

A METHOD FOR SIMULATING ELECTRICALLY MOVING  
TEMPERATURE FIELDS

L. A. Kozdoba

UDC 536.12:681.142.334

A simplified method, by comparison with that of [1], is given for simulating moving temperature fields using ohmic resistance electric network simulators.

The general form of the Kirchhoff–Fourier equation for nonstationary thermal conductivity in a moving medium is

$$\operatorname{div}(\lambda \operatorname{grad} T) - c\gamma \left( \frac{\partial T}{\partial t} + v_x \frac{\partial T}{\partial x} + v_y \frac{\partial T}{\partial y} + v_z \frac{\partial T}{\partial z} \right) + \omega = 0, \quad (1)$$

where  $v_x, v_y, v_z$  are velocity components of the motion of the medium at a given point.

The equation of the temperature field, referred to coordinates moving with velocity  $v_0$  in the direction of the x-axis when  $\lambda, c, \gamma, v_0 = -v_x$  are constants, is a particular case of (1) but is of sufficiently wide application in studying heat-transfer processes in welding and melting, continuous sequence heat treatment, mechanical material treatment, etc. With the above assumptions, Eq. (1) can be written

$$\frac{\partial^2 T}{\partial x^2} + \frac{\partial^2 T}{\partial y^2} + \frac{\partial^2 T}{\partial z^2} - \frac{1}{a} \frac{\partial T}{\partial t} + \frac{v_0}{a} \frac{\partial T}{\partial x} + \frac{\omega}{\lambda} = 0. \quad (2)$$

The initial and boundary conditions of the I–IV kind can be written in the same way as for a fixed coordinate system.

In [1] a method was given for simulating the solution of (2) electrically on ohmic resistance networks (R-networks), when the following substitution was used to deduce the parameters of the R-network:

$$T = u\varphi, \quad (3)$$

where

$$\varphi = \exp\left(-\frac{v_0 x}{2a}\right). \quad (4)$$

Then, from (2) we can obtain

$$\frac{\partial^2 u}{\partial x^2} + \frac{\partial^2 u}{\partial y^2} + \frac{\partial^2 u}{\partial z^2} - \frac{v^2}{4a^2} u - \frac{1}{a} \frac{\partial u}{\partial t} + \frac{\omega}{\lambda\varphi} = 0. \quad (5)$$

If we put

$$T = U\Phi, \quad (6)$$

where

$$\Phi = \exp\left(-\frac{v_0 x}{2a} - \frac{v_0^2 t}{4a}\right), \quad (7)$$

from (2) we can obtain the simpler equation

---

Kiev Institute of Technical Thermophysics, Academy of Sciences of the Ukrainian SSR. Translated from *Inzhenerno-Fizicheskii Zhurnal*, Vol. 18, No. 1, pp. 167–171, January, 1970. Original article submitted March 31, 1969.

© 1972 Consultants Bureau, a division of Plenum Publishing Corporation, 227 West 17th Street, New York, N. Y. 10011. All rights reserved. This article cannot be reproduced for any purpose whatsoever without permission of the publisher. A copy of this article is available from the publisher for \$15.00.

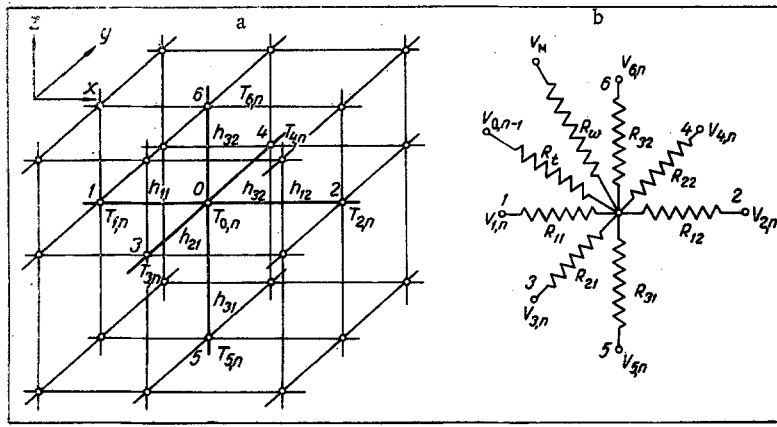


Fig. 1. Diagram of an elementary volume (a) and the nodes of an R-network for simulating the solution of (2) electrically.

