A = pre_A_y[i];
    std::vector<double> &b = pre_b_y[i];

    // First treat all as an internal volume
    A.addTopRow(a_p(i, 0) * w_inv, -a_n(i, 0));
    A.addBottomRow(-a_s(i, Ny - 1), a_p(i, Ny - 1) * w_inv);
    for (unsigned int j = 1; j < Ny - 1; ++j)
        A.addMiddleRow(j, -a_s(i, j), a_p(i, j) * w_inv, -a_n(i, j));

    // Left Dirichlet
    if (i == 0)
        for (unsigned int j = 0; j < Ny; ++j)
            b[j] += T_L * a_w(i, j);
    // Top Dirichlet
    b[Ny - 1] += T_T * a_n(i, Ny - 1);
    // Bottom Dirichlet
    b[0] += T_B * a_s(i, 0);
}

void Conduction2D::solveRow(unsigned int j) {
    // Copy pre-filled Ax = b for this row
    A_x = pre_A_x[j];
    b_x = pre_b_x[j];
}

```

```

// RHS contribution from volumes above and below
if (j > 0)
    for (unsigned int i = 0; i < Nx; ++i)
        b_x[i] += T(i, j - 1) * a_s(i, j);
if (j < Ny - 1)
    for (unsigned int i = 0; i < Nx; ++i)
        b_x[i] += T(i, j + 1) * a_n(i, j);

// Relax, solve, and store solution (which is in b_x)
if (w_inv != 1)
    for (unsigned int i = 0; i < Nx; ++i)
        b_x[i] += (w_inv - 1.0) * a_p(i, j) * T(i, j);
A_x.solveTDMA(b_x);
T.setRow(j, b_x);
}

void Conduction2D::solveColumn(unsigned int i) {
    // Copy pre-filled Ax = b for this row
    A_y = pre_A_y[i];
    b_y = pre_b_y[i];

    // RHS contribution from volumes left and right
    if (i > 0)
        for (unsigned int j = 0; j < Ny; ++j)
            b_y[j] += T(i - 1, j) * a_w(i, j);
    if (i < Nx - 1)
        for (unsigned int j = 0; j < Ny; ++j)
            b_y[j] += T(i + 1, j) * a_e(i, j);

    // Relax, solve, and store solution (which is in b_y)
    if (w_inv != 1)
        for (unsigned int j = 0; j < Ny; ++j)
            b_y[j] += (w_inv - 1.0) * a_p(i, j) * T(i, j);
    A_y.solveTDMA(b_y);
    T.setColumn(i, b_y);
}

double Conduction2D::computeResidual() const {
    double R = 0.0, val = 0.0;
    // Sum over all CVs
    for (unsigned int i = 0; i < Nx; ++i)
        for (unsigned int j = 0; j < Ny; ++j) {
            // Main diagonal contribution and pre-computed RHS (from BC)
            val = a_p(i, j) * T(i, j) - pre_b_y[i][j];
            // Not on left boundary
            if (i > 0)
                val -= a_w(i, j) * T(i - 1, j);
            // Not on right boundary
            if (i < Nx - 1)
                val -= a_e(i, j) * T(i + 1, j);
            // Not on bottom boundary
            if (j > 0)
                val -= a_s(i, j) * T(i, j - 1);
            // Not top boundary
            if (j < Ny - 1)
                val -= a_n(i, j) * T(i, j + 1);
            R += std::abs(val);
        }
}

```

```

    return R;
}

```

Matrix.h

```

#ifndef MATRIX
#define MATRIX

#define NDEBUG
#include <cassert>
#include <vector>

/**
 * Class that holds a Nx x Ny matrix with common matrix operations.
 */
class Matrix {
public:
    Matrix(unsigned int Nx, unsigned int Ny, double v = 0)
        : Nx(Nx), Ny(Ny), M(Nx, std::vector<double>(Ny, v)) {}

