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Non-Fourier Heat Transfer with Phonons and Electrons in a Circular Thin Layer Surrounding a Hot Nanodevice

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Abstract: A nonlocal model for heat transfer with phonons and electrons is applied to infer the steady-state radial temperature profile in a circular layer surrounding an inner hot component. Such a profile, following by the numerical solution of the heat equation, predicts that the temperature behaves in an anomalous way, since for radial distances from the heat source smaller than the mean-free path of phonons and electrons, it increases for increasing distances. The compatibility of this temperature behavior with the second law of thermodynamics is investigated by calculating numerically the local entropy production as a function of the radial distance. It turns out that such a production is positive and strictly decreasing with the radial distance.

Keywords: temperature hump; nonlocal model for heat transfer; heat conduction with phonons and electrons

PACS Classifications: 44.10.+i; 66.70.-f; 05.70.Ln

1. Introduction

Nanotechnology has changed our vision, expectations and abilities to control all of the material world. It offers a huge amount of potential applications and solutions that in the recent past were only possible in the realm of scientific fiction. Its usefulness has been known for many years, especially

in microelectronics and, more recently, in manufacturing of nano-electromechanical-systems (NEMS): systems with less space, material, energy consumption and usually higher performance can be actually realized. In particular, current research in nanoscaled electronic devices focuses the attention on carbon nanotubes, quantum wires and quantum dots, where the additional confinement of electrons is used to make compact high-performance components.

From the theoretical point of view, the use of nanotechnologies requires revising some well-known theories, as for example the classical Fourier law. Experimental evidence, in fact, clearly shows that it is completely inefficient to describe accurately heat transport at the nanometer length scale [1–7]. Several theories have been developed to describe heat transport in nanostructured materials [8–15].

In order to analyze thermoelectric effects, in [16] the authors developed a hydrodynamic model accounting for the possibility that both electrons and phonons may contribute to the heat transport. In particular, they assumed that the local heat flux \mathbf{q} just arises from two contributions: the partial heat flux \mathbf{q}_p due to the phonons and the partial heat flux \mathbf{q}_e due to the electrons, in such a way that:

$$\mathbf{q} = \mathbf{q}_p + \mathbf{q}_e \quad (1)$$

The aforementioned model can be applied to describe situations in which both a voltage at the ends of a conductor is created by a temperature difference (Seebeck effect) and a heat flux is generated by an electric current (Peltier effect). However, for common materials, these direct conversions are not always efficient, since they only show limited thermoelectric properties, *i.e.*, very small values of the Seebeck and Peltier material parameters. In such a case, for example, the Seebeck effect is only able to generate an electric current that is negligible, and in the absence of nonlinear effects, the evolution of the partial heat fluxes \mathbf{q}_p and \mathbf{q}_e in Equation (1) is ruled by the following equations:

$$\tau_p \dot{\mathbf{q}}_p + \mathbf{q}_p = -\lambda_p \nabla T + \ell_p^2 (\nabla^2 \mathbf{q}_p + 2\nabla \nabla \cdot \mathbf{q}_p) \quad (2a)$$

$$\tau_e \dot{\mathbf{q}}_e + \mathbf{q}_e = -\lambda_e \nabla T + \ell_e^2 (\nabla^2 \mathbf{q}_e + 2\nabla \nabla \cdot \mathbf{q}_e) \quad (2b)$$

wherein $(\tau_p; \ell_p; \lambda_p)$ and $(\tau_e; \ell_e; \lambda_e)$ are the relaxation time, the mean-free path and the thermal conductivity of phonons and electrons, respectively, and T is the absolute temperature. These equations account for nonlocal effects through the mean-free paths ℓ_p and ℓ_e , and for memory effects through the relaxation times τ_p and τ_e . They are still far from the standard form of the equations of linear heat-conduction theory with nonlocal and memory effects [17], since they contain information on the partial heat fluxes, relaxation times, thermal conductivities and mean-free paths.