$$\frac{\partial^2 U}{\partial x^2} + \frac{\partial^2 U}{\partial y^2} + \frac{\partial^2 U}{\partial z^2} - \frac{1}{a} \frac{\partial U}{\partial t} + \frac{\omega}{\lambda \Phi} = 0, \quad (8)$$

i.e., there is no additional heat sink ( $(v_0^2/4a^2)u$ ), which simplifies the construction and use of an electrical model. The boundary conditions of the I-IV kind correspondingly take the form

$$U_{s,n} = U_{me,n}, \quad (9)$$

$$\frac{\partial U_s}{\partial n} \Phi_s + \frac{\partial \Phi_s}{\partial n} U_s + \frac{q}{\lambda} = 0, \quad (10)$$

$$\frac{\partial U_s}{\partial n} \Phi_s + \frac{\partial \Phi_s}{\partial n} U_s + \frac{\alpha}{\lambda} (U_{me} \Phi_{me} - U_s \Phi_s) = 0, \quad (11)$$

$$\lambda \left( \frac{\partial U_s}{\partial n} \Phi_s + \frac{\partial \Phi_s}{\partial n} U_s \right) + \lambda_s \left( \frac{\partial U_{me}}{\partial n} \Phi_s + \frac{\partial \Phi_s}{\partial n} U_s \right) = 0, \quad (12)$$

where the subscripts s and me refer to the surface and the medium. The remaining notation is in common use.

If we write (8)-(12) in finite difference form, as in [2], and compare with the expressions of Kirchhoff's law for the currents at the nodes of R-networks, we obtain an expression for calculating the parameters of the R-network which solves the system (8)-(12).

From the values of U which have been found, we can use (6) and (7) to determine T.

The determination of T can be simplified if the network parameters are computed so that we can obtain the values of T directly from the R-network.

We write Eq. (8) in finite difference form and for U we substitute its value from (6) (at the same time we multiply all the terms of the finite difference equations by  $\prod_{i=1}^3 \sum_{k=1}^2 h_{ik}$  and cancel common factors)

$$\begin{aligned} & \frac{2 \left( T_{1,n} \frac{\Phi_{0,n}}{\Phi_{1,n}} - T_{0,n} \right) \prod_{i=2,3} \sum_{k=1}^2 h_{ik}}{h_{11} \Phi_{0,n}} + \frac{2 \left( T_{2,n} \frac{\Phi_{0,n}}{\Phi_{2,n}} - T_{0,n} \right) \prod_{i=2,3} \sum_{k=1}^2 h_{ik}}{h_{12} \Phi_{0,n}} \\ & + \frac{2 (T_{3,n} - T_{0,n}) \prod_{i=1,3} \sum_{k=1}^2 h_{ik}}{h_{21} \Phi_{0,n}} + \frac{2 (T_{4,n} - T_{0,n}) \prod_{i=1,3} \sum_{k=1}^2 h_{ik}}{h_{22} \Phi_{0,n}} \\ & + \frac{2 (T_{5,n} - T_{0,n}) \prod_{i=1,2} \sum_{k=1}^2 h_{ik}}{h_{31} \Phi_{0,n}} + \frac{2 (T_{6,n} - T_{0,n}) \prod_{i=1,2} \sum_{k=1}^2 h_{ik}}{h_{32} \Phi_{0,n}} \\ & + \frac{\left( T_{0,n-1} \frac{\Phi_{0,n}}{\Phi_{0,n-1}} - T_{0,n} \right) \prod_{i=1}^3 \sum_{k=1}^2 h_{ik}}{\alpha \delta t \Phi_{0,n}} + \frac{\omega \prod_{i=1}^3 \sum_{k=1}^2 h_{ik}}{\lambda \Phi_{0,n}} = 0, \quad (13) \end{aligned}$$

where the  $h_{ik}$  are space intervals on both sides of the node 0 (see Fig. 1); the axis  $x - i = 1$ ; the axis  $y - i = 2$ ; the axis  $z - i = 3$ . The time is  $t = n\delta t$ . In general the time interval  $\delta t$  may change during the solution.

The corresponding finite difference expressions for the boundary conditions of the II and III kind are:

$$\frac{T_s - T_m \frac{\Phi_{s,n}}{\Phi_{m,n}}}{h_m} + \frac{\partial \Phi_s}{\partial n_m} \frac{T_s}{\Phi_{s,n}} + \frac{q}{\lambda} = 0, \quad (14)$$

$$\frac{T_s - T_m \frac{\Phi_{s,n}}{\Phi_{m,n}}}{h_m} + \frac{\partial \Phi_s}{\partial n_m} \frac{T_s}{\Phi_{s,n}} + \frac{\alpha}{\lambda} (T_m - T_s) = 0, \quad (15)$$

where  $h_m$  is the space interval along the normal to the surface ( $h_m$  is equal to that  $h_{ik}$  which approaches the node on the surface from the side of the body).