    // Const operator for getting the (i, j) element
    const double &operator()(unsigned int i, unsigned int j) const {
        assert(i < Nx && j < Ny);
        return M[i][j];
    }
    // Operator for getting the (i, j) element
    double &operator()(unsigned int i, unsigned int j) {
        assert(i < Nx && j < Ny);
        return M[i][j];
    }
    // Operator for setting the entire matrix to a value
    void operator=(double v) {
        for (unsigned int j = 0; j < Ny; ++j)
            setRow(j, v);
    }

    // Saves the matrix in csv format
    void save(const std::string filename, unsigned int precision = 12) const {
        std::ofstream f;
        f.open(filename);
        for (unsigned int j = 0; j < Ny; ++j) {
            for (unsigned int i = 0; i < Nx; ++i) {
                if (i > 0)
                    f << ",";
                f << std::setprecision(precision) << M[i][j];
            }
            f << std::endl;
        }
        f.close();
    }

    // Set the j-th row to v
    void setRow(unsigned int j, double v) {
        assert(j < Ny);
        for (unsigned int i = 0; i < Nx; ++i)
            M[i][j] = v;
    }
}

```

```

// Set the i-th column to v
void setColumn(unsigned int i, double v) {
    assert(i < Nx);
    for (unsigned int j = 0; j < Ny; ++j)
        M[i][j] = v;
}

// Set the j-th row to vs
void setRow(unsigned int j, std::vector<double> &vs) {
    assert(j < Ny && vs.size() == Nx);
    for (unsigned int i = 0; i < Nx; ++i)
        M[i][j] = vs[i];
}

// Set the i-th column to vs
void setColumn(unsigned int i, std::vector<double> &vs) {
    assert(i < Nx && vs.size() == Ny);
    for (unsigned int j = 0; j < Ny; ++j)
        M[i][j] = vs[j];
}

private:
    // The size of this matrix
    const unsigned int Nx, Ny;

    // Matrix storage
    std::vector<std::vector<double> > M;
};

#endif /* MATRIX_H */

```

TriDiagonal.h

```

#ifndef TRIDIAGONAL_H
#define TRIDIAGONAL_H

#define NDEBUG
#include <cassert>

/**
 * Class that holds a tri-diagonal matrix and is able to perform TDMA in place
 * with a given RHS.
 */
class TriDiagonal {
public:
    TriDiagonal(unsigned int N, double v = 0)
        : N(N), A(N, v), B(N, v), C(N - 1, v) {}

    // Operator for setting the entire matrix to a value
    void operator=(TriDiagonal & from) {
        assert(from.getN() == N);
        A = from.getA();
        B = from.getB();
        C = from.getC();
    }

    // Gets the value of the (i, j) entry
    const double operator()(unsigned int i, unsigned int j) const {

```

```

    assert(i < N && j > i - 2 && j < i + 2);
    if (j == i - 1)
        return A[i];
    else if (j == i)
        return B[i];
    else if (j == i + 1)
        return C[i];
    else {
        std::cerr << "( " << i << ", " << j << ") out of TriDiagonal system";
        std::terminate();
    }
}

// Adders for the top, middle, and bottom rows
void addTopRow(double b, double c) {
    B[0] += b;
    C[0] += c;
}

void addMiddleRow(unsigned int i, double a, double b, double c) {
    assert(i < N - 1 && i != 0);
    A[i] += a;
    B[i] += b;
    C[i] += c;
}

void addBottomRow(double a, double b) {
    A[N - 1] += a;
    B[N - 1] += b;
}

// Getters for the raw vectors
const std::vector<double> &getA() const { return A; }
const std::vector<double> &getB() const { return B; }
const std::vector<double> &getC() const { return C; }

// Getter for the size
unsigned int getN() { return N; }

// Solves the system  $Ax = d$  in place where  $d$  eventually stores the solution
void solveTDMA(std::vector<double> &d) {
    // Forward sweep
    double tmp = 0;
    for (unsigned int i = 1; i < N; ++i) {
        tmp = A[i] / B[i - 1];
        B[i] -= tmp * C[i - 1];
        d[i] -= tmp * d[i - 1];
    }