The system above must be coupled to the local balance of energy, which, in the absence of heat sources and with negligible electric current, reads:

$$c_v \dot{T} + \nabla \cdot \mathbf{q} = 0 \quad (3)$$

with c_v the specific heat of the conductor.

Although more general situations, allowing for different temperatures for phonons and electrons, could be considered [18,19], in the present paper, we limit ourselves to the case in which the equations above hold. As we will see later, this assumption allows us to conserve a sufficient generality and reduces the calculations to a simpler level. Under the hypotheses above, here we study the radial heat transfer

from a hot source surrounded by a circular layer and point out some non-standard characteristics of the temperature behavior.

The paper displays the following layout.

In Section 2, we evaluate the system of Equations (2) and (3) in steady states, in order to obtain the differential equation for the temperature profile. In the special case of a two-dimensional system with a radial symmetry, that equation is solved numerically, and the influence of nonlocal effects on the solution is pointed out.

In Section 3, we investigate the thermodynamic compatibility of the results obtained in Section 2 by calculating numerically the local entropy production as a function of the distance from the heat source. We show that such a quantity is positive everywhere in the system, according to the prediction of the second law of thermodynamics.

In Section 4, we draw the main conclusions and discuss possible developments of the present investigation.

2. Radial Heat Transfer from a Hot Source

In integrated circuits, the number of transistors that can be placed on a silicon chip has increased exponentially, since the developments in nanotechnology allow the transistors to shrink in size. With more transistors integrated on a single chip, running at a faster clock rate, designers have been able to improve the system performance effectively over the past few decades. Unfortunately, the design complexity, as well as the increase in the computational rate enhances the amount of produced heat per unit volume greatly, which has to be dissipated, thus being a limiting factor in current computer miniaturization. Being able to remove the surplus of produced heat may be of crucial interest in practical applications. Therefore, in the present section, we analyze the possible consequences of accounting for nonlocal effects in Equation (2). In particular, we assume that a small transistor (sketched as a hot cylindrical device) acts as a steady heat source and derive the temperature profile in the circular surrounding zone (the radius of which is r_l), wherein both the phonons and the electrons are the heat carriers (see Figure 1 for a qualitative sketch of the system we are considering). In order to emphasize the physical problem, we further assume that the inner hot device can only exchange heat with the surrounding circular layer. This is tantamount to assume that the system at hand (hot device and circular layer) is thermally isolated.

2.1. The Physical Model

In what follows, we represent the surrounding layer as a homogeneous and isotropic two-dimensional system and denote by r the radial distance of its points from the center of the hot component (see Figure 1). Since we do not have information on the constitutive nature of the hot spot, in principle, we cannot exclude that heat propagates anisotropically inside it. Thus, there will be a thin annular region, starting at $r = r_0$, in which the transition from anisotropic to isotropic propagation may occur. For the sake of simplicity, we suppose that the dimension of this thin annulus is negligible and that the isotropic propagation starts at $r = r_0$. Thus, in steady states, the heat flux in the circular layer has only the radial component $q = q(r)$.

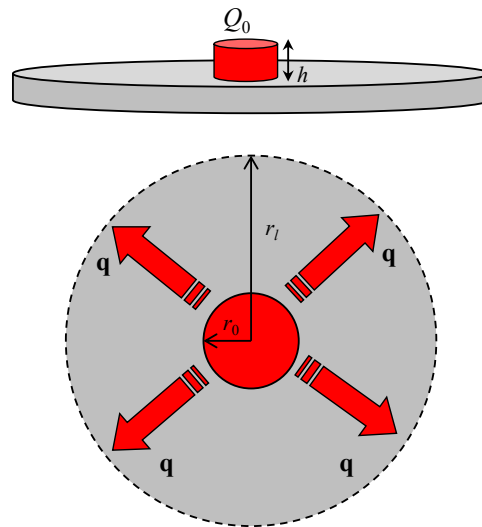


Figure 1. Cylindrical nanodevice connected to a circular surrounding layer. The internal device (*i.e.*, the red hot spot in the figure) is characterized by a radius r_0 and a thickness h . It acts as a steady heat source, with Q_0 being the total heat produced per unit time. In the circular hot spot, we assumed that only the phonons are the heat carriers. The surrounding layer (grey in the figure) is characterized, instead, by an outer radius r_l . In it, both the phonons and the electrons contribute to the heat transfer. Outside that layer, instead, the electrons stop being the heat carriers again.