If there is to be similarity between (13)-(15) and the expressions for Kirchhoff's law for the currents flowing into the corresponding nodes of the R-network, it is necessary that the parameters of the R-network should be as follows:

$$\begin{aligned} R_{11} &= \frac{h_{11} \Phi_{0,n}}{2 \prod_{i=2,3} \sum_{k=1}^2 h_{ik}} R_N, & R_{32} &= \frac{h_{32} \Phi_{0,n}}{2 \prod_{i=1,2} \sum_{k=1}^2 h_{ik}} R_N, \\ R_t &= \frac{a \delta t \Phi_{0,n}}{\prod_{i=1}^3 \sum_{k=1}^2 h_{ik}} R_N, & R_w &= \frac{(V_M - V_{0,n}) K \lambda \Phi_{0,n}}{w \prod_{i=1}^3 \sum_{k=1}^2 h_{ik}} R_N, \\ R'_q &= \frac{(V_M - V_{0,n}) K \lambda R_{ih}}{h_{ik} q}, & R''_q &= \frac{\Phi_s R_{ih}}{\frac{\partial \Phi_s}{\partial n_m} h_{ik}}, \\ R'_\alpha &= \frac{\lambda R_{ih}}{\alpha h_{ik}}, & R''_\alpha &= \frac{\Phi_s R_{ih}}{\frac{\partial \Phi_s}{\partial n_m} h_{ik}}, \end{aligned} \quad (16)$$

where the resistances  $R_{11}, \dots, R_{32}$  correspond to the space intervals  $h_{11}, \dots, h_{32}$  (see Fig. 1) and simulate the thermal heat conduction resistance;  $R_t$  is the resistance through which passes the current simulating heat absorption or emission by the due to the heat capacity of an elementary volume  $\prod_{i=1}^3 \sum_{k=1}^2 h_{ik}$  in a time interval  $\delta t$ . One end of the resistance  $R_t$  is attached to the node, the other to a voltage divider from which the voltage  $V_{0,n-1}$  is taken.

At the second end of the resistance  $R_w$ , voltage  $V_M$  is applied to  $R'_q$ . If  $V_M \gg V_{0,n}$ , when  $w, q$  are heat sources, or  $V_M \ll V_{0,n}$ , when  $w, q$  are heat sinks, instead of  $(V_M - V_{0,n})$  in the expressions for  $R_w, R_q$  we may put  $V_M$ .  $K$  is the scale of the transition from temperature to voltage:

$$K = \frac{T_{\max} - T_{\min}}{V_{\max} - V_{\min}}.$$

For given boundary conditions of the I kind, voltages are applied to the nodes corresponding to  $T_s$ . At the free end of the resistance  $R'_\alpha$  a voltage is applied corresponding to the temperature  $T_{me}$  of the medium. At the free ends of the resistances  $R''_q$  and  $R''_\alpha$  the minimal voltage  $V_M = 0$  is applied. If  $T_{me} = 0$  then

$$R_\alpha = \left[ \frac{1}{R'_\alpha} + \frac{1}{R''_\alpha} \right] = R_{ih} \left[ h_m \left( \frac{\alpha}{\lambda} + \frac{1}{\Phi_s} \frac{\partial \Phi_s}{\partial n_m} \right) \right]^{-1}$$

For given boundary conditions of the IV kind, the networks simulating the bodies in contact are joined to each other.

If  $\partial \Phi_s / \partial n_m = 0$ , which occurs if the normal to the surface lies in the plane perpendicular to the  $x$ -axis, the boundary conditions of the II and III kind are simulated using only the resistances  $R'_\alpha$  and  $R'_q$ .

The accuracy of the solution depends on the ratios  $\Phi_0/\Phi_1$  and  $\Phi_0/\Phi_2$ ; the closer they are to unity, the higher is the accuracy. The ratios  $\Phi_0/\Phi_1$  and  $\Phi_0/\Phi_2 \rightarrow 1$  for  $h_{1k} = \text{const}$  and  $x \rightarrow \infty$  or  $h_{1k} \rightarrow 0$  and  $x = \text{const}$ , i.e., for network nodes with small values of  $x$  we have to choose appropriately small values of  $h_{1k}$  and for nodes with large values of  $x$ , the intervals  $h_1$  and  $h_2$  can be increased.

A similar discussion holds for  $\Phi_{0,n-1}/\Phi_{0,n}$ .

In solving problems with moving concentrated heat sources the origin is linked to the position of the sources.

The quasistationary situation, when  $\partial T/\partial t = 0$  was discussed in [1].

In a similar manner the parameters of the R-network can be deduced for the original set of equations in nondimensional form.

If it cannot be assumed that  $\lambda$ ,  $c$ ,  $\gamma$ ,  $v$  are constant, the solution of Eq. (1) in the most general formulation can be obtained by complicating the technique of solving on an R-network or by a combined model using the general method of [2], or on a special quasianalog model developed in [3, 4].

#### LITERATURE CITED

1. L.A.Kozdoba and V.I.Makhnenko, *Inzh.-Fiz. Zh.*, 4, No. 11 (1961).
2. L.A.Kozdoba, *The Electrical Simulation of Temperature Fields for Components in Marine Power Plants* [in Russian], Sudostroenie, Leningrad (1964).
3. A.V.Temnikov, B.M.Gavrilov, and V.V.Basov, *Izv. Vuzov, Priborostroenie*, No. 5 (1966).
4. A.N.Reznikov, A.V.Temnikov, N.V.Diligenskii, and B.M.Gavrilov, in: *Mathematical Simulation and the Theory of Electrical Circuits* [in Russian], No. 5, Naukova Dumka, Kiev (1967).