# Chapter 2

# **Finite Element Approximation**

#### 2.1 Piecewise linear approximation

**One-dimensional piecewise linear approximation** Let us approximate one-dimensional function f(x) over region  $P_iP_j$  in one-dimensional space. The function takes values  $f_i$ ,  $f_j$  at points  $P_i$ ,  $P_j$ . Let  $x_i$ ,  $x_j$  be coordinates of points  $P_i$ ,  $P_j$ . Let P be any point of which coordinate is given by x. Let us introduce the following two functions

$$N_{i,j}(x) = \frac{PP_j}{P_i P_j} = \frac{x_j - x}{x_j - x_i}$$
(2.1.1a)

$$N_{j,i}(x) = \frac{\mathbf{P}_i \mathbf{P}}{\mathbf{P}_i \mathbf{P}_j} = \frac{x - x_i}{x_j - x_i}$$
(2.1.1b)

Noting that

$$N_{i,j}(x) = \begin{cases} 1 & x = x_i \\ 0 & x = x_j \end{cases}, \qquad N_{j,i}(x) = \begin{cases} 0 & x = x_i \\ 1 & x = x_j \end{cases}$$

linear approximation of function f(x) over region  $P_iP_j$  is described as follows:

$$L_{i,j}(x) = f_i N_{i,j}(x) + f_j N_{j,i}(x).$$
(2.1.2)

This function is linear since both  $N_{i,j}(x)$  and  $N_{j,i}(x)$  are linear. Also, this function satisfies

$$L_{i,j}(x_i) = f_i N_{i,j}(x_i) + f_j N_{j,i}(x_i) = f_i$$
  

$$L_{i,j}(x_j) = f_i N_{i,j}(x_j) + f_j N_{j,i}(x_j) = f_j$$

concluding that function  $L_{i,j}(x)$  provides one-dimensional approximation over  $P_i P_j$ .

**Two-dimensional piecewise linear approximation** Let us approximate two-dimensional function f(x, y) over triangle region  $\triangle P_i P_j P_k$  in two-dimensional space. The function takes values  $f_i$ ,  $f_j$ ,  $f_k$  at points  $P_i$ ,  $P_j$ ,  $P_k$ . Let  $(x_i, y_i)$  be coordinates of point  $P_i$ ,  $(x_j, y_j)$  be coordinates of point  $P_j$ , and  $(x_k, y_k)$  be coordinates of point  $P_k$ . Let P(x, y) be any point of

which coordinate is given by (x, y). Let us introduce the following three functions

$$N_{i,j,k}(x,y) = \frac{\triangle PP_j P_k}{\triangle P_i P_j P_k} = \frac{(y_j - y_k)x - (x_j - x_k)y + (x_j y_k - x_k y_j)}{2\triangle P_i P_j P_k}$$
(2.1.3a)

$$N_{j,k,i}(x,y) = \frac{\triangle \mathbf{P}_i \mathbf{P} \mathbf{P}_k}{\triangle \mathbf{P}_i \mathbf{P}_i \mathbf{P}_k} = \frac{(y_k - y_i)x - (x_k - x_i)y + (x_k y_i - x_i y_k)}{2\triangle \mathbf{P}_i \mathbf{P}_k \mathbf{P}_k}$$
(2.1.3b)

$$N_{k,i,j}(x,y) = \frac{\triangle P_i P_j P}{\triangle P_i P_j P_k} = \frac{(y_i - y_j)x - (x_i - x_j)y + (x_i y_j - x_j y_i)}{2\triangle P_i P_j P_k}$$
(2.1.3c)

where

$$2\triangle \mathbf{P}_i\mathbf{P}_j\mathbf{P}_k = (x_iy_j - x_jy_i) + (x_jy_k - x_ky_j) + (x_ky_i - x_iy_k).$$

Noting that

$$N_{i,j,k}(x,y) = \begin{cases} 1 & \text{at } \mathbf{P}_i \\ 0 & \text{at } \mathbf{P}_j, \mathbf{P}_k \end{cases}$$
$$N_{j,k,i}(x,y) = \begin{cases} 1 & \text{at } \mathbf{P}_j \\ 0 & \text{at } \mathbf{P}_k, \mathbf{P}_i \end{cases}$$
$$N_{k,i,j}(x,y) = \begin{cases} 1 & \text{at } \mathbf{P}_k \\ 0 & \text{at } \mathbf{P}_i, \mathbf{P}_j \end{cases}$$

linear approximation of f(x, y) over region  $\triangle P_i P_j P_k$  is described as follows:

$$L_{i,j,k}(x,y) = f_i N_{i,j,k}(x,y) + f_j N_{j,k,i}(x,y) + f_k N_{k,i,j}(x,y)$$
(2.1.4)

This function is linear since  $N_{i,j,k}(x,y)$ ,  $N_{j,k,i}(x,y)$ , and  $N_{k,i,j}(x,y)$  are linear. Also, this function satisfies

$$L_{i,j,k}(x_i, y_i) = f_i, \qquad L_{i,j,k}(x_j, y_j) = f_j, \qquad L_{i,j,k}(x_k, y_k) = f_k$$

concluding that function  $L_{i,j,k}(x,y)$  provides two-dimensional approximation over triangle region  $\triangle \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k$ .

**Three-dimensional piecewise linear approximation** Let us approximate three-dimensional function f(x, y, z) over tetrahedron region  $\Diamond P_i P_j P_k P_l$  in three-dimensional space. The function takes values  $f_i$ ,  $f_j$ ,  $f_k$ ,  $f_l$  at points  $P_i$ ,  $P_j$ ,  $P_k$ ,  $P_l$ . Let  $(x_i, y_i, z_i)$  be coordinates of point  $P_i$ ,  $(x_j, y_j, z_j)$  be coordinates of point  $P_j$ ,  $(x_k, y_k, z_k)$  be coordinates of point  $P_k$ , and  $(x_l, y_l, z_l)$  be coordinates of point  $P_l$ . Let P(x, y, z) be any point of which coordinate is given by (x, y, z). Let us introduce the following four functions

$$N_{i,j,k,l}(x,y,z) = \frac{\Diamond \mathrm{PP}_j \mathrm{P}_k \mathrm{P}_l}{\Diamond \mathrm{P}_i \mathrm{P}_j \mathrm{P}_k \mathrm{P}_l}$$
(2.1.5a)

$$N_{j,k,l,i}(x,y,z) = \frac{\Diamond \mathbf{P}_i \mathbf{P}_k \mathbf{P}_l}{\Diamond \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k \mathbf{P}_l}$$
(2.1.5b)

$$N_{k,l,i,j}(x,y,z) = \frac{\Diamond \mathbf{P}_i \mathbf{P}_l \mathbf{P}_l}{\Diamond \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k \mathbf{P}_l}$$
(2.1.5c)

$$N_{l,i,j,k}(x,y,z) = \frac{\Diamond \mathbf{P}_i \mathbf{P}_i \mathbf{P}_k \mathbf{P}}{\Diamond \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k \mathbf{P}_l}$$
(2.1.5d)

Then, linear approximation of f(x, y, z) over region  $\langle P_i P_j P_k P_l$  is described as follows:

$$L_{i,j,k,l}(x,y,z) = f_i N_{i,j,k,l}(x,y,z) + f_j N_{j,k,l,i}(x,y,z) + f_k N_{k,l,i,j}(x,y,z) + f_l N_{l,i,j,k}(x,y,z)$$

This function  $L_{i,j,k,l}(x, y, z)$  provides three-dimensional approximation over tetrahedron region  $\langle \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k \mathbf{P}_l$ .

#### 2.2 One-dimensional finite element approximation

Strain potential and kinetic energies are formulated as integral forms over one-, two-, or three-dimensional regions. It is difficult or impossible to analytically calculate such integrals. *Finite element approximation* provides methods to calculate the integrals numerically. Finite element approximation employs divide-and-conquer approach, which is outlined as follows:

- **Step 1** Obtain integral form with respect to unknown functions.
- Step 2 Divide the integral into a finite number of integrals over small regions.
- **Step 3** Approximate unknown functions to calculate integrals over small regions.
- **Step 4** Sum up the calculated integrals over small regions.

Recall that strain potential energy of a one-dimensional soft robot is given by eq. (1.5.3), that is:

$$U = \int_0^L \frac{1}{2} E \varepsilon^2 A \, \mathrm{d}x = \int_0^L \frac{1}{2} E A \left(\frac{\partial u}{\partial x}\right)^2 \mathrm{d}x \tag{2.2.1}$$

This integral U includes one unknown function u(x), which should be obtained. The above integral over region [0, L] can be divided into, for example, integrals over four small regions:

$$\int_0^L = \int_{x_1}^{x_2} + \int_{x_2}^{x_3} + \int_{x_3}^{x_4} + \int_{x_4}^{x_5}$$

Applying piecewise linear approximation, we analytically or numerically calculate individual integrals over small regions, resulting that we can obtain integral U.

Finite element approximation of strain potential energy Let us detail the above procedure. Divide region [0, L] into a finite number of small regions. Here we divide the region into four equal regions. Width of the small regions is h = L/4. End points of small regions are referred to as *nodal points*. Here we have five nodal points. Let us describe the nodal points as  $x_1 = 0$ ,  $x_2 = h$ ,  $x_3 = 2h$ ,  $\cdots$ ,  $x_5 = L$ . Dividing integral interval [0, L] into small regions, we have

$$U = \int_{x_1}^{x_2} \frac{1}{2} EA\left(\frac{\mathrm{d}u}{\mathrm{d}x}\right)^2 \mathrm{d}x + \int_{x_2}^{x_3} \frac{1}{2} EA\left(\frac{\mathrm{d}u}{\mathrm{d}x}\right)^2 \mathrm{d}x + \dots + \int_{x_4}^{x_5} \frac{1}{2} EA\left(\frac{\mathrm{d}u}{\mathrm{d}x}\right)^2 \mathrm{d}x.$$
(2.2.2)

We apply piecewise linear approximation (eq. (2.1.2)) to function u(x) over small region  $[x_i, x_j]$ . Piecewise linear approximation of the function is described as follows:

$$u(x) = u_i N_{i,j}(x) + u_j N_{j,i}(x), \quad x \in [x_i, x_j]$$
(2.2.3)

where  $u_i$ ,  $u_j$  represent displacements at nodal points  $P(x_i)$ ,  $P(x_j)$ . Through this approximation, function u(x) can be described by five parameters  $u_1, u_2, \dots, u_5$ .

