

Designing a Program and Subroutines

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Note: Subroutines are commonly called, depending on the programming language, modules, subprograms, methods, and functions.

Top-down design (sometimes called stepwise refinement) is used to break down an algorithm into subroutines.

Top-Down Design Process:

- The overall task of the program is broken down into a series of subtasks.
- Each of the subtasks is examined to determine whether it can be further broken down into more subtasks. This step is repeated until no more subtasks can be identified.
- Once all of the subtasks have been identified, they are written in code.

Three main tools for designing a program and its subroutines:

1. **Hierarchy Chart** – or a structure chart, a top-level visual representation of the main program and the relationships between subroutines.
2. **Flowcharts** – a diagram that graphically depicts the steps that take place in a program.
3. **Pseudocode** – or “fake code” is an informal language that has no syntax rules, it is a “mock-up” program. Each statement in the pseudocode represents an operation that can be performed in any high-level language.

Top-Down Design
Program: Solve the Heat Equation Numerically
(Solving_PDEs_numerically_Parabolic_Crank_Nicolson_Handwritten_2021-07-02_152842.pdf)

Overall Task:

Develop a general-purpose MATLAB program to solve the one-dimensional heat equation (partial differential equation - PDE) using the Crank-Nicolson method.

Steps that must be taken to perform the task:

1. Mathematical modeling – represent all the important features of the system; see the figure below for the model.
2. Derivation of governing equations. Re-write the second-order PDE as a finite difference equation.
3. Solution of the governing equations. Solve the system of equations using MATLAB, to obtain a series of spatial distributions corresponding to the state (Temp.) of the node at each time.
4. Interpretation of results. The solution of the governing equations gives the temperature distribution of the element of the system. Plot these responses.

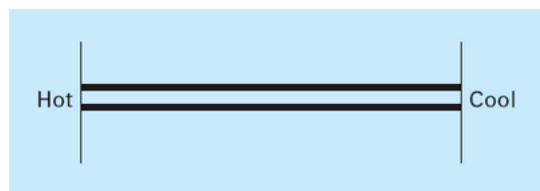


FIGURE 30.1
A thin rod, insulated at all points except at its ends.

$$[T(x, t)] = \text{heatEq_1D}(k, L, dx, dt, T(0), T(L), q0)$$

Main Program

[Output: temperature(space, time); Input: coefficient of thermal diffusivity, domain, element size, temporal step size, boundary conditions, initial conditions – heat input?]

Note: Hierarchy charts does not show the steps that are taken inside a subroutine; they do not reveal any details about how subroutines work.

1. Hierarchy Chart

Analysis Procedure Background:
Step 1: Mathematical Modeling
And
Step 2: Derivation of Governing Equations

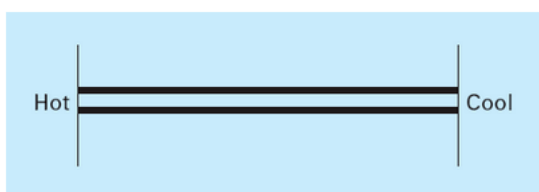
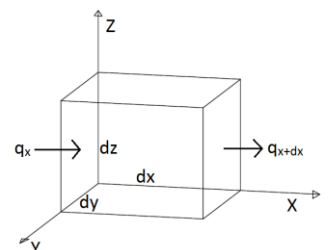


FIGURE 30.1
A thin rod, insulated at all points except at its ends.



Conservation of heat principle is used to develop a heat balance for the differential element:
(rate of heat in) = (rate of heat out) + (rate of heat stored)

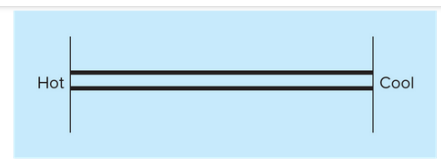


FIGURE 30.1 A thin rod, insulated at all points except at its ends.

$$q(x)\Delta y\Delta z\Delta t - q(x+\Delta x)\Delta y\Delta z\Delta t = \Delta x\Delta y\Delta z\rho C\Delta T$$