To obtain the radial dependence of q , one may observe that in the absence of transversal exchanges, the total heat flowing across each concentric circular area is constant. Therefore, by considering two different concentric circular areas at a radial distance from the source equal to r and $r + dr$, the radial profile of the local heat flux (in a first-order approximation in dr) must be such that:

$$2\pi r q(r) = 2\pi [r + dr] q(r + dr) \approx 2\pi [r + dr] \left[q(r) + \frac{dq}{dr} dr \right] \Rightarrow r \frac{dq}{dr} + q(r) = 0 \tag{4}$$

The last relation in Equation (4) may be easily integrated and yields:

$$q(r) \equiv q_p(r) + q_e(r) = \frac{\Gamma}{r} \tag{5}$$

with $\Gamma = Q_0 / (2\pi h)$ as a constant value, being Q_0 the constant heat produced by the hot device per unit time.

In steady states, from Equations (1) and (3), we get:

$$\nabla \cdot (\mathbf{q}_p + \mathbf{q}_e) = 0 \tag{6}$$

On the other hand, since in our approach, \mathbf{q}_p and \mathbf{q}_e are independent state variables [4,16], the condition $\nabla \cdot \mathbf{q}_p = -\nabla \cdot \mathbf{q}_e$ cannot hold in general, but can only describe a particular situation. In the general case, instead, from Equation (6), it follows that both vectors \mathbf{q}_p and \mathbf{q}_e are divergence free. As a consequence, Equation (2) reduces to:

$$\mathbf{q}_p = -\lambda_p \nabla T + \ell_p^2 \nabla^2 \mathbf{q}_p \tag{7a}$$

$$\mathbf{q}_e = -\lambda_e \nabla T + \ell_e^2 \nabla^2 \mathbf{q}_e \tag{7b}$$

which, by means of Equation (5), produces the following evolution equation for the phonon contribution to the local heat flux:

$$\left(\frac{\lambda_e}{\lambda_p} \ell_p^2 + \ell_e^2\right) \frac{d^2 q_p}{dr^2} + \frac{1}{r} \left(\frac{\lambda_e}{\lambda_p} \ell_p^2 + \ell_e^2\right) \frac{dq_p}{dr} - \left(\frac{\lambda_e}{\lambda_p} + 1\right) q_p = \frac{\Gamma}{r} \left(\frac{\ell_e^2}{r^2} - 1\right) \tag{8}$$

This second-order differential equation can be solved by prescribing two Dirichlet-type boundary conditions. In fact, if we assume that both in the internal device ($r < r_0$) and outside of the circular layer ($r > r_l$) the phonons are the sole heat carriers, and since in steady states, Γ takes a constant value, Equation (5) yields the following boundary conditions:

$$q_p(r_0) = \frac{\Gamma}{r_0} \tag{9a}$$

$$q_p(r_l) = \frac{\Gamma}{r_l} \tag{9b}$$

Once the q_p profile (and consequently, from Equation (5), the q_e profile, too) in the surrounding circular layer has been obtained, then the solution of the following first-order differential equation:

$$\frac{dT}{dr} = \left(\frac{\ell_p^2}{\lambda_p}\right) \frac{d^2 q_p}{dr^2} + \frac{1}{r} \left(\frac{\ell_p^2}{\lambda_p}\right) \frac{dq_p}{dr} - \frac{q_p}{\lambda_p} \tag{10}$$

allows us to obtain the radial temperature behavior in the neighborhood of the hot device. In this case, the boundary condition:

$$T(r_0) = T_0 \tag{11}$$

T_0 being the temperature of the hot device, is sufficient to obtain the radial temperature profile.