Let us substitute the above piecewise linear approximation into individual integrals over small regions. For sake of simplicity, assume that Young's modulus E and cross-sectional area A are constants. Substituting piecewise linear approximation given in eq. (2.2.3) into integral over small region  $[x_i, x_j]$ , we have

$$\int_{x_i}^{x_j} \frac{1}{2} EA\left(\frac{\mathrm{d}u}{\mathrm{d}x}\right)^2 \mathrm{d}x = \frac{1}{2} \begin{bmatrix} u_i & u_j \end{bmatrix} \frac{EA}{h} \begin{bmatrix} 1 & -1 \\ -1 & 1 \end{bmatrix} \begin{bmatrix} u_i \\ u_j \end{bmatrix}$$

(see Problem 5 in Chapter 2). Consequently, we have

$$U = \frac{1}{2} \begin{bmatrix} u_1 & u_2 \end{bmatrix} \frac{EA}{h} \begin{bmatrix} 1 & -1 \\ -1 & 1 \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \end{bmatrix}$$
$$+ \frac{1}{2} \begin{bmatrix} u_2 & u_3 \end{bmatrix} \frac{EA}{h} \begin{bmatrix} 1 & -1 \\ -1 & 1 \end{bmatrix} \begin{bmatrix} u_2 \\ u_3 \end{bmatrix} + \cdots$$
$$+ \frac{1}{2} \begin{bmatrix} u_4 & u_5 \end{bmatrix} \frac{EA}{h} \begin{bmatrix} 1 & -1 \\ -1 & 1 \end{bmatrix} \begin{bmatrix} u_4 \\ u_5 \end{bmatrix}$$

which directly yields

$$U = \frac{1}{2} \begin{bmatrix} u_1 & u_2 & u_3 & u_4 & u_5 \end{bmatrix} \frac{EA}{h} \begin{bmatrix} 1 & -1 & & & \\ -1 & 2 & -1 & & \\ & -1 & 2 & -1 & \\ & & -1 & 2 & -1 \\ & & & -1 & 1 \end{bmatrix} \begin{bmatrix} u_1 \\ u_2 \\ u_3 \\ u_4 \\ u_5 \end{bmatrix}.$$

Introducing nodal displacement vector

$$\boldsymbol{u}_{\mathrm{N}} = \begin{bmatrix} u_{1} \\ u_{2} \\ u_{3} \\ u_{4} \\ u_{5} \end{bmatrix}$$
(2.2.4)

and  ${\it stiffness}\ {\it matrix}$ 

$$K = \frac{EA}{h} \begin{bmatrix} 1 & -1 & & \\ -1 & 2 & -1 & \\ & -1 & 2 & -1 \\ & & -1 & 2 & -1 \\ & & & -1 & 1 \end{bmatrix},$$
 (2.2.5)

strain potential energy is described by the following quadratic form:

$$U = \frac{1}{2} \boldsymbol{u}_{\mathrm{N}}^{\mathsf{T}} K \boldsymbol{u}_{\mathrm{N}}.$$
(2.2.6)

Note that K is a band matrix.

Let us calculate strain potential energy of a one-dimensional soft robot with non-uniform cross-sectional area. Let function A(x) denote the cross-sectional area at P(x). Assume that Young's modulus E is constant. Recalling that du/dx takes a constant value  $(-u_i + u_j)/h$  in small region  $[x_i, x_j]$ , strain potential energy over the region is given as

$$\int_{x_i}^{x_j} \frac{1}{2} EA(x) \left(\frac{\mathrm{d}u}{\mathrm{d}x}\right)^2 \mathrm{d}x = \frac{1}{2} E\left(\frac{-u_i + u_j}{h}\right)^2 \int_{x_i}^{x_j} A(x) \mathrm{d}x$$
$$= \frac{1}{2} \begin{bmatrix} u_i & u_j \end{bmatrix} \frac{E}{h^2} \begin{bmatrix} V_{i,j} & -V_{i,j} \\ -V_{i,j} & V_{i,j} \end{bmatrix} \begin{bmatrix} u_i \\ u_j \end{bmatrix}$$

where

$$V_{i,j} = \int_{x_i}^{x_j} A(x) \,\mathrm{d}x \tag{2.2.7}$$

represents the volume of the three-dimensional region specified by small region  $[x_i, x_j]$ . Thus, when region [0, L] is divided into four small regions, stiffness matrix is described as follows:

$$K = \frac{E}{h^2} \begin{bmatrix} V_{1,2} & -V_{1,2} \\ -V_{1,2} & V_{1,2} + V_{2,3} & -V_{2,3} \\ & -V_{2,3} & V_{2,3} + V_{3,4} & -V_{3,4} \\ & & -V_{3,4} & V_{3,4} + V_{4,5} & -V_{4,5} \\ & & & -V_{4,5} & V_{4,5} \end{bmatrix}.$$
 (2.2.8)

This matrix K is also a band matrix.

Let us reformulate the above calculation. Assume again that Young's modulus E and cross-sectional area A are constants. Potential energy stored in region  $[x_i, x_j]$  is given by

$$U_{i,j} = \frac{1}{2} \begin{bmatrix} x_i & x_j \end{bmatrix} K_{i,j} \begin{bmatrix} x_i \\ x_j \end{bmatrix}$$
(2.2.9)

where

$$K_{i,j} = \frac{EA}{h} \begin{bmatrix} 1 & -1\\ -1 & 1 \end{bmatrix}$$
(2.2.10)

We obtain stiffness matrix K (eq. (2.2.5)) by synthesizing matrices  $K_{1,2}$ ,  $K_{2,3}$ ,  $K_{3,4}$ , and  $K_{4,5}$ . Let us introduce operator  $\oplus$  to describe this synthesizing:

$$K = K_{1,2} \oplus K_{2,3} \oplus K_{3,4} \oplus K_{4,5} \tag{2.2.11}$$

This equation implies that summing up all contributions of  $K_{1,2}$  through  $K_{4,5}$  yields stiffness matrix K. Note that

(1, 1)-th element of  $K_{i,j}$  contributes to (i, i)-th element of K, (1, 2)-th element of  $K_{i,j}$  contributes to (i, j)-th element of K, (2, 1)-th element of  $K_{i,j}$  contributes to (j, i)-th element of K, (2, 2)-th element of  $K_{i,j}$  contributes to (j, j)-th element of K.

We simply describe these contributions as

 $(1,2) \times (1,2)$  elements of  $K_{i,j}$  contribute to  $(i,j) \times (i,j)$  elements of K.

Namely, operator × denotes direct product:  $(1, 2) \times (1, 2)$  implies (1, 1), (1, 2), (2, 1), (2, 2) while  $(i, j) \times (i, j)$  implies (i, i), (i, j), (j, i), (j, j).

**Finite element approximation of kinetic energy** Let us calculate kinetic energy of a one-dimensional soft robot given by eq. (1.6.2), that is:

$$T = \int_0^L \frac{1}{2} \rho A \left(\frac{\partial u}{\partial t}\right)^2 \mathrm{d}x = \int_0^L \frac{1}{2} \rho A \dot{u}^2 \,\mathrm{d}x \tag{2.2.12}$$

For sake of simplicity, assume that density  $\rho$  and cross-sectional area A are constants. Dividing integral region [0, L] into four equal regions, we have

$$T = \int_{x_1}^{x_2} \frac{1}{2} \rho A \dot{u}^2 dx + \int_{x_2}^{x_3} \frac{1}{2} \rho A \dot{u}^2 dx + \dots + \int_{x_4}^{x_5} \frac{1}{2} \rho A \dot{u}^2 dx$$
(2.2.13)

Piecewise linear approximation of function u(x,t) over small region  $[x_i, x_j]$  is described as follows:

$$u(x,t) = u_i(t) N_{i,j}(x) + u_j(t) N_{j,i}(x), \quad x \in [x_i, x_j]$$
(2.2.14)

Note that  $u_i$ ,  $u_j$  depend on time t whereas functions  $N_{i,j}(x)$ ,  $N_{j,i}(x)$  are not. Differentiating the above equation with respect time t yields

$$\dot{u}(x,t) = \dot{u}_i(t) N_{i,j}(x) + \dot{u}_j(t) N_{j,i}(x), \quad x \in [x_i, x_j]$$
(2.2.15)

Applying the above equation into integral over small region  $[x_i, x_j]$ , we have

$$\int_{x_i}^{x_j} \frac{1}{2} \rho A \dot{u}^2 \mathrm{d}x = \frac{1}{2} \begin{bmatrix} \dot{u}_i & \dot{u}_j \end{bmatrix} \rho A h \begin{bmatrix} 1/3 & 1/6 \\ 1/6 & 1/3 \end{bmatrix} \begin{bmatrix} \dot{u}_i \\ \dot{u}_j \end{bmatrix}$$

(see Problem 4 in Chapter 2). Consequently,

$$T = \frac{1}{2} \begin{bmatrix} \dot{u}_1 & \dot{u}_2 \end{bmatrix} \rho Ah \begin{bmatrix} 1/3 & 1/6 \\ 1/6 & 1/3 \end{bmatrix} \begin{bmatrix} \dot{u}_1 \\ \dot{u}_2 \end{bmatrix} \\ + \frac{1}{2} \begin{bmatrix} \dot{u}_2 & \dot{u}_3 \end{bmatrix} \rho Ah \begin{bmatrix} 1/3 & 1/6 \\ 1/6 & 1/3 \end{bmatrix} \begin{bmatrix} \dot{u}_2 \\ \dot{u}_3 \end{bmatrix} + \cdots \\ + \frac{1}{2} \begin{bmatrix} \dot{u}_4 & \dot{u}_5 \end{bmatrix} \rho Ah \begin{bmatrix} 1/3 & 1/6 \\ 1/6 & 1/3 \end{bmatrix} \begin{bmatrix} \dot{u}_4 \\ \dot{u}_5 \end{bmatrix}$$

which directly yields

$$T = \frac{1}{2} \dot{\boldsymbol{u}}_{\mathrm{N}}^{\mathsf{T}} M \dot{\boldsymbol{u}}_{\mathrm{N}}$$
(2.2.16)

where  $\dot{\boldsymbol{u}}_{\mathrm{N}} = [\dot{u}_1, \dot{u}_2, \cdots, \dot{u}_5]^{\top}$  and

$$M = \rho Ah \cdot \frac{1}{6} \begin{bmatrix} 2 & 1 & & \\ 1 & 4 & 1 & \\ & 1 & 4 & 1 \\ & & 1 & 4 & 1 \\ & & & 1 & 2 \end{bmatrix}$$
(2.2.17)

Matrix M is referred to as a *inertia matrix*. Note that M is a band matrix. Sum of all elements of M coincides with the total mass, implying that the inertia matrix defines its distribution. The above calculation is simply described as

$$M = M_{1,2} \oplus M_{2,3} \oplus M_{3,4} \oplus M_{4,5} \tag{2.2.18}$$

where

$$M_{i,j} = \frac{\rho A h}{6} \begin{bmatrix} 2 & 1\\ 1 & 2 \end{bmatrix}$$
(2.2.19)

denotes a partial inertia matrix corresponding to region  $[x_i, x_j]$ .

Let us calculate kinetic energy of a one-dimensional soft robot with non-uniform crosssectional area. Let function A(x) denote the cross-sectional area at P(x). Assume that density  $\rho$  is constant. Kinetic energy over small region  $[x_i, x_j]$  is then given by

$$\int_{x_i}^{x_j} \frac{1}{2} \rho A \dot{u}^2 \mathrm{d}x = \frac{1}{2} \begin{bmatrix} \dot{u}_i & \dot{u}_j \end{bmatrix} \rho \begin{bmatrix} \bar{V}_{i,j}^{i,i} & \bar{V}_{i,j}^{i,j} \\ \bar{V}_{i,j}^{i,j} & \bar{V}_{i,j}^{j,j} \end{bmatrix} \begin{bmatrix} \dot{u}_i \\ \dot{u}_j \end{bmatrix}$$



Figure 2.1: Approximation of two-dimensional region

where

$$\bar{V}_{i,j}^{i,i} = \int_{x_i}^{x_j} A(x) \{N_{i,j}(x)\}^2 \mathrm{d}x, \qquad \bar{V}_{i,j}^{j,j} = \int_{x_i}^{x_j} A(x) \{N_{j,i}(x)\}^2 \mathrm{d}x,$$
$$\bar{V}_{i,j}^{i,j} = \int_{x_i}^{x_j} A(x) N_{i,j}(x) N_{j,i}(x) \mathrm{d}x$$

Thus, when region [0, L] is divided into four small regions, inertia matrix is described as follows:

$$M = \rho \begin{bmatrix} \bar{V}_{1,2}^{1,1} & \bar{V}_{1,2}^{1,2} \\ \bar{V}_{1,2}^{1,2} & \bar{V}_{1,2}^{2,2} + \bar{V}_{2,3}^{2,2} & \bar{V}_{2,3}^{2,3} \\ & \bar{V}_{2,3}^{2,3} & \bar{V}_{2,3}^{3,3} + \bar{V}_{3,4}^{3,3} & \bar{V}_{3,4}^{3,4} \\ & & \bar{V}_{3,4}^{3,4} & \bar{V}_{3,4}^{4,4} + \bar{V}_{4,5}^{4,4} & \bar{V}_{4,5}^{4,5} \\ & & & \bar{V}_{4,5}^{4,5} & \bar{V}_{4,5}^{5,5} \end{bmatrix}.$$

This matrix M is also a band matrix. Note that  $\overline{V}_{i,j}^{i,i} + \overline{V}_{i,j}^{j,j} + 2\overline{V}_{i,j}^{i,j} = V_{i,j}$ , implying that the inertia matrix defines the distribution of the total mass.

