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Steady State Conduction with Boundary Convection

Let us now modify our one-dimensional, steady-state heat conduction model by replacing the applied heat flux boundary condition with a convective heat transfer, hence

$$L(T) = -\frac{d}{dx} \left(k(x) \frac{dT}{dx} \right) - s(x) = 0 \qquad \Omega \in (0, L)$$
 (1a)

$$I(T) = k(x)\frac{dT}{dn} + h(T - T_r) = 0 \qquad \partial\Omega_1 \in 0$$
 (1b)

$$T = T_b \qquad \partial \Omega_2 \in L$$
 (1c)

Step 1: Assume an approximate solution exists

$$T(x) \approx T^{N}(x) = \sum_{a=1}^{N} \Psi_{a}(x) Q_{a} = \Psi_{a}(x) Q_{a} \quad \text{for} \quad 1 \le a \le N$$
 (2)

Step 2: What are suitable trial functions and how do we solve for expansion coefficients?

Step 3: Solve for expansion coefficients by

a. Defining the error in the approximation

$$e^{N}(x) = T(x) - T^{N}(x)$$
(3a)

b. Measuring the error

$$L(e^{N}) = L(T) - L(T^{N}) = -L(T^{N})$$
(3b)

c. Minimizing the error

$$WS^{N} = \int_{\Omega} \Phi_{\mathbf{b}}(x) L(T^{N}) dx \equiv 0 \quad \text{for} \quad 1 \le \mathbf{b} \le N$$
 (3c)

d. Optimizing the minimization

$$GWS^{N} = \int_{\Omega} \Psi_{\mathbf{b}}(x) L(T^{N}) dx = 0 \qquad \text{for} \quad 1 \le \mathbf{b} \le N$$
 (3d)

e. Applying the operator

$$GWS^{N} = \int_{\Omega} \Psi_{b}(x) \left(-\frac{d}{dx} \left(k(x) \frac{dT^{N}}{dx} \right) - s(x) \right) dx = 0$$
 (3e)

f. Integrate by parts

$$GWS^{N} = \int_{\Omega} \frac{d\Psi_{\boldsymbol{b}}(x)}{dx} k(x) \frac{dT^{N}}{dx} dx - \int_{\Omega} \Psi_{\boldsymbol{b}}(x) s(x) dx - \Psi_{\boldsymbol{b}}(x) k(x) \frac{dT^{N}}{dx} \bigg|_{\partial\Omega} = 0 \quad \text{for} \quad 1 \le \boldsymbol{b} \le N$$

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We must now pay close attention to our boundary conditions before moving forward.

Expanding the boundary condition term

$$k(x)\frac{dT^{N}}{dx}\bigg|_{\partial\Omega} = \left(+k(x)\frac{dT^{N}}{dx}\bigg|_{\partial\Omega_{2}}\right) - \left(-k(x)\frac{dT^{N}}{dx}\bigg|_{\partial\Omega_{1}}\right)$$
$$= k(x)\frac{dT^{N}}{dx}\bigg|_{\partial\Omega_{2}} - h(T^{N} - T_{r})\bigg|_{\partial\Omega_{1}}$$

Substituting back

$$GWS^{N} = \int_{\Omega} \frac{d\Psi_{\boldsymbol{b}}(x)}{dx} k(x) \frac{dT^{N}}{dx} dx - \int_{\Omega} \Psi_{\boldsymbol{b}}(x) s(x) dx + \Psi_{\boldsymbol{b}}(x) h(T^{N} - T_{r}) \Big|_{\partial \Omega_{1}} - \Psi_{\boldsymbol{b}}(x) k(x) \frac{dT^{N}}{dx} \Big|_{\partial \Omega_{2}} = 0$$
 (3f)

g. Substitute the series expansion

$$GWS^{N} = \int_{\Omega} \frac{d\Psi_{b}(x)}{dx} k(x) \frac{d\Psi_{a}(x)}{dx} dx Q_{a} - \int_{\Omega} \Psi_{b}(x) s(x) dx + \Psi_{b}(x) h(\Psi_{a}(x)Q_{a} - T_{r})\Big|_{\partial\Omega_{1}} - \Psi_{b}(x) k(x) \frac{dT^{N}}{dx}\Big|_{\partial\Omega_{2}} = 0$$