    // Backward sweep
    d[N - 1] /= B[N - 1];
    for (unsigned int i = N - 2; i != std::numeric_limits<unsigned int>::max(); --i) {
        d[i] -= C[i] * d[i + 1];
        d[i] /= B[i];
    }
}

protected:
    // Matrix size (N x N)
    unsigned int N;

```

```

// Left/main/right diagonal storage
std::vector<double> A, B, C;
};

#endif /* TRIDIAGONAL_H */

```

plots.py

```

import numpy as np
import matplotlib.pyplot as plt
import glob
import os

plt.rc('text', usetex=True)
plt.rc('font', family='serif')

# Problem a
fig = plt.figure()
fig.set_figwidth(9)
fig.set_figheight(4)
alphas = [1, 1.05, 1.1, 1.15, 1.2, 1.25, 1.3, 1.35]
for alpha in alphas:
    residuals = np.loadtxt('results/a/w{}.csv'.format(alpha))
    iterations = list(range(1, len(residuals) + 1))
    plt.semilogy(iterations, residuals, '-', linewidth=1,
                  label=r'$\alpha = {:.2f}'.format(alpha))
plt.xlabel(r'Iteration count')
plt.ylabel(r'Residual')
plt.legend()
plt.tight_layout()
plt.grid()
fig.savefig('results/a.pdf')

# Problem b number of iterations
fig, ax = plt.subplots(2, 2)
fig.set_figwidth(11)
fig.set_figheight(8)
alphas = [1, 1.05, 1.1, 1.15, 1.2, 1.25, 1.3]
Ns = [[21, 25], [31, 41]]
for i in range(2):
    for j in range(2):
        N = Ns[i][j]
        for alpha in alphas:
            residuals = np.loadtxt('results/b-iterations/N{}_w{}.csv'.format(N, alpha))
            iterations = list(range(1, len(residuals) + 1))
            ax[i][j].semilogy(iterations, residuals, '-', linewidth=1,
                              label=r'$\alpha = {:.2f}'.format(alpha))
        ax[i][j].set_title('{} CVs'.format(N * N))
        ax[i][j].grid()
        if i == 1: ax[i][j].set_xlabel('Iteration count')
        if j == 0: ax[i][j].set_ylabel('Residual')
handles, labels = ax[1][1].get_legend_handles_labels()
lgd = ax[1][1].legend(handles, labels, loc='lower center', bbox_to_anchor=(-0.2, -0.34),
                      ncol=4, fontsize=9)
fig.savefig('results/b-iterations.pdf', bbox_inches='tight', bbox_extra_artists=(lgd,))

```



```

# Problem b temps
fig = plt.figure()
fig.set_figwidth(9)
fig.set_figheight(4)
Ns = [15**2, 21**2, 26**2, 31**2, 41**2]
for file in glob.glob('results/b-temps/*.csv'):
    alpha = float(os.path.splitext(file)[0].split("/")[-1])
    temps = np.loadtxt(file)
    plt.plot(Ns, temps, '-.', linewidth=1, label=r'$\alpha = $ {:.2f}'.format(alpha))
plt.xlabel(r'Number of CVs')
plt.ylabel(r'Center plate temperature ( $\theta_c$ )')
plt.ticklabel_format(useOffset=False)
plt.legend()
plt.tight_layout()
plt.grid()
fig.savefig('results/b-temps.pdf')

# Problem c
T = np.loadtxt('results/c/solution.csv', delimiter=',')
fig = plt.figure()
fig.set_figwidth(5.5)
fig.set_figheight(4)
x, y = np.meshgrid(np.linspace(0, 0.5, num=41), np.linspace(0, 0.5, num=41))
plot = plt.contourf(x, y, T, levels=np.linspace(50, 100, num=25))
cbar = fig.colorbar(plot)
cbar.set_ticks(np.arange(50, 101, 10))
cbar.ax.set_ylabel(r'$T$ ( $\theta_c$ )')
plt.xlabel('$x$ (m)')
plt.ylabel('$y$ (m)')
plt.tight_layout()
fig.savefig('results/c.pdf')

```