#### 2.3 Two-dimensional finite element approximation

Strain potential energy and kinetic energy of a two-dimensional soft robot are formulated by integrals over two-dimensional region S, which is often described by an irregular shape, making analytical calculation of integrals difficult or impossible. Let us approximate twodimensional region S (Fig. 2.1(a)) by a set of small triangles (Fig. 2.1(b)). Then, integral over two-dimensional region S can be approximated by the sum of integrals over small triangles:

$$\int_{S} \approx \sum_{\Delta \mathbf{P}_{i} \mathbf{P}_{j} \mathbf{P}_{k}} \int_{\Delta \mathbf{P}_{i} \mathbf{P}_{j} \mathbf{P}_{k}}$$

Here we apply piecewise linear approximation to individual integrals over small triangles so that we can analytically or numerically calculate the integrals.

**Finite element approximation of kinetic energy** Let us calculate kinetic energy of a two-dimensional soft robot given by eq. (1.6.4), that is:

$$T = \int_{S} \frac{1}{2} \rho \, \dot{\boldsymbol{u}}^{\top} \dot{\boldsymbol{u}} \, h \, \mathrm{d}S \tag{2.3.1}$$



Figure 2.2: Example of rectangle region

First, we calculate integral over triangle region  $\triangle P_i P_j P_k$ :

$$T_{i,j,k} = \int_{\Delta \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k} \frac{1}{2} \rho \, \dot{\boldsymbol{u}}^\top \dot{\boldsymbol{u}} \, h \, \mathrm{d}S \tag{2.3.2}$$

Piecewise linear approximation of function u over triangle region  $\triangle P_i P_j P_k$  is described as follows:

$$u = u_i N_{i,j,k} + u_j N_{j,k,i} + u_k N_{k,i,j}.$$
(2.3.3)

Noting that  $u_i, u_j, u_k$  depend on time while  $N_{i,j,k}, N_{j,k,i}, N_{k,i,j}$  do not, we have

$$\dot{\boldsymbol{u}} = \dot{\boldsymbol{u}}_i \, N_{i,j,k} + \dot{\boldsymbol{u}}_j \, N_{j,k,i} + \dot{\boldsymbol{u}}_k \, N_{k,i,j} \tag{2.3.4}$$

which directly yields

$$\dot{\boldsymbol{u}}^{\top} \dot{\boldsymbol{u}} = \\ \begin{bmatrix} \dot{\boldsymbol{u}}_{i}^{\top} & \dot{\boldsymbol{u}}_{j}^{\top} & \dot{\boldsymbol{u}}_{k}^{\top} \end{bmatrix} \begin{bmatrix} \{N_{i,j,k}\}^{2} I_{2\times 2} & N_{i,j,k} N_{j,k,i} I_{2\times 2} & N_{i,j,k} N_{k,i,j} I_{2\times 2} \\ N_{i,j,k} N_{j,k,i} I_{2\times 2} & \{N_{j,k,i}\}^{2} I_{2\times 2} & N_{j,k,i} N_{k,i,j} I_{2\times 2} \\ N_{i,j,k} N_{k,i,j} I_{2\times 2} & N_{j,k,i} N_{k,i,j} I_{2\times 2} & \{N_{k,i,j}\}^{2} I_{2\times 2} \end{bmatrix} \begin{bmatrix} \dot{\boldsymbol{u}}_{i} \\ \dot{\boldsymbol{u}}_{j} \\ \dot{\boldsymbol{u}}_{k} \end{bmatrix}$$

where  $I_{2\times 2}$  represents  $2 \times 2$  identical matrix. For sake of simplicity, assume that density  $\rho$  and thickness h are constants. Then,

$$T_{i,j,k} = \frac{1}{2} \begin{bmatrix} \dot{\boldsymbol{u}}_{i}^{\top} & \dot{\boldsymbol{u}}_{j}^{\top} & \dot{\boldsymbol{u}}_{k}^{\top} \end{bmatrix} \rho h \begin{bmatrix} (\triangle/6)I_{2\times2} & (\triangle/12)I_{2\times2} & (\triangle/12)I_{2\times2} \\ (\triangle/12)I_{2\times2} & (\triangle/6)I_{2\times2} & (\triangle/12)I_{2\times2} \\ (\triangle/12)I_{2\times2} & (\triangle/6)I_{2\times2} \end{bmatrix} \begin{bmatrix} \dot{\boldsymbol{u}}_{i} \\ \dot{\boldsymbol{u}}_{j} \\ \dot{\boldsymbol{u}}_{k} \end{bmatrix}$$
$$= \frac{1}{2} \begin{bmatrix} \dot{\boldsymbol{u}}_{i}^{\top} & \dot{\boldsymbol{u}}_{j}^{\top} & \dot{\boldsymbol{u}}_{k}^{\top} \end{bmatrix} \frac{\rho h \triangle}{12} \begin{bmatrix} 2I_{2\times2} & I_{2\times2} & I_{2\times2} \\ I_{2\times2} & 2I_{2\times2} & I_{2\times2} \\ I_{2\times2} & I_{2\times2} & 2I_{2\times2} \end{bmatrix} \begin{bmatrix} \dot{\boldsymbol{u}}_{i} \\ \dot{\boldsymbol{u}}_{j} \\ \dot{\boldsymbol{u}}_{k} \end{bmatrix}$$
(2.3.5)

where  $\triangle = \triangle P_i P_j P_k$  (see Problem 6). Matrix

$$M_{i,j,k} = \frac{\rho h \Delta}{12} \begin{bmatrix} 2I_{2\times2} & I_{2\times2} & I_{2\times2} \\ I_{2\times2} & 2I_{2\times2} & I_{2\times2} \\ I_{2\times2} & I_{2\times2} & 2I_{2\times2} \end{bmatrix}$$
(2.3.6)

is referred to as *partial inertia matrix*. Note that the sum of all blocks of matrix  $M_{i,j,k}$  is equal to  $\rho h \triangle I_{2 \times 2}$ , which denotes the mass of this triangular element.

Let us calculate the total kinetic energy over rectangle region  $\Box P_1P_3P_6P_4$  shown in Fig. 2.2. This region consists of four triangle regions:  $\triangle P_1P_2P_4$ ,  $\triangle P_2P_3P_5$ ,  $\triangle P_5P_4P_2$ , and  $\triangle P_6P_5P_3$ . For sake of simplicity, assume that  $\rho h \triangle / 12$  is constantly equal to 1. Then, partial inertia matrices are given as

$$M_{1,2,4} = M_{2,3,5} = M_{5,4,2} = M_{6,5,3} = \begin{bmatrix} 2I_{2\times2} & I_{2\times2} & I_{2\times2} \\ I_{2\times2} & 2I_{2\times2} & I_{2\times2} \\ I_{2\times2} & I_{2\times2} & 2I_{2\times2} \end{bmatrix}.$$

Let  $u_{\rm N}$  be a collective vector consisting of all displacement vectors at nodal points:

$$\boldsymbol{u}_{\mathrm{N}} = \begin{bmatrix} \boldsymbol{u}_{1} \\ \boldsymbol{u}_{2} \\ \vdots \\ \boldsymbol{u}_{6} \end{bmatrix}$$
(2.3.7)

which is referred to as *nodal displacement vector*. The total kinetic energy is then described by a quadratic form with respect to  $\dot{\boldsymbol{u}}_{N}$ :

$$T = \frac{1}{2} \dot{\boldsymbol{u}}_{\mathrm{N}}^{\top} \boldsymbol{M} \, \dot{\boldsymbol{u}}_{\mathrm{N}},$$

where M is referred to as *inertia matrix*. Noting that

 $(1,2,3) \times (1,2,3)$  blocks of  $M_{1,2,4}$  contribute to  $(1,2,4) \times (1,2,4)$  blocks of M,

namely,

(1,1), (1,2), (1,3) blocks of  $M_{1,2,4}$  contribute to (1,1), (1,2), (1,4) blocks of M, (2,1), (2,2), (2,3) blocks of  $M_{1,2,4}$  contribute to (2,1), (2,2), (2,4) blocks of M, (3,1), (3,2), (3,3) blocks of  $M_{1,2,4}$  contribute to (4,1), (4,2), (4,4) blocks of M,

we find contribution of  $M_{1,2,4}$  to M as follows:

$2I_{2\times 2}$	$I_{2\times 2}$	$I_{2\times 2}$	-	]
$I_{2\times 2}$	$2I_{2\times 2}$	$I_{2\times 2}$		
$I_{2\times 2}$	$I_{2\times 2}$	$2I_{2\times 2}$		.

Similarly,

 $(1,2,3) \times (1,2,3)$  blocks of  $M_{5,4,2}$  contribute to  $(5,4,2) \times (5,4,2)$  blocks of M,

namely,

(1, 1), (1, 2), (1, 3) blocks of  $M_{5,4,2}$  contribute to (5, 5), (5, 4), (5, 2) blocks of M, (2, 1), (2, 2), (2, 3) blocks of  $M_{5,4,2}$  contribute to (4, 5), (4, 4), (4, 2) blocks of M, (3, 1), (3, 2), (3, 3) blocks of  $M_{5,4,2}$  contribute to (2, 5), (2, 4), (2, 2) blocks of M,

we find contribution of  $M_{5,4,2}$  to M as follows:

ſ	-				
		$2I_{2\times 2}$	$I_{2\times 2}$	$I_{2\times 2}$	
		$I_{2\times 2}$	$2I_{2\times 2}$	$I_{2\times 2}$	.
		$I_{2\times 2}$	$I_{2\times 2}$	$2I_{2\times 2}$	
L	_				

Summing up all contributions, we finally have

$$M = \begin{bmatrix} 2I_{2\times2} & I_{2\times2} & I_{2\times2} \\ I_{2\times2} & 6I_{2\times2} & I_{2\times2} & 2I_{2\times2} & 2I_{2\times2} \\ & I_{2\times2} & 4I_{2\times2} & 2I_{2\times2} & I_{2\times2} \\ I_{2\times2} & 2I_{2\times2} & 4I_{2\times2} & I_{2\times2} \\ & & 2I_{2\times2} & 2I_{2\times2} & I_{2\times2} & 6I_{2\times2} & I_{2\times2} \\ & & & I_{2\times2} & & I_{2\times2} & 2I_{2\times2} \end{bmatrix}$$

This inertia matrix M consists of  $6^2~2\times 2$  blocks and is a sparse matrix. We simply describe the above calculation as

$$M = M_{1,2,4} \oplus M_{2,3,5} \oplus M_{5,4,2} \oplus M_{6,5,3}.$$
(2.3.8)

.

Operator  $\oplus$  works block-wise. In general, inertia matrix is described as

$$M = \bigoplus_{i,j,k} M_{i,j,k} \tag{2.3.9}$$

where i, j, k represent nodal point numbers of each triangle.