Step 4: Select Lagrange polynomials as our trial functions

Step5: Discretize the problem domain

$$\Psi_{\mathbf{a}}(x)Q_{\mathbf{a}} = \mathbf{S}_{e}\left\{\left[N_{k}(\mathbf{z}_{i})\right]^{T}\left\{Q\right\}_{e}\right\}$$
(5a)

$$\Psi_{\mathbf{h}} = S_{e}(\{N_{k}(\mathbf{z}_{i})\}) \tag{5b}$$

Step 6: Form discrete *GWS* h

$$GWS^{h} = S_{e} \left(\int_{\Omega_{e}} \frac{d\{N_{k}\}}{dx} k(x) \frac{d\{N_{k}\}^{T}}{dx} d\overline{x} \{Q\}_{e} - \int_{\Omega_{e}} \{N_{k}\} s(x) d\overline{x} + \{N_{k}\} h(\{N_{k}\}^{T} \{Q\}_{e} - T_{r})_{\partial\Omega_{1}} - \{N_{k}\} k(x) \frac{dT^{N}}{dx} \Big|_{\partial\Omega_{2}} = \{0\} \right)$$

a. Get all the unknowns on the left

$$GWS^{h} = S_{e} \left(\int_{\Omega_{e}} k(x) \frac{d\{N_{k}\}}{dx} \frac{d\{N_{k}\}^{T}}{dx} d\overline{x} \{Q\}_{e} + h\{N_{k}\}\{N_{k}\}^{T} \{Q\}_{e} \Big|_{\partial\Omega_{1}} \right) = \int_{\Omega_{e}} \{N_{k}\}s(x) d\overline{x} + hT_{r}\{N_{k}\}_{\partial\Omega_{1}} + \{N_{k}\}k(x) \frac{dT^{N}}{dx} \Big|_{\partial\Omega_{2}}$$

b. Give each term a name and decide what to do with distributed and boundary terms

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Integrated terms:

$$[DIFF]_{e} = \int_{\Omega_{e}} k(x) \frac{d\{N_{k}\}}{dx} \frac{d\{N_{k}\}^{T}}{dx} d\overline{x}$$

$$\approx \{K\}_{e}^{T} \int_{\Omega_{e}} \{N_{k}\} \frac{d\{N_{k}\}}{dx} \frac{d\{N_{k}\}^{T}}{dx} d\overline{x}$$

$$= ()()_{e} \{K\}_{e}^{T} (-1)[A3011k]\{\}_{e}$$

$$(6a)$$

$$\left\{ \operatorname{SRC} \right\}_{e} \equiv \int_{\Omega_{e}} \left\{ N_{k} \right\} s(x) d\overline{x}$$

$$\approx \int_{\Omega_{e}} \left\{ N_{k} \right\} \left\{ N_{k} \right\}^{T} d\overline{x} \left\{ S \right\}_{e}$$

$$= \left(\right) \left(\right)_{e} \left\{ \right\}_{e}^{T} \left(1 \right) \left[\operatorname{A200k} \right] \left\{ S \right\}_{e}$$

$$(6b)$$

Boundary terms:

$$[HBC]_e = h\{N_k\}\{N_k\}^T\{Q\}_e\Big|_{\partial\Omega_1}$$
(6c)

$$\left\{ \mathsf{HTR} \right\}_e \equiv h T_r \left\{ N_k \right\}_{\partial \Omega_1} \tag{6d}$$

$$\left\{ BFLX \right\}_e = \left\{ N_k \right\} k(x) \frac{dT^N}{dx} \bigg|_{\partial \Omega_2}$$
 (6e)

What new stuff do we have?