Dividing by the volume of the element ($= \Delta x\Delta y\Delta z$) and Δt gives

$$\frac{q(x) - q(x+\Delta x)}{\Delta x} = \rho C \frac{\Delta T}{\Delta t}$$

Taking the limit yields

$$\frac{\partial q}{\partial x} = \rho C \frac{\partial T}{\partial t}$$

Substituting Fourier's law of heat conduction [Eq. (29.4)] results in

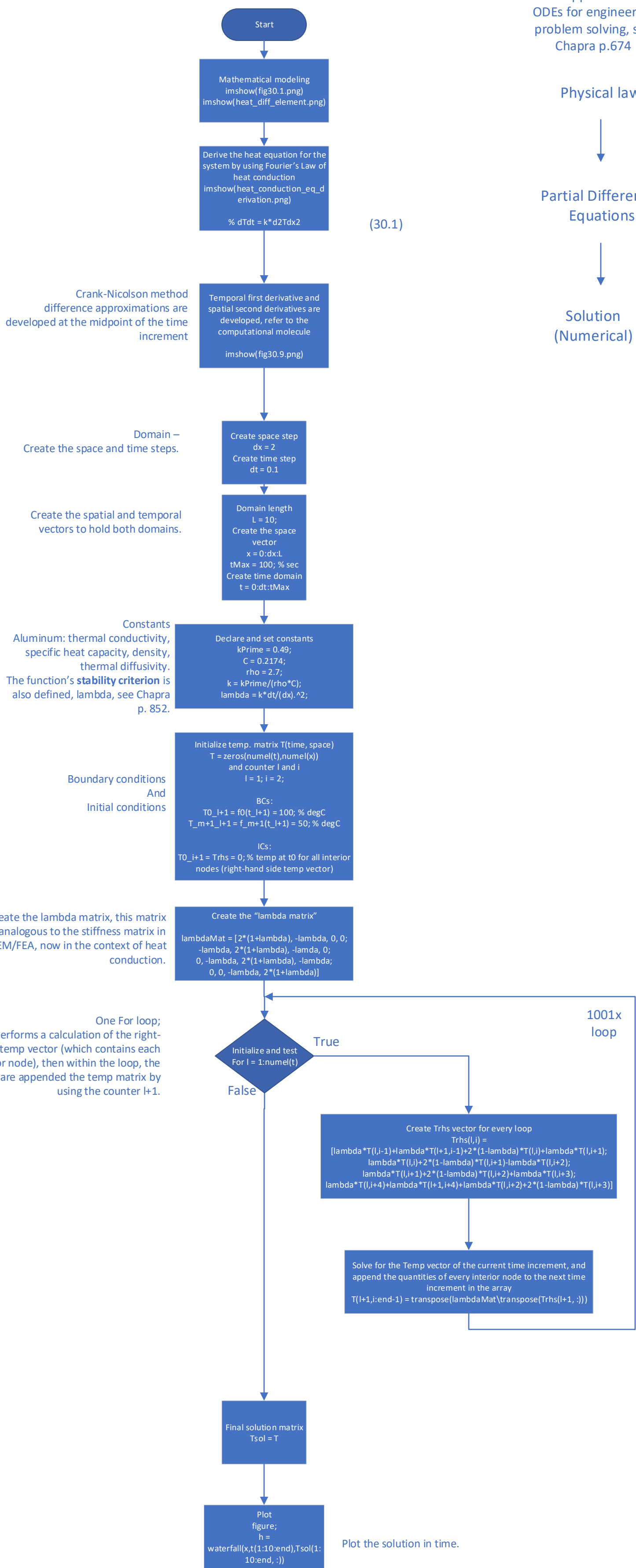
$$k \frac{\partial^2 T}{\partial x^2} = \frac{\partial T}{\partial t}$$

which is the *heat-conduction equation*.

See Chapra, p. 840.