Finite element approximation of strain potential energy We apply the above calculation to strain potential energy. First, let us calculate strain potential energy stored in small triangle region  $\Delta P_i P_j P_k$ :

$$U_{i,j,k} = \int_{\triangle \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k} \frac{1}{2} \,\boldsymbol{\varepsilon}^\top (\lambda I_\lambda + \mu I_\mu) \boldsymbol{\varepsilon} \, h \, \mathrm{d}S.$$
(2.3.10)

Piecewise linear approximation of function  $\boldsymbol{u}$  over triangle region  $\triangle \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k$  is described as  $\boldsymbol{u} = \boldsymbol{u}_i N_{i,j,k} + \boldsymbol{u}_j N_{j,k,i} + \boldsymbol{u}_k N_{k,i,j}$ . Introducing collective vectors  $\boldsymbol{\gamma}_u = [u_i, u_j, u_k]^{\top}$  and  $\boldsymbol{\gamma}_v = [v_i, v_j, v_k]^{\top}$ , we find

$$\frac{\partial u}{\partial x} = \boldsymbol{a}^{\top} \boldsymbol{\gamma}_{u}, \qquad \frac{\partial u}{\partial y} = \boldsymbol{b}^{\top} \boldsymbol{\gamma}_{u}, \qquad \frac{\partial v}{\partial x} = \boldsymbol{a}^{\top} \boldsymbol{\gamma}_{v}, \qquad \frac{\partial v}{\partial y} = \boldsymbol{b}^{\top} \boldsymbol{\gamma}_{v}$$

where

$$\boldsymbol{a} = \frac{1}{2\Delta} \begin{bmatrix} y_j - y_k \\ y_k - y_i \\ y_i - y_j \end{bmatrix}, \qquad \boldsymbol{b} = \frac{-1}{2\Delta} \begin{bmatrix} x_j - x_k \\ x_k - x_i \\ x_i - x_j \end{bmatrix}$$
(2.3.11)

(see Problem 2). Then, strain vector is given as

$$\boldsymbol{\varepsilon} = \begin{bmatrix} \boldsymbol{a}^{\top} \boldsymbol{\gamma}_{u} \\ \boldsymbol{b}^{\top} \boldsymbol{\gamma}_{v} \\ \boldsymbol{b}^{\top} \boldsymbol{\gamma}_{u} + \boldsymbol{a}^{\top} \boldsymbol{\gamma}_{v} \end{bmatrix}$$
(2.3.12)

Substituting the above equation into eq. (2.3.10), we have

$$U_{i,j,k} = \frac{1}{2} \begin{bmatrix} \boldsymbol{\gamma}_{u}^{\top} & \boldsymbol{\gamma}_{v}^{\top} \end{bmatrix} \lambda \begin{bmatrix} \boldsymbol{a}\boldsymbol{a}^{\top} & \boldsymbol{a}\boldsymbol{b}^{\top} \\ \boldsymbol{b}\boldsymbol{a}^{\top} & \boldsymbol{b}\boldsymbol{b}^{\top} \end{bmatrix} h \triangle \begin{bmatrix} \boldsymbol{\gamma}_{u} \\ \boldsymbol{\gamma}_{v} \end{bmatrix} \\ + \frac{1}{2} \begin{bmatrix} \boldsymbol{\gamma}_{u}^{\top} & \boldsymbol{\gamma}_{v}^{\top} \end{bmatrix} \mu \begin{bmatrix} 2\boldsymbol{a}\boldsymbol{a}^{\top} + \boldsymbol{b}\boldsymbol{b}^{\top} & \boldsymbol{b}\boldsymbol{a}^{\top} \\ \boldsymbol{a}\boldsymbol{b}^{\top} & 2\boldsymbol{b}\boldsymbol{b}^{\top} + \boldsymbol{a}\boldsymbol{a}^{\top} \end{bmatrix} h \triangle \begin{bmatrix} \boldsymbol{\gamma}_{u} \\ \boldsymbol{\gamma}_{v} \end{bmatrix}$$
(2.3.13)

(see Problem 7). Then, we have

$$U_{i,j,k} = \frac{1}{2} \boldsymbol{\gamma}^{\top} \left( \lambda H_{\lambda} + \mu H_{\mu} \right) \boldsymbol{\gamma}$$
(2.3.14)

where

$$egin{aligned} oldsymbol{\gamma} &= \left[ egin{aligned} oldsymbol{\gamma}_u \ oldsymbol{\gamma}_v \end{array} 
ight], \quad &H_\lambda = \left[ egin{aligned} oldsymbol{aa}^ op oldsymbol{ab}^ op \ oldsymbol{ba}^ op oldsymbol{bb}^ op \end{array} 
ight] hrangle, \ &H_\mu = \left[ egin{aligned} oldsymbol{2aa}^ op + oldsymbol{bb}^ op \ oldsymbol{ab}^ op \ oldsymbol{ab}^ op \ oldsymbol{2bb}^ op + oldsymbol{aa}^ op \end{array} 
ight] hrangle, \end{aligned}$$

The above equation is a quadratic form with respect to  $\boldsymbol{\gamma} = [u_i, u_j, u_k, v_i, v_j, v_k]^\top$ . Let us permutate rows and columns of  $H_{\lambda}$  and  $H_{\mu}$  so that  $U_{i,j,k}$  is described by a quadratic form with respect to  $\boldsymbol{u}_{i,j,k} = [u_i, v_i, u_j, v_j, u_k, v_k]^\top$ . Namely, let 1, 4, 2, 5, 3, 6 rows and columns of  $H_{\lambda}$  be 1, 2, 3, 4, 5, 6 rows and columns of  $J_{\lambda}^{i,j,k}$ . Similarly, let 1, 4, 2, 5, 3, 6 rows and columns of  $H_{\mu}$  be 1, 2, 3, 4, 5, 6 rows and columns of  $J_{\mu}^{i,j,k}$ . Then, we have

$$\boldsymbol{\gamma}^{\top} H_{\lambda} \boldsymbol{\gamma} = \boldsymbol{u}_{i,j,k}^{\top} J_{\lambda}^{i,j,k} \boldsymbol{u}_{i,j,k}, \qquad \boldsymbol{\gamma}^{\top} H_{\mu} \boldsymbol{\gamma} = \boldsymbol{u}_{i,j,k}^{\top} J_{\mu}^{i,j,k} \boldsymbol{u}_{i,j,k}$$

Matrices  $J_{\lambda}^{i,j,k}$  and  $J_{\mu}^{i,j,k}$  are referred to as *partial connection matrices*. Once coordinates of  $P_i, P_j, P_k$  are given, we can calculate partial connection matrices  $J_{\lambda}^{i,j,k}$  and  $J_{\mu}^{i,j,k}$ .

Finally, we find strain potential energy stored in  $\triangle P_i P_j P_k$ :

$$U_{i,j,k} = \frac{1}{2} \boldsymbol{u}_{i,j,k}^{\top} K_{i,j,k} \, \boldsymbol{u}_{i,j,k}$$
(2.3.15)

where

$$K_{i,j,k} = \lambda J_{\lambda}^{i,j,k} + \mu J_{\mu}^{i,j,k}$$
(2.3.16)

is referred to as *partial stiffness matrix*.

Summing up all strain potential energies over small triangle regions, we obtain the total strain potential energy described as

$$U = \frac{1}{2} \boldsymbol{u}_{\mathrm{N}}^{\top} K \, \boldsymbol{u}_{\mathrm{N}}$$
(2.3.17)

where

$$K = \bigoplus_{i,j,k} K_{i,j,k}$$
(2.3.18)

is referred to as *stiffness matrix*. Assuming that Lamé's constants  $\lambda$  and  $\mu$  are uniform over the region, stiffness matrix is described as

$$K = \bigoplus_{i,j,k} \left( \lambda J_{\lambda}^{i,j,k} + \mu J_{\mu}^{i,j,k} \right) = \lambda \bigoplus_{i,j,k} J_{\lambda}^{i,j,k} + \mu \bigoplus_{i,j,k} J_{\mu}^{i,j,k}$$

which directly yields

$$K = \lambda J_{\lambda} + \mu J_{\mu} \tag{2.3.19}$$

where

$$J_{\lambda} = \bigoplus_{i,j,k} J_{\lambda}^{i,j,k}, \qquad J_{\mu} = \bigoplus_{i,j,k} J_{\mu}^{i,j,k}$$

are referred to as *connection matrices*.

**Example** Let us calculate partial connection matrices of triangle  $P_1P_2P_4$  shown in Fig. 2.2. Vectors  $\boldsymbol{a}, \boldsymbol{b}$  are given by  $\boldsymbol{a} = [-1, 1, 0]^{\top}$  and  $\boldsymbol{b} = [-1, 0, 1]^{\top}$ . Assuming h = 2, we have

$H_{\lambda} =$	$\begin{bmatrix} 1 \\ -1 \\ 0 \\ 1 \end{bmatrix}$	-1 1 0	$\begin{array}{c} 0 \\ 0 \\ 0 \\ 0 \\ \end{array}$	$     \begin{array}{c}       1 \\       -1 \\       0 \\       1     \end{array} $	$\begin{array}{c} 0 \\ 0 \\ 0 \\ 0 \\ \end{array}$	-1 1 0 -1	],	$H_{\mu} = $	$\begin{bmatrix} 3\\ -2\\ -1\\ \hline 1 \end{bmatrix}$		$-1 \\ 0 \\ 1 \\ -1$	$     \begin{array}{c}       1 \\       0 \\       -1 \\       3     \end{array} $	$-1 \\ 0 \\ 1 \\ -1$	$\begin{bmatrix} 0\\0\\0\\\hline-2 \end{bmatrix}$
	$\begin{bmatrix} 1\\ 0\\ -1 \end{bmatrix}$	-1 0 1	0 0 0	$     \begin{array}{c}       1 \\       0 \\       -1     \end{array} $	0 0 0	$-1 \\ 0 \\ 1 \\ -1$			$\begin{bmatrix} 1\\ -1\\ 0 \end{bmatrix}$	0 0 0	$\begin{array}{c} -1 \\ 1 \\ 0 \end{array}$		-1 1 0	$\begin{bmatrix} -2\\0\\2 \end{bmatrix}$

Permuting rows and columns of the above matrices, we find

$$J_{\lambda}^{1,2,4} = \begin{bmatrix} 1 & 1 & -1 & 0 & 0 & -1 \\ 1 & 1 & -1 & 0 & 0 & -1 \\ \hline -1 & -1 & 1 & 0 & 0 & 1 \\ 0 & 0 & 0 & 0 & 0 & 0 \\ \hline 0 & 0 & 0 & 0 & 0 & 0 \\ \hline -1 & -1 & 1 & 0 & 0 & 1 \end{bmatrix}$$
$$J_{\mu}^{1,2,4} = \begin{bmatrix} 3 & 1 & -2 & -1 & -1 & 0 \\ 1 & 3 & 0 & -1 & -1 & -2 \\ \hline -2 & 0 & 2 & 0 & 0 & 0 \\ \hline -1 & -1 & 0 & 1 & 1 & 0 \\ \hline -1 & -1 & 0 & 1 & 1 & 0 \\ \hline 0 & -2 & 0 & 0 & 0 & 2 \end{bmatrix}$$

Let us calculate partial connection matrices of triangle  $P_5P_4P_2$  shown in Fig.2.2. Vectors  $\boldsymbol{a}, \boldsymbol{b}$  are given by  $\boldsymbol{a} = [-1, 1, 0]^{\top}$  and  $\boldsymbol{b} = [-1, 0, 1]^{\top}$ . Thus, assuming h = 2, we find  $J_{\lambda}^{5,4,2} = J_{\lambda}^{1,2,4}$  and  $J_{\mu}^{5,4,2} = J_{\mu}^{1,2,4}$ . Partial connection matrices are invariant with respect to translation displacement. As a result, under the same assumption, we have

$$J_{\lambda}^{1,2,4} = J_{\lambda}^{2,3,5} = J_{\lambda}^{5,4,2} = J_{\lambda}^{6,5,3}, \qquad J_{\mu}^{1,2,4} = J_{\mu}^{2,3,5} = J_{\mu}^{5,4,2} = J_{\mu}^{6,5,3}$$

Let us calculate connection matrices  $J_{\lambda}$  and  $J_{\mu}$  of rectangle region  $\Box P_1P_3P_6P_4$  shown in Fig. 2.2. Noting that