2.2. Results for a Bi₂Te₃ Surrounding Layer

In order to investigate the influence of nonlocal effects on the radial heat transfer in a practical situation, here, we assume that the circular surrounding layer in Figure 1 is made of a p -doped sample of Bi₂Te₃ at the average temperature of 300 K. The values of the different material functions involved in our problem are summarized in Table 1 [20,21]. In that table, we also quote the radius of the internal device (*i.e.*, r_0) and that of the circular layer (*i.e.*, r_l).

Table 1. Values of the different quantities involved in the problem of the radial heat transfer described by the differential Equations (8) and (10).

λ_e (W m ⁻¹ K ⁻¹)	λ_p (W m ⁻¹ K ⁻¹)	ℓ_e (m)	ℓ_p (m)	r_0 (m)	r_l (m)
2.4	1.6	9.1×10^{-10}	3.0×10^{-9}	10^{-9}	10^{-8}

In principle, the solutions of Equations (8) and (10) are strictly related to the particular experimental situations one is trying to analyze, as for example the amount of Q_0 , as well as the value of T_0 . In order to reduce these dependencies in the computation and to obtain results that are as general as possible, it is convenient to consider the following nondimensional version of Equation (8):

$$\left(\alpha \text{Kn}_p^2 + \text{Kn}_e^2\right) \frac{d^2 q_*^p}{dr_*^2} + \frac{\left(\alpha \text{Kn}_p^2 + \text{Kn}_e^2\right) dq_*^p}{r_*} - (\alpha + 1) q_*^p = \frac{\beta}{r_*} \left(\frac{\text{Kn}_e^2}{r_*^2} - 1\right) \tag{12}$$

wherein the subscript * indicates the (unitary) nondimensional quantities in such a way that $q_p = q_0 q_*^p$, $r = r_0 r_*$, and:

$$(1) \alpha = \frac{\lambda_e}{\lambda_p}$$

$$(2) \text{Kn}_p = \frac{\ell_p}{r_0}$$

$$(3) \text{Kn}_e = \frac{\ell_e}{r_0}$$

$$(4) \beta = \frac{Q_0}{2\pi h r_0 q_0}$$

Moreover, since, for the sake of computation, we assumed $r_l = 10r_0$, the boundary conditions in Equations (9) become:

$$q_*^p(1) = \beta \quad (13a)$$

$$q_*^p(10) = \frac{\beta}{10} \quad (13b)$$

In Figure 2, we plot the theoretical profile of q_*^p arising from the solution of the boundary value problem in Equations (12) and (13), when the surrounding layer is made of a p -doped sample of Bi_2Te_3 at the average temperature of 300 K. In the bottom side of the same figure, we also plot the theoretical behavior of q_*^e , arising from Equation (5), namely,

$$q_*^e = \frac{1}{r_*} - q_*^p$$

We refer to Table 1 for the different parameters involved in the problem, whereas in Equation (12), we have set:

$$\beta = 1 \quad (14)$$

as it naturally arises from Equation (9a).

From the analysis of Figure 2, it follows that in any point of the circular layer q_*^e is always one order of magnitude smaller than q_*^p . This means that for the particular problem we are analyzing, the phonons yield the main contribution to the local heat flux \mathbf{q} . It seems also worth observing that q_*^p attains its maximum in correspondence to the heat source (*i.e.*, when $r_* = 1$) and its minimum when $r_* \in]8; 9[$, whereas q_*^e has the minimum when $r_* \in]1; 2[$ and the maximum when $r_* \approx 6$. Although these features seem to be related to the boundary conditions used, it will be seen in Section 3 that they are also connected to the entropy production in the circular layer. Thus, a deeper investigation of the role of boundary conditions seems to be necessary.