Interpolated thermal conductivity

Examining the three bases within the integral for [DIFF]_e, we see an *incorrect* matrix multiplication!

$$\{K\}_{e}^{T} \int_{\Omega} \{N_{k}\} \frac{d\{N_{k}\}}{dx} \frac{d\{N_{k}\}^{T}}{dx} d\overline{x}$$

The undifferentiated basis, however, multiplies the (constant) nodal values of conductivity and yields a scalar on the element $\{K\}_e^T\{N_k\}$. Assuming a linear basis, $[\mathsf{DIFF}]_e$ becomes

$$\left\{K\right\}_{e}^{T} \int_{\Omega_{e}} \left\{N_{1}\right\} \frac{1}{l_{e}^{2}} \begin{bmatrix} 1 & -1 \\ -1 & 1 \end{bmatrix} d\overline{x}$$

Distributing the undifferentiated basis yields the hypermatrix

$$\{K\}_{e}^{T} \int_{\Omega_{e}} \frac{1}{t_{e}^{2}} \begin{bmatrix} \{N_{1}\} & -\{N_{1}\} \\ -\{N_{1}\} & \{N_{1}\} \end{bmatrix} d\overline{x}$$

where it is understood that $\{K\}_e^T$ multiplies every value of $\{N_1\}$

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Boundary terms

Drawing a simple, three-element domain, we note that boundary conditions are applied exactly on the boundaries. They are not interpolated over the interior. Our GWS^h boundary terms $[HBC]_e$, $\{HTR\}_e$, and $\{BFLX\}_e$ must therefore employ a physically consistent, reduced basis function. Sticking with the linear basis for demonstration, the distribution of basis functions becomes

Examining the leftmost element, the leftmost boundary conditions are applied via $z_1 = 1$ and $z_2 = 0$ at the leftmost node. The leftmost boundary conditions are applied nowhere else, hence $z_1 = z_2 = 0$ on all other nodes.

Examining the rightmost element, the rightmost boundary conditions are applied via $z_1 = 0$ and $z_2 = 1$ at the rightmost node. The rightmost boundary conditions are applied nowhere else, hence $z_1 = z_2 = 0$ on all other nodes.

Hence, on the boundaries, our basis functions reduce to a binary switch - "on" at the appropriate node and "off" at all other nodes. This is conveniently expressed with the Kronecker delta function

$$\mathbf{d}_{ij} = 1$$
 for $i = j$
 $\mathbf{d}_{ij} = 0$ for $i \neq j$

The boundary basis functions thus become **linear**

$$\left\{ N_{1}\right\} _{\partial\Omega}\equiv\left\{ \begin{matrix} \boldsymbol{d}_{e1}\\ \boldsymbol{d}_{eM} \end{matrix}\right\}$$

quadratic

$$\left\{ N_2 \right\}_{\partial \Omega} \equiv \left\{ \begin{matrix} \boldsymbol{d}_{e1} \\ \boldsymbol{0} \\ \boldsymbol{d}_{eM} \end{matrix} \right\}$$

cubic

$$\left\{ N_{3} \right\}_{\partial \Omega} \equiv \left\{ \begin{matrix} \boldsymbol{d}_{e1} \\ 0 \\ 0 \\ \boldsymbol{d}_{eM} \end{matrix} \right\}$$

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Thus, the square matrix $[HBC]_e$ will "assemble" to an nnodes x nnodes matrix of zeros with the convective heat transfer coefficient h in either the (1,1) position (left BC) or the (nnodes, nnodes) position (right BC). The column vector $\{HTR\}_e$ will assemble to an nnodes x 1 column of zeros with reference heat transfer hT_r in either the (1,1) position or the (nnodes, 1) position.

Why can we ignore the column vector $\{BFLX\}_{e}$?