 $(1,2,3) \times (1,2,3)$  blocks of  $J_{\lambda}^{1,2,4}$  contribute to  $(1,2,4) \times (1,2,4)$  blocks of  $J_{\lambda}$ ,

namely,

(1,1), (1,2), (1,3) blocks of  $J_{\lambda}^{1,2,4}$  contribute to (1,1), (1,2), (1,4) blocks of  $J_{\lambda}$ , (2,1), (2,2), (2,3) blocks of  $J_{\lambda}^{1,2,4}$  contribute to (2,1), (2,2), (2,4) blocks of  $J_{\lambda}$ , (3,1), (3,2), (3,3) blocks of  $J_{\lambda}^{1,2,4}$  contribute to (4,1), (4,2), (4,4) blocks of  $J_{\lambda}$ ,

we obtain contribution of  $J_{\lambda}^{1,2,4}$  to  $J_{\lambda}$ . Noting that

 $(1,2,3) \times (1,2,3)$  blocks of  $J_{\lambda}^{5,4,2}$  contribute to  $(5,4,2) \times (5,4,2)$  blocks of  $J_{\lambda}$ , namely,

(1,1), (1,2), (1,3) blocks of  $J_{\lambda}^{5,4,2}$  contribute to (5,5), (5,4), (5,2) blocks of  $J_{\lambda}$ , (2,1), (2,2), (2,3) blocks of  $J_{\lambda}^{5,4,2}$  contribute to (4,5), (4,4), (4,2) blocks of  $J_{\lambda}$ , (3,1), (3,2), (3,3) blocks of  $J_{\lambda}^{5,4,2}$  contribute to (2,5), (2,4), (2,2) blocks of  $J_{\lambda}$ ,

we obtain contribution of  $J_{\lambda}^{5,4,2}$  to  $J_{\lambda}$ . Summing up all contributions, we finally have

	1	1	-1	0			0	-1				-
	1	1	-1	0			0	-1				
	-1	-1	2	1	-1	0	0	1	0	-1		
	0	0	1	2	-1	0	1	0	-1	-2		
			-1	-1	1	0			0	1	0	0
τ			0	0	0	1			1	0	-1	-1
$J \chi = 1$												
• A	0	0	0	1			1	0	-1	-1		
• 7	$0 \\ -1$	$0 \\ -1$	$\begin{array}{c} 0 \\ 1 \end{array}$	$\begin{array}{c} 1\\ 0\end{array}$			$\begin{array}{c} 1\\ 0\end{array}$	$\begin{array}{c} 0 \\ 1 \end{array}$	$     \begin{array}{c}       -1 \\       0     \end{array} $	$-1 \\ 0$		
~ /	$0 \\ -1$	$0 \\ -1$	$\begin{array}{c} 0 \\ 1 \\ 0 \end{array}$	$     \begin{array}{c}       1 \\       0 \\       -1     \end{array} $	0	1	$     \begin{array}{c}       1 \\       0 \\       -1     \end{array} $	$\begin{array}{c} 0 \\ 1 \\ \hline 0 \end{array}$			-1	-1
~ ~	$0 \\ -1$	$0 \\ -1$	$\begin{array}{c} 0 \\ 1 \\ 0 \\ -1 \end{array}$	$ \begin{array}{r}1\\0\\-1\\-2\end{array} $	0 1	1 0	$     \begin{array}{r}       1 \\       0 \\       -1 \\       -1     \end{array} $	$\begin{array}{c} 0 \\ 1 \\ \hline 0 \\ 0 \\ \end{array}$	$\begin{array}{r} -1 \\ 0 \\ \hline 2 \\ 1 \end{array}$	$ \begin{array}{r} -1 \\ 0 \\ \hline 1 \\ 2 \end{array} $	$-1 \\ 0$	$-1 \\ 0$
~ ~	0 	0 -1	$     \begin{array}{c}       0 \\       1 \\       0 \\       -1     \end{array} $	$     \begin{array}{c}       1 \\       0 \\       -1 \\       -2     \end{array} $	0 1 0		$     \begin{array}{c}       1 \\       0 \\       -1 \\       -1 \\       -1     \end{array} $	0 1 0 0	$     \begin{array}{r}       -1 \\       0 \\       2 \\       1 \\       -1     \end{array} $	$     \begin{array}{r}       -1 \\       0 \\       1 \\       2 \\       0     \end{array} $	$-1 \\ 0 \\ 1$	$-1 \\ 0 \\ 1$

We simply describe the above calculation as

$$J_{\lambda} = J_{\lambda}^{1,2,4} \oplus J_{\lambda}^{2,3,5} \oplus J_{\lambda}^{5,4,2} \oplus J_{\lambda}^{6,5,3}.$$
 (2.3.20)

Operator  $\oplus$  works block-wise. Similarly, we have

$$J_{\mu} = J_{\mu}^{1,2,4} \oplus J_{\mu}^{2,3,5} \oplus J_{\mu}^{5,4,2} \oplus J_{\mu}^{6,5,3},$$
(2.3.21)

which yields

	3	1	-2	-1			-1	0				1
	1	3	0	-1			-1	-2				
	-2	0	6	1	-2	-1	0	1	-2	-1		
	-1	-1	1	6	0	-1	1	0	-1	-4		
	-		-2	0	3	0			0	1	-1	-1
τ			-1	-1	0	3			1	0	0	-2
$J_{\mu} =$	-1	-1	0	1			3	0	-2	0		
	0	-2	1	0			0	3	-1	-1		
			-2	-1	0	1	-2	-1	6	1	-2	0
			-1	-4	1	0	0	-1	1	6	-1	-1
					-1	0			-2	-1	3	1
					-1	-2			0	-1	1	3

Matrices  $J_{\lambda}$  and  $J_{\mu}$  are sparse matrices.

### 2.4 Three-dimensional finite element approximation

Strain potential energy and kinetic energy of a three-dimensional soft robot are formulated by integrals over three-dimensional region V, which is often described by an irregular shape, making analytical calculation of integrals difficult or impossible. Let us approximate threedimensional region V by a set of small tetrahedra. Then, integral over three-dimensional region V can be approximated by the sum of integrals over small tetrahedra:

$$\int_{V} \approx \sum_{\mathbf{Q} \mathbf{P}_{i} \mathbf{P}_{j} \mathbf{P}_{k} \mathbf{P}_{l}} \int_{\mathbf{Q} \mathbf{P}_{i} \mathbf{P}_{j} \mathbf{P}_{k} \mathbf{P}_{l}}$$

Here we apply piecewise linear approximation so that we can analytically or numerically calculate individual integrals over small tetrahedra.

Finite element approximation of kinetic energy Let us calculate kinetic energy over tetrahedron region  $\langle P_i P_j P_k P_l \rangle$ :

$$T_{i,j,k,l} = \int_{\Diamond \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k \mathbf{P}_l} \frac{1}{2} \rho \, \dot{\boldsymbol{u}}^\top \dot{\boldsymbol{u}} \, \mathrm{d}V$$
(2.4.1)

Piecewise linear approximation of function  $\boldsymbol{u}$  over tetrahedron region  $\langle P_i P_j P_k P_l$  is described as follows:

$$\boldsymbol{u} = \boldsymbol{u}_i \, N_{i,j,k,l} + \boldsymbol{u}_j \, N_{j,k,l,i} + \boldsymbol{u}_k \, N_{k,l,i,j} + \boldsymbol{u}_l \, N_{l,k,i,j}.$$
(2.4.2)

Differentiating the above equation with respect to time t, we have

$$\dot{\boldsymbol{u}} = \dot{\boldsymbol{u}}_i \, N_{i,j,k,l} + \dot{\boldsymbol{u}}_j \, N_{j,k,l,i} + \dot{\boldsymbol{u}}_k \, N_{k,l,i,j} + \dot{\boldsymbol{u}}_l \, N_{l,k,i,j}.$$
(2.4.3)

For sake of simplicity, assume that density  $\rho$  is constant. Letting  $I_{3\times 3}$  represent  $3\times 3$  identical matrix, we have

$$T_{i,j,k,l} = \frac{1}{2} \begin{bmatrix} \dot{\boldsymbol{u}}_{i}^{\top} & \dot{\boldsymbol{u}}_{j}^{\top} & \dot{\boldsymbol{u}}_{k}^{\top} & \dot{\boldsymbol{u}}_{l}^{\top} \end{bmatrix} \frac{\rho \Diamond}{20} \begin{bmatrix} 2I_{3\times3} & I_{3\times3} & I_{3\times3} & I_{3\times3} \\ I_{3\times3} & 2I_{3\times3} & I_{3\times3} & I_{3\times3} \\ I_{3\times3} & I_{3\times3} & I_{3\times3} & I_{3\times3} \\ I_{3\times3} & I_{3\times3} & I_{3\times3} & I_{3\times3} \end{bmatrix} \begin{bmatrix} \dot{\boldsymbol{u}}_{i} \\ \dot{\boldsymbol{u}}_{k} \\ \dot{\boldsymbol{u}}_{l} \end{bmatrix}$$
(2.4.4)

where  $\Diamond = \Diamond \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k \mathbf{P}_l$  (see Problem 8). Matrix

$$M_{i,j,k,l} = \frac{\rho \Diamond}{20} \begin{bmatrix} 2I_{3\times3} & I_{3\times3} & I_{3\times3} & I_{3\times3} \\ I_{3\times3} & 2I_{3\times3} & I_{3\times3} & I_{3\times3} \\ I_{3\times3} & I_{3\times3} & 2I_{3\times3} & I_{3\times3} \\ I_{3\times3} & I_{3\times3} & I_{3\times3} & 2I_{3\times3} \end{bmatrix}$$
(2.4.5)

is referred to as *partial inertia matrix*. Note that the sum of all blocks of matrix  $M_{i,j,k,l}$  is equal to  $\rho \Diamond I_{3\times 3}$ , which denotes the mass of this tetrahedron element.

Summing up all kinetic energies over small tetrahedron regions, we obtain the total kinetic energy described as

$$T = \frac{1}{2} \dot{\boldsymbol{u}}_{\mathrm{N}}^{\top} M \, \dot{\boldsymbol{u}}_{\mathrm{N}},$$

where M is referred to as *inertia matrix*.

Finite element approximation of strain potential energy We calculate strain potential energy stored in small tetrahedron region  $\langle P_i P_j P_k P_l$ . Introducing collective vectors  $\gamma_u = [u_i, u_j, u_k, u_l]^\top$ ,  $\gamma_v = [v_i, v_j, v_k, v_k]^\top$ , and  $\gamma_w = [w_i, w_j, w_k, w_k]^\top$ , we find

$$\begin{aligned} \frac{\partial u}{\partial x} &= \boldsymbol{a}^{\top} \boldsymbol{\gamma}_{u}, \qquad \frac{\partial u}{\partial y} = \boldsymbol{b}^{\top} \boldsymbol{\gamma}_{u}, \qquad \frac{\partial u}{\partial z} = \boldsymbol{c}^{\top} \boldsymbol{\gamma}_{u}, \\ \frac{\partial v}{\partial x} &= \boldsymbol{a}^{\top} \boldsymbol{\gamma}_{v}, \qquad \frac{\partial v}{\partial y} = \boldsymbol{b}^{\top} \boldsymbol{\gamma}_{v}, \qquad \frac{\partial v}{\partial z} = \boldsymbol{c}^{\top} \boldsymbol{\gamma}_{v}, \\ \frac{\partial w}{\partial x} &= \boldsymbol{a}^{\top} \boldsymbol{\gamma}_{w}, \qquad \frac{\partial w}{\partial y} = \boldsymbol{b}^{\top} \boldsymbol{\gamma}_{w}, \qquad \frac{\partial w}{\partial z} = \boldsymbol{c}^{\top} \boldsymbol{\gamma}_{w} \end{aligned}$$

where

$$\boldsymbol{a} = \frac{1}{6\Diamond} \begin{bmatrix} -a_{j,k,l} \\ a_{k,l,i} \\ -a_{l,i,j} \\ a_{i,j,k} \end{bmatrix}, \quad \boldsymbol{b} = \frac{1}{6\Diamond} \begin{bmatrix} -b_{j,k,l} \\ b_{k,l,i} \\ -b_{l,i,j} \\ b_{i,j,k} \end{bmatrix}, \quad \boldsymbol{c} = \frac{1}{6\Diamond} \begin{bmatrix} -c_{j,k,l} \\ c_{k,l,i} \\ -c_{l,i,j} \\ c_{i,j,k} \end{bmatrix}$$

with

$$\begin{aligned} a_{j,k,l} &= (y_j z_k - y_k z_j) + (y_k z_l - y_l z_k) + (y_l z_j - y_j z_l) \\ b_{j,k,l} &= (z_j x_k - z_k x_j) + (z_k x_l - z_l x_k) + (z_l x_j - z_j x_l) \\ c_{j,k,l} &= (x_j y_k - x_k y_j) + (x_k y_l - x_l y_k) + (x_l y_j - x_j y_l) \end{aligned}$$