Step 1:
Mathematical
ModelingStep 2:
Derivation of Governing
EquationsStep 3:
Solution of the Governing
EquationsStep 4:
Interpretation of Results

2. Flowchart



Sequence of events in the application of ODEs for engineering problem solving, see Chapra p.674

Physical law

Partial Differential Equations

Solution (Numerical)

3. Pseudocode

% Start
% Step 1 Mathematical Modeling

% Display image that represents system model
imshow(fig30.1.png)
imshow(heat_diff_element.png)

% Step 2: Derivation of Governing Equations

% Fourier's Law of heat conduction
imshow(heat_conduction_eq_derivation.png)
% dT/dt = k*dt^2/dx^2

% Step 3: Solution of the Governing Equations

% Crank-Nicolson method
% diff approximations are developed at the midpoint of the time increment
imshow(fig30.9.png) % computational molecule

% Domain
dx = 2; % space step
dt = 0.1; % time step
L = 10; % domain length
x = 0:dx:L; % space vector
tMax = 100; % temporal domain
t = 0:dt:tMax; % time vector

% Declare and set constants, aluminum material
kPrime = 0.49; % thermal conductivity
C = 0.2174; % specific heat capacity
rho = 2.7; % density
k = kPrime/(rho*C);

% Function's stability criterion
lambda = k*dt/(dx.^2);

% Boundary conditions and Initial conditions

% Initialize Temp matrix, which will be a function of node and time step, T(time, space), and counters l and i
T = zeros(numel(t), numel(x)) % zeros for all time x all nodes

l = 1; i = 2; % time step number one is initialized at element 1, node number is initialized at node #2, because the spatial second derivative is determined at the midpoint (current node) by averaging the difference approximations at the beginning and end of the time increment; this is done through averaging temp quantities for three nodes, and therefore the first approximation will be of nodes 1, 2, and 3 – which is why we initialize i at 2. Essentially, we are integrating forward in time.

% BCs:

T(:, x(1)) = 100; % degC, T as a function of all time and node 1 (only), that is, the left end is fixed at the value given
T(:, x(end)) = 50; % degC, T will be set to the value for all time to the last (end) node in the system, that is, the right end is fixed

% ICs:

% Temp data for the interior nodes, this is the RHS temp vector
intNodes = numel(x) - 2; % calculates the number of interior nodes
Trhs = zeros(numel(t), intNodes); % Declare, set, and initialize Trhs, to hold the RHS temp vector (INTERIOR) nodes only

% Create the lambda matrix

% The lambda matrix is analogous to the stiffness matrix in FEA, but in the context of heat transfer (obv)

lambdaMat = [2*(1+lambda), -lambda, 0, 0;
-lambda, 2*(1+lambda), -lambda, 0;
0, -lambda, 2*(1+lambda), -lambda;
0, 0, -lambda, 2*(1+lambda)]

% Calculate the solution Temp for all time, and all nodes, then append the temp-node value to the Temp matrix using counter l=1.
% Noting that t=0 sec starts at l=1, and from ICs the interior nodes temp = 0; and t=0.1 sec is l=2.
for l = 1:numel(t)

% Node 2 temp of Temp matrix, time +1:
Trhs(l+1, i-1) = lambda*T(l,i-1)+lambda*T(l+1,i-1)+2*(1-lambda)*T(l,i)+lambda*T(l,i+1);
% Node 3:
Trhs(l+1, i) = lambda*T(l,i)+2*(1-lambda)*T(l,i+1)-lambda*T(l,i+2);
% Node 4:
Trhs(l+1, i+1) = lambda*T(l,i+1)+2*(1-lambda)*T(l,i+2)+lambda*T(l,i+3);
% Node 5:
Trhs(l+1, i+2) = lambda*T(l,i+4)+lambda*T(l+1,i+4)+lambda*T(l,i+2)+2*(1-lambda)*T(l,i+3);

% Solve the system of equations, outputting Temp for each time-node combination, append the quantities of every interior node to the next time increment in the array.
T(l+1,:end-1) = transpose(lambdaMat*transpose(Trhs(l+1,:)));

end

% Final solution matrix
Tsol = T

% Step 4: Interpretation of the Results

% Plot solution in time
figure;
h = waterfall(x, t(1:10:end), Tsol(1:10:end))

% End program ☺