(see Problem 3). Strain vector is given as

$$arepsilon = \left[egin{array}{c} a^ op \gamma_u \ b^ op \gamma_v \ c^ op \gamma_w \ c^ op \gamma_w + b^ op \gamma_w \ a^ op \gamma_w + c^ op \gamma_u \ b^ op \gamma_u + a^ op \gamma_v \end{array}
ight]$$

Then, strain potential energy stored in  $\langle P_i P_i P_k P_l$  is given by

$$U_{i,j,k,l} = \frac{1}{2} \boldsymbol{\gamma}^{\top} \left( \lambda H_{\lambda} + \mu H_{\mu} \right) \boldsymbol{\gamma}$$
(2.4.6)

where

$$\begin{split} \boldsymbol{\gamma} &= \left[ \begin{array}{c} \boldsymbol{\gamma}_{u} \\ \boldsymbol{\gamma}_{v} \\ \boldsymbol{\gamma}_{w} \end{array} \right], \quad \boldsymbol{H}_{\lambda} &= \left[ \begin{array}{cc} \boldsymbol{a}\boldsymbol{a}^{\top} & \boldsymbol{a}\boldsymbol{b}^{\top} & \boldsymbol{a}\boldsymbol{c}^{\top} \\ \boldsymbol{b}\boldsymbol{a}^{\top} & \boldsymbol{b}\boldsymbol{b}^{\top} & \boldsymbol{b}\boldsymbol{c}^{\top} \\ \boldsymbol{c}\boldsymbol{a}^{\top} & \boldsymbol{c}\boldsymbol{b}^{\top} & \boldsymbol{c}\boldsymbol{c}^{\top} \end{array} \right] \boldsymbol{\Diamond}, \\ \boldsymbol{H}_{\mu} &= \left[ \begin{array}{cc} 2\boldsymbol{a}\boldsymbol{a}^{\top} + \boldsymbol{b}\boldsymbol{b}^{\top} + \boldsymbol{c}\boldsymbol{c}^{\top} & \boldsymbol{b}\boldsymbol{a}^{\top} & \boldsymbol{c}\boldsymbol{c}^{\top} \\ \boldsymbol{a}\boldsymbol{b}^{\top} & 2\boldsymbol{b}\boldsymbol{b}^{\top} + \boldsymbol{c}\boldsymbol{c}^{\top} + \boldsymbol{a}\boldsymbol{a}^{\top} & \boldsymbol{c}\boldsymbol{b}^{\top} \\ \boldsymbol{a}\boldsymbol{c}^{\top} & \boldsymbol{b}\boldsymbol{c}^{\top} & 2\boldsymbol{c}\boldsymbol{c}^{\top} + \boldsymbol{a}\boldsymbol{a}^{\top} + \boldsymbol{b}\boldsymbol{b}^{\top} \end{array} \right] \boldsymbol{\Diamond}. \quad (2.4.7) \end{split}$$

Let us permutate rows and columns of  $H_{\lambda}$  and  $H_{\mu}$  so that  $U_{i,j,k,l}$  is described by a quadratic form of  $\boldsymbol{u}_{i,j,k,l} = [\boldsymbol{u}_i^{\top}, \boldsymbol{u}_j^{\top}, \boldsymbol{u}_k^{\top}, \boldsymbol{u}_l^{\top}]^{\top}$ . Namely, let 1, 5, 9, 2, 6, 10, 3, 7, 11, 4, 8, 12 rows and columns of  $H_{\lambda}$  be 1 through 12 rows and columns of  $J_{\lambda}^{i,j,k,l}$ . Similarly, let 1, 5, 9, 2, 6, 10, 3, 7, 11, 4, 8, 12 rows and columns of  $H_{\mu}$  be 1 through 12 rows and columns of  $J_{\mu}^{i,j,k,l}$ . Then, we have

$$\boldsymbol{\gamma}^{ op} H_{\lambda} \boldsymbol{\gamma} = \boldsymbol{u}_{i,j,k,l}^{ op} \, J_{\lambda}^{i,j,k,l} \, \boldsymbol{u}_{i,j,k,l}, \qquad \boldsymbol{\gamma}^{ op} H_{\mu} \boldsymbol{\gamma} = \boldsymbol{u}_{i,j,k,l}^{ op} \, J_{\mu}^{i,j,k,l} \, \boldsymbol{u}_{i,j,k,l}.$$

Matrices  $J_{\lambda}^{i,j,k,l}$  and  $J_{\mu}^{i,j,k,l}$  are referred to as *partial connection matrices*. Once coordinates of  $P_i$ ,  $P_j$ ,  $P_k$ ,  $P_l$  are given, we can calculate partial connection matrices  $J_{\lambda}^{i,j,k,l}$  and  $J_{\mu}^{i,j,k,l}$ . Finally, we find strain potential energy stored in  $\Diamond P_i P_j P_k P_l$ :

$$U_{i,j,k,l} = \frac{1}{2} \boldsymbol{u}_{i,j,k,l}^{\top} K_{i,j,k,l} \, \boldsymbol{u}_{i,j,k,l}$$
(2.4.8)

where

$$K_{i,j,k,l} = \lambda J_{\lambda}^{i,j,k,l} + \mu J_{\mu}^{i,j,k,l}$$
(2.4.9)

which is referred to as *partial stiffness matrix*.

Summing up all strain potential energies over small tetrahedron regions, we obtain the total strain potential energy described as

$$U = \frac{1}{2} \boldsymbol{u}_{\mathrm{N}}^{\mathsf{T}} K \, \boldsymbol{u}_{\mathrm{N}} \tag{2.4.10}$$



Figure 2.3: Example of regular quadrangular pyramid

where

$$K = \bigoplus_{i,j,k,l} K_{i,j,k,l} \tag{2.4.11}$$

is referred to as stiffness matrix. Assuming that Lamé's constants  $\lambda$  and  $\mu$  are uniform over the region, stiffness matrix is described as

$$K = \lambda J_{\lambda} + \mu J_{\mu} \tag{2.4.12}$$

where matrices  $J_{\lambda}$  and  $J_{\mu}$  are referred to as *connection matrices*.

Let us calculate connection matrices of a regular quadrangular pyramid (Fig. 2.3). The base of the pyramid is square  $\Box P_1 P_2 P_3 P_4$  and the apex of the pyramid is  $P_5$ . Coordinates of vertices are given by  $\boldsymbol{x}_1 = [0, 0, 0]^{\top}$ ,  $\boldsymbol{x}_2 = [2, 0, 0]^{\top}$ ,  $\boldsymbol{x}_3 = [2, 2, 0]^{\top}$ ,  $\boldsymbol{x}_4 = [0, 2, 0]^{\top}$ , and  $\boldsymbol{x}_5 = [1, 1, 1]^{\top}$ . The piramid consists of two tetraheda:  $\langle P_1 P_2 P_3 P_5$  and  $\langle P_3 P_4 P_1 P_5$ .

Partial connection matrices of  $\langle P_1 P_2 P_3 P_5$  are as follows:

Note that  $(1,2,3,4) \times (1,2,3,4)$  blocks of  $J^{1,2,3,5}_{\lambda}$  and  $J^{1,2,3,5}_{\mu}$  contribute to  $(1,2,3,5) \times (1,2,3,5)$  blocks of  $J_{\lambda}$  and  $J_{\mu}$ 

Partial connection matrices of  $\Diamond P_3 P_4 P_1 P_5$  are as follows:

	1	0	-1	-1	1	0 0	-1	-1	0 (	) 2	1	
	0	0	0	0	0	0 0	0	0	0 (	0 0		
	-1	0	1	1	-1	0 0	1	1	0 (	) -2	2	
	-1	0	1	1	-1	0 0	1	1	0 (	) -2	2	
	1	0	-1	-1	1	0 0	-1	-1	0 (	) 2		
$_{7^{3,4,1,5}}$ 1	0	0	0	0	0	0 0	0	0	0 (	0 0	1	
$J_{\lambda} = \frac{1}{4}$	0	0	0	0	0	0 0	0	0	0 (	0 0	_	
	-1	0	1	1	$^{-1}$	0 0	1	1	0 (	) -2	2	
	-1	0	1	1	$^{-1}$	0 0	1	1	0 (	) -2	2	
	0	0	0	0	0	0 0	0	0	0 (	0 0	_	
	0	0	0	0	0	0 0	0	0	0 (	0 0		
	2	0	-2	-2	2	0 0	-2	-2	0 (	) 4		
	3	0	-1	-2	0	1	1	0	0	-2	0	0
	$\begin{bmatrix} 3\\ 0 \end{bmatrix}$	$\begin{array}{c} 0 \\ 2 \end{array}$	$-1 \\ 0$	$\begin{vmatrix} -2\\ 1 \end{vmatrix}$	$0 \\ -1$	$1 \\ -1$	$\begin{vmatrix} 1 \\ -1 \end{vmatrix}$	$\begin{array}{c} 0 \\ 1 \end{array}$	$\begin{array}{c} 0 \\ 1 \end{array}$	$\begin{vmatrix} -2\\ 0 \end{vmatrix}$	$0 \\ -2$	0 0
	$\begin{bmatrix} 3 \\ 0 \\ -1 \end{bmatrix}$	$\begin{array}{c} 0 \\ 2 \\ 0 \end{array}$	$-1 \\ 0 \\ 3$	$\begin{vmatrix} -2\\ 1\\ 0 \end{vmatrix}$	$     \begin{array}{c}       0 \\       -1 \\       0     \end{array} $	$     1 \\     -1 \\     -1 $	$\begin{vmatrix} 1\\ -1\\ -1 \end{vmatrix}$	$egin{array}{c} 0 \ 1 \ 0 \end{array}$	$\begin{array}{c} 0 \\ 1 \\ 2 \end{array}$	$\begin{vmatrix} -2 \\ 0 \\ 2 \end{vmatrix}$	$\begin{array}{c} 0 \\ -2 \\ 0 \end{array}$	$\begin{array}{c} 0 \\ 0 \\ -4 \end{array}$
	$\begin{bmatrix} 3\\0\\-1\\-2\end{bmatrix}$	$\begin{array}{c} 0 \\ 2 \\ 0 \\ \hline 1 \end{array}$	$\begin{array}{r} -1 \\ 0 \\ 3 \\ \hline 0 \end{array}$	$\begin{vmatrix} -2\\1\\0\\3 \end{vmatrix}$			$\begin{vmatrix} 1 \\ -1 \\ -1 \\ -1 \end{vmatrix}$	0 1 0 0	$\begin{array}{c} 0\\ 1\\ 2\\ \end{array}$	$\begin{array}{ c c } -2 \\ 0 \\ 2 \\ 0 \\ \end{array}$	$\begin{array}{c} 0 \\ -2 \\ 0 \\ \hline 0 \end{array}$	$\begin{array}{c} 0 \\ 0 \\ -4 \\ \hline 0 \end{array}$
	$ \begin{bmatrix} 3 \\ 0 \\ -1 \\ -2 \\ 0 \end{bmatrix} $	$     \begin{array}{c}       0 \\       2 \\       0 \\       \hline       1 \\       -1     \end{array} $	$\begin{array}{c} -1 \\ 0 \\ 3 \\ \hline 0 \\ 0 \\ \end{array}$	$\begin{vmatrix} -2\\ 1\\ 0\\ -1 \end{vmatrix}$	$     \begin{array}{c}       0 \\       -1 \\       0 \\       -1 \\       3     \end{array} $	$ \begin{array}{c} 1 \\ -1 \\ -1 \\ 0 \\ 0 \end{array} $	$\begin{vmatrix} 1 \\ -1 \\ -1 \\ -1 \\ 1 \end{vmatrix}$	$\begin{array}{c} 0\\ 1\\ 0\\ \hline 0\\ -2 \end{array}$	$\begin{array}{c} 0\\ 1\\ 2\\ \hline 0\\ 0\\ \end{array}$	$\begin{vmatrix} -2\\0\\2\\0\\0 \end{vmatrix}$	$\begin{array}{c} 0\\ -2\\ 0\\ \hline 0\\ 0\\ 0 \end{array}$	$\begin{array}{c} 0 \\ 0 \\ -4 \\ \hline 0 \\ 0 \\ \end{array}$
1 <sup>3,4,1,5</sup> _ 1	$ \begin{bmatrix} 3 \\ 0 \\ -1 \\ -2 \\ 0 \\ 1 \end{bmatrix} $	$\begin{array}{c} 0 \\ 2 \\ 0 \\ \hline 1 \\ -1 \\ -1 \end{array}$	$ \begin{array}{r} -1 \\ 0 \\ 3 \\ \hline 0 \\ 0 \\ -1 \end{array} $	$ \begin{array}{c c} -2 \\ 1 \\ 0 \\ -1 \\ 0 \end{array} $		$ \begin{array}{c} 1 \\ -1 \\ -1 \\ 0 \\ 0 \\ 2 \end{array} $	$ \begin{array}{c c} 1 \\ -1 \\ -1 \\ 1 \\ 1 \\ 1 \end{array} $	$     \begin{array}{c}       0 \\       1 \\       0 \\       -2 \\       -1     \end{array} $	$     \begin{array}{c}       0 \\       1 \\       2 \\       0 \\       0 \\       -1 \\     \end{array} $	$ \begin{array}{ c c } -2 \\ 0 \\ 2 \\ 0 \\ 0 \\ -2 \\ \end{array} $	$\begin{array}{c} 0\\ -2\\ 0\\ \hline 0\\ 0\\ 2\\ \end{array}$	$\begin{array}{c} 0\\ 0\\ -4\\ \hline 0\\ 0\\ 0\\ 0\\ \end{array}$
$J^{3,4,1,5}_{\mu} = \frac{1}{4}$	$ \begin{array}{c} 3 \\ 0 \\ -1 \\ -2 \\ 0 \\ 1 \\ -1 \\ -2 \\ 0 \\ 1 \end{array} $	$\begin{array}{c} 0 \\ 2 \\ 0 \\ \hline 1 \\ -1 \\ -1 \\ -1 \\ -1 \end{array}$	$ \begin{array}{r} -1 \\ 0 \\ 3 \\ \hline 0 \\ 0 \\ -1 \\ \hline -1 \end{array} $	$ \begin{array}{c c} -2 \\ 1 \\ 0 \\ -1 \\ 0 \\ -1 \end{array} $	$ \begin{array}{c} 0 \\ -1 \\ 0 \\ \hline -1 \\ 3 \\ 0 \\ \hline 1 \end{array} $	$ \begin{array}{cccc} 1 \\ -1 \\ 0 \\ 0 \\ 2 \\ 1 \end{array} $	$ \begin{array}{c c} 1 \\ -1 \\ -1 \\ 1 \\ 1 \\ 2 \\ \end{array} $	$ \begin{array}{c} 0 \\ 1 \\ 0 \\ -2 \\ -1 \\ 0 \end{array} $	$ \begin{array}{c} 0 \\ 1 \\ 2 \\ 0 \\ 0 \\ -1 \\ 0 \end{array} $	$ \begin{array}{c c} -2 \\ 0 \\ 2 \\ 0 \\ 0 \\ -2 \\ -2 \end{array} $	$ \begin{array}{c} 0 \\ -2 \\ 0 \\ 0 \\ 2 \\ 0 \end{array} $	$\begin{array}{c} 0\\ 0\\ -4\\ 0\\ 0\\ 0\\ 0\\ 0\\ \end{array}$
$J^{3,4,1,5}_{\mu} = \frac{1}{4}$	$   \begin{bmatrix}     3 \\     0 \\     -1 \\     -2 \\     0 \\     1 \\     0   \end{bmatrix}   $	$ \begin{array}{c} 0 \\ 2 \\ 0 \\ -1 \\ -1 \\ -1 \\ 1 \end{array} $	$ \begin{array}{r} -1 \\ 0 \\ 3 \\ 0 \\ -1 \\ -1 \\ 0 \\ \end{array} $	$ \begin{array}{c c} -2 \\ 1 \\ 0 \\ -1 \\ 0 \\ -1 \\ 0 \end{array} $	$\begin{array}{c} 0 \\ -1 \\ 0 \\ \hline -1 \\ 3 \\ 0 \\ \hline 1 \\ -2 \end{array}$	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	$ \begin{array}{c c} 1 \\ -1 \\ -1 \\ 1 \\ 1 \\ 2 \\ 0 \\ \end{array} $	$ \begin{array}{c} 0 \\ 1 \\ 0 \\ -2 \\ -1 \\ 0 \\ 3 \end{array} $	$\begin{array}{c} 0 \\ 1 \\ 2 \\ 0 \\ 0 \\ -1 \\ 0 \\ 1 \end{array}$	$ \begin{array}{c c} -2 \\ 0 \\ 2 \\ 0 \\ -2 \\ 0 \end{array} $	$ \begin{array}{c} 0 \\ -2 \\ 0 \\ 0 \\ 2 \\ 0 \\ -2 \end{array} $	$ \begin{array}{c} 0 \\ 0 \\ -4 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \end{array} $
$J^{3,4,1,5}_{\mu} = \frac{1}{4}$	$ \begin{array}{c} 3 \\ 0 \\ -1 \\ -2 \\ 0 \\ 1 \\ 0 \\ 0 \end{array} $	$\begin{array}{c} 0 \\ 2 \\ 0 \\ \hline 1 \\ -1 \\ -1 \\ -1 \\ 1 \\ 1 \end{array}$	$ \begin{array}{r} -1 \\ 0 \\ 3 \\ 0 \\ 0 \\ -1 \\ 0 \\ 2 \end{array} $	$ \begin{array}{c c} -2 \\ 1 \\ 0 \\ -1 \\ 0 \\ -1 \\ 0 \\ 0 \end{array} $	$ \begin{array}{c} 0 \\ -1 \\ 0 \\ \hline -1 \\ 0 \\ \hline -1 \\ 3 \\ 0 \\ \hline 1 \\ -2 \\ 0 \\ \end{array} $	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	$ \begin{array}{c c} 1 \\ -1 \\ -1 \\ 1 \\ 1 \\ 2 \\ 0 \\ 0 \\ \end{array} $	$\begin{array}{c} 0 \\ 1 \\ 0 \\ -2 \\ -1 \\ 0 \\ 3 \\ 1 \end{array}$	$\begin{array}{c} 0 \\ 1 \\ 2 \\ 0 \\ 0 \\ -1 \\ 0 \\ 1 \\ 3 \end{array}$	$ \begin{array}{c c} -2 \\ 0 \\ 2 \\ 0 \\ -2 \\ 0 \\ 0 \\ 0 \end{array} $	$\begin{array}{c} 0 \\ -2 \\ 0 \\ 0 \\ 2 \\ 0 \\ -2 \\ -2 \end{array}$	$ \begin{array}{c} 0 \\ -4 \\ 0 \\ 0 \\ 0 \\ 0 \\ -4 \\ \end{array} $
$J^{3,4,1,5}_{\mu} = \frac{1}{4}$	$ \begin{array}{c} 3 \\ 0 \\ -1 \\ -2 \\ 0 \\ 1 \\ 0 \\ 0 \\ -2 \end{array} $	$\begin{array}{c} 0 \\ 2 \\ 0 \\ -1 \\ -1 \\ -1 \\ 1 \\ 1 \\ 0 \end{array}$	$ \begin{array}{c} -1 \\ 0 \\ 3 \\ 0 \\ -1 \\ -1 \\ 0 \\ 2 \\ 2 \end{array} $	$ \begin{array}{c c} -2 \\ 1 \\ 0 \\ -1 \\ 0 \\ -1 \\ 0 \\ 0 \\ 0 \\ 0 \end{array} $	$ \begin{array}{c} 0 \\ -1 \\ 0 \\ -1 \\ 3 \\ 0 \\ 1 \\ -2 \\ 0 \\ 0 \end{array} $	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	$ \begin{array}{c c} 1 \\ -1 \\ -1 \\ 1 \\ 1 \\ 2 \\ 0 \\ 0 \\ -2 \\ \end{array} $	$\begin{array}{c} 0 \\ 1 \\ 0 \\ -2 \\ -1 \\ 0 \\ 3 \\ 1 \\ 0 \\ \end{array}$	$ \begin{array}{c} 0 \\ 1 \\ 2 \\ 0 \\ -1 \\ 0 \\ 1 \\ 3 \\ 0 \end{array} $	$ \begin{array}{c c} -2 \\ 0 \\ 2 \\ 0 \\ -2 \\ 0 \\ 0 \\ 4 \\ \end{array} $	$\begin{array}{c} 0 \\ -2 \\ 0 \\ 0 \\ 2 \\ 0 \\ -2 \\ -2 \\ 0 \\ \end{array}$	$\begin{array}{c} 0 \\ 0 \\ -4 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ -4 \\ 0 \\ \end{array}$
$J^{3,4,1,5}_{\mu} = \frac{1}{4}$	$ \begin{bmatrix} 3 \\ 0 \\ -1 \\ -2 \\ 0 \\ 1 \\ 0 \\ 0 \\ -2 \\ 0 \end{bmatrix} $	$\begin{array}{c} 0 \\ 2 \\ 0 \\ -1 \\ -1 \\ -1 \\ 1 \\ 1 \\ 0 \\ -2 \end{array}$	$ \begin{array}{c} -1 \\ 0 \\ 3 \\ 0 \\ -1 \\ 0 \\ 2 \\ 0 \\ \end{array} $	$ \begin{array}{c c} -2 \\ 1 \\ 0 \\ -1 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \end{array} $	$ \begin{array}{c} 0 \\ -1 \\ 0 \\ -1 \\ 0 \\ -1 \\ 0 \\ -1 \\ 3 \\ 0 \\ 1 \\ -2 \\ 0 \\ 0 \\ 0 \\ 0 \\ 0 \end{array} $	$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	$ \begin{array}{c c} 1 \\ -1 \\ -1 \\ 1 \\ 1 \\ 2 \\ 0 \\ 0 \\ -2 \\ 0 \\ \end{array} $	$ \begin{array}{c} 0 \\ 1 \\ 0 \\ -2 \\ -1 \\ 0 \\ 3 \\ 1 \\ 0 \\ -2 \end{array} $	$\begin{array}{c} 0 \\ 1 \\ 2 \\ 0 \\ 0 \\ -1 \\ 0 \\ 1 \\ 3 \\ 0 \\ -2 \end{array}$	$ \begin{array}{c c} -2 \\ 0 \\ 2 \\ 0 \\ -2 \\ 0 \\ 0 \\ 4 \\ 0 \end{array} $	$\begin{array}{c} 0 \\ -2 \\ 0 \\ 0 \\ 2 \\ 0 \\ -2 \\ -2 \\ 0 \\ 4 \\ \end{array}$	$ \begin{array}{c} 0 \\ -4 \\ 0 \\ 0 \\ 0 \\ 0 \\ -4 \\ 0 \\ 0 \\ 0 \end{array} $

Note that  $(1,2,3,4) \times (1,2,3,4)$  blocks of  $J^{3,4,1,5}_{\lambda}$  and  $J^{3,4,1,5}_{\mu}$  contribute to  $(3,4,1,5) \times (3,4,1,5)$  blocks of  $J_{\lambda}$  and  $J_{\mu}$ 

Synthesizing the above partial connection matrices yields the following connection matrices:

	[ 1	0	1	-1	1	0	0	-1	1	0	0	0	0	0	$-2^{-1}$
	0	1	1	0	0	0	-1	0	1	1	$^{-1}$	0	0	0	-2
	1	1	2	-1	1	0	-1	-1	2	1	-1	0	0	0	-4
	-1	0	-1	1	-1	0	0	1	-1				0	0	2
	1	0	1	-1	1	0	0	-1	1				0	0	-2
	0	0	0	0	0	0	0	0	0				0	0	0
1	0	-1	-1	0	0	0	1	0	-1	-1	1	0	0	0	2
$J_{\lambda} = \frac{1}{4}$	-1	0	-1	1	-1	0	0	1	-1	0	0	0	0	0	2
4	1	1	2	-1	1	0	-1	-1	2	1	-1	0	0	0	-4
	0	1	1				-1	0	1	1	-1	0	0	0	-2
	0	-1	-1				1	0	-1	-1	1	0	0	0	2
	0	0	0				0	0	0	0	0	0	0	0	0
	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	0	0	0	0	0	0	0	0	0	0	0	0	0	0	0
	-2	-2	-4	2	-2	0	2	2	-4	-2	2	0	0	0	8

	5	0	1	-2	0	-1	2	-1	-1	-1	1	1	-4	0	0 -
	0	5	1	1	-1	1	-1	2	-1	0	-2	$^{-1}$	0	-4	0
	1	1	6	0	0	-1	1	1	4	0	0	-1	-2	-2	-8
	-2	1	0	3	-1	0	-1	0	0				0	0	0
	0	-1	0	-1	3	0	1	-2	0				0	0	0
	-1	1	-1	0	0	2	-1	1	-1				2	-2	0
1	2	-1	1	-1	1	-1	5	0	-1	-2	0	1	-4	0	0
$J_{\mu} = \frac{1}{4}$	-1	2	1	0	-2	1	0	5	-1	1	-1	$^{-1}$	0	-4	0
4	-1	-1	4	0	0	-1	-1	-1	6	0	0	-1	2	2	-8
	-1	0	0				-2	1	0	3	-1	0	0	0	0
	1	-2	0				0	-1	0	-1	3	0	0	0	0
	1	-1	-1				1	-1	-1	0	0	2	-2	2	0
	-4	0	-2	0	0	2	-4	0	2	0	0	-2	8	0	0
	0	-4	-2	0	0	-2	0	-4	2	0	0	2	0	8	0
	0	0	-8	0	0	0	0	0	-8	0	0	0	0	0	16

Since either tetrahedron does not include both  $P_2$  and  $P_4$ , (2,4) and (4,2) blocks of the connection matrices are zero matrices.

#### 2.5 Implementation

Two-dimensional finite element calculation was implemented on MATLAB. Classes Body, Triangle, and NodalPoint were introduced. Class Body defines a two-dimensional body, which consists of an array of triangles and an array of nodal points. Class Triangle specifies a triangle, including three numbers of nodal points. Class NodalPoint defines a nodal point, including its two coordinates.

For example, rectangle region in Fig. 2.2 consists of 6 nodal points and 4 triangles. Coordinates of individual nodal points are listed as

points =  $\begin{bmatrix} x_1 & x_2 & x_3 & x_4 & x_5 & x_6 \end{bmatrix} = \begin{bmatrix} 0 & 1 & 2 & 0 & 1 & 2 \\ 0 & 0 & 0 & 1 & 1 & 1 \end{bmatrix}$ 

Nodal point numbers for individual triangular elements are listed as

$$\text{triangles} = \begin{bmatrix} 1 & 2 & 4 \\ 2 & 3 & 5 \\ 5 & 4 & 2 \\ 6 & 5 & 3 \end{bmatrix},$$

which implies that  $\triangle_1 = \triangle P_1 P_2 P_4$ ,  $\triangle_2 = \triangle P_2 P_3 P_5$ ,  $\triangle_3 = \triangle P_5 P_4 P_2$ , and  $\triangle_4 = \triangle P_6 P_5 P_3$ . The rectangle region is then given by

elastic = Body(6, points, 4, triangles, thickness);

where thickness specifies thickness h of the two-dimensional body.

Instance of class Triangle includes geometric propertices such as nodal point numbers, area, and thickness as well as physical parameters such as density and Lamé's constants. Class Triangle involves the following methods:

**partial\_derivaties** calculating partial derivatives  $\partial u/\partial x$ ,  $\partial u/\partial y$ ,  $\partial v/\partial x$ ,  $\partial v/\partial y$ **calculate\_Cauchy\_strain** calculating Cauchy strain in the triangle **partial\_strain\_potential\_energy** strain potential energy stored in the triangle calculate\_Green\_strain calculating Green strain in the triangle partial\_strain\_potential\_energy\_Green\_strain strain energy using Green strain partial\_gravitational\_potential\_energy gravitational energy stored in the triangle partial\_stiffness\_matrix calculating partial stiffness matrix  $K_{i,j,k}$ partial\_inertia\_matrix calculating partial inertia matrix  $M_{i,j,k}$ partial\_gravitational\_vector calculating partial gravitational vector  $g_{i,j,k}$ 

Class Body involves the following methods:

total\_strain\_potential\_energy calculating strain energy stored in the body total\_strain\_potential\_energy\_Green\_strain strain energy using Green strain total\_gravitational\_potential\_energy gravitational energy stored in the body calculate\_stiffness\_matrix calculating stiffness matrix K calculate\_inertia\_matrix calculating inertia matrix M calculate\_gravitational\_vector calculating gravitational vector g constraint\_matrix constraint matrix when specified nodal points are fixed draw draw the shape of the body

Assuming that density  $\rho$  and Lamé's constants  $\lambda$ ,  $\mu$  are uniform over the region, the following specifies these parameters:

elastic = elastic.mechanical\_parameters(rho, lambda, mu);

The following calculates the stiffness and inertia matrices:

```
elastic = elastic.calculate_stiffness_matrix;
elastic = elastic.calculate_inertia_matrix;
```

The stiffness and inertia matrices are then referred by

M = elastic.Inertia\_Matrix; K = elastic.Stiffness\_Matrix;

which can be applied to static and dynamic calculation of the motion and deformation of a soft body.

Three-dimensional finite element calculation was implemented on MATLAB. Classes Body, Tetrahedron, and NodalPoint were introduced. Class Body defines a three-dimensional body, which consists of an array of tetrahedra and an array of nodal points. Class Tetrahedron specifies a tetrahedron, including four numbers of nodal points. Class NodalPoint defines a nodal point, including its three coordinates.

For example, a regular quadrangular pyramid (Fig. 2.3) consists of 5 nodal points and 2 tetrahedra. Coordinates of individual nodal points are listed as

	0	2	2	0	1
points = $\begin{bmatrix} x_1 & x_2 & x_3 & x_4 & x_5 \end{bmatrix}$ =	0	0	2	2	1
	0	0	0	0	1

Nodal point numbers for individual tetrehedron elements are listed as

 $\text{tetrahedra} = \left[ \begin{array}{rrrr} 1 & 2 & 3 & 5 \\ 3 & 4 & 1 & 5 \end{array} \right],$ 

which implies that  $\Diamond_1 = \Diamond P_1 P_2 P_3 P_5$  and  $\Diamond_2 = \Diamond P_3 P_4 P_1 P_5$ . The quadrangular pyramid is then given by

elastic = Body(5, points, 2, tetrahedra);

followed by methods to define physical parameters and calculate inertia and stiffness matrices.

## Problems

- 1. Show eqs. (2.1.3a)(2.1.3b)(2.1.3c).
- 2. Calculate partial derivatives of piecewise linear approximation  $L_{i,j,k}(x,y)$  given in eq. (2.1.4) with respect to x, y.
- 3. Calculate partial derivatives of piecewise linear approximation  $L_{i,j,k,l}(x, y, z)$  given in eq. (2.1.6) with respect to x, y, z.
- 4. Show the following equations:

$$\int_{x_i}^{x_j} N_{i,j}(x) N_{i,j}(x) dx = \int_{x_i}^{x_j} N_{j,i}(x) N_{j,i}(x) dx = \frac{1}{3}(x_j - x_i)$$
$$\int_{x_i}^{x_j} N_{i,j}(x) N_{j,i}(x) dx = \int_{x_i}^{x_j} N_{j,i}(x) N_{i,j}(x) dx = \frac{1}{6}(x_j - x_i)$$

Letting  $L_{i,j}(x) = f_i N_{i,j}(x) + f_j N_{j,i}(x)$ , show

$$\int_{x_i}^{x_j} \left\{ L_{i,j}(x) \right\}^2 \, \mathrm{d}x = \left[ \begin{array}{cc} f_i & f_j \end{array} \right] \frac{x_j - x_i}{6} \left[ \begin{array}{cc} 2 & 1 \\ 1 & 2 \end{array} \right] \left[ \begin{array}{cc} f_i \\ f_j \end{array} \right]$$

5. Show the following equations:

$$\int_{x_i}^{x_j} N'_{i,j}(x) N'_{i,j}(x) \, \mathrm{d}x = \int_{x_i}^{x_j} N'_{j,i}(x) N'_{j,i}(x) \, \mathrm{d}x = \frac{1}{x_j - x_i}$$
$$\int_{x_i}^{x_j} N'_{i,j}(x) N'_{j,i}(x) \, \mathrm{d}x = \int_{x_i}^{x_j} N'_{j,i}(x) N'_{i,j}(x) \, \mathrm{d}x = \frac{-1}{x_j - x_i}$$

Letting  $L_{i,j}(x) = f_i N_{i,j}(x) + f_j N_{j,i}(x)$ , show

$$\int_{x_i}^{x_j} \left\{ L'_{i,j}(x) \right\}^2 \, \mathrm{d}x = \left[ \begin{array}{cc} f_i & f_j \end{array} \right] \frac{1}{x_j - x_i} \left[ \begin{array}{cc} 1 & -1 \\ -1 & 1 \end{array} \right] \left[ \begin{array}{cc} f_i \\ f_j \end{array} \right]$$

6. Show the following equations:

$$\int_{\Delta} N_{i,j,k}^2 \, \mathrm{d}S = \int_{\Delta} N_{j,k,i}^2 \, \mathrm{d}S = \int_{\Delta} N_{k,i,j}^2 \, \mathrm{d}S = \frac{\Delta}{6}$$
$$\int_{\Delta} N_{i,j,k} \, N_{j,k,i} \, \mathrm{d}S = \int_{\Delta} N_{j,k,i} \, N_{k,i,j} \, \mathrm{d}S = \int_{\Delta} N_{k,i,j} \, N_{i,j,k} \, \mathrm{d}S = \frac{\Delta}{12}$$

where  $\triangle = \triangle \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k$ .

- 7. Show eq. (2.3.13).
- 8. Show the following equations:

$$\int_{\Diamond} N_{i,j,k,l}^{2} \, \mathrm{d}V = \dots = \frac{\Diamond}{10}$$
$$\int_{\Diamond} N_{i,j,k,l} N_{j,k,l,i} \, \mathrm{d}V = \dots = \frac{\Diamond}{20}$$

where  $\Diamond = \Diamond \mathbf{P}_i \mathbf{P}_j \mathbf{P}_k \mathbf{P}_l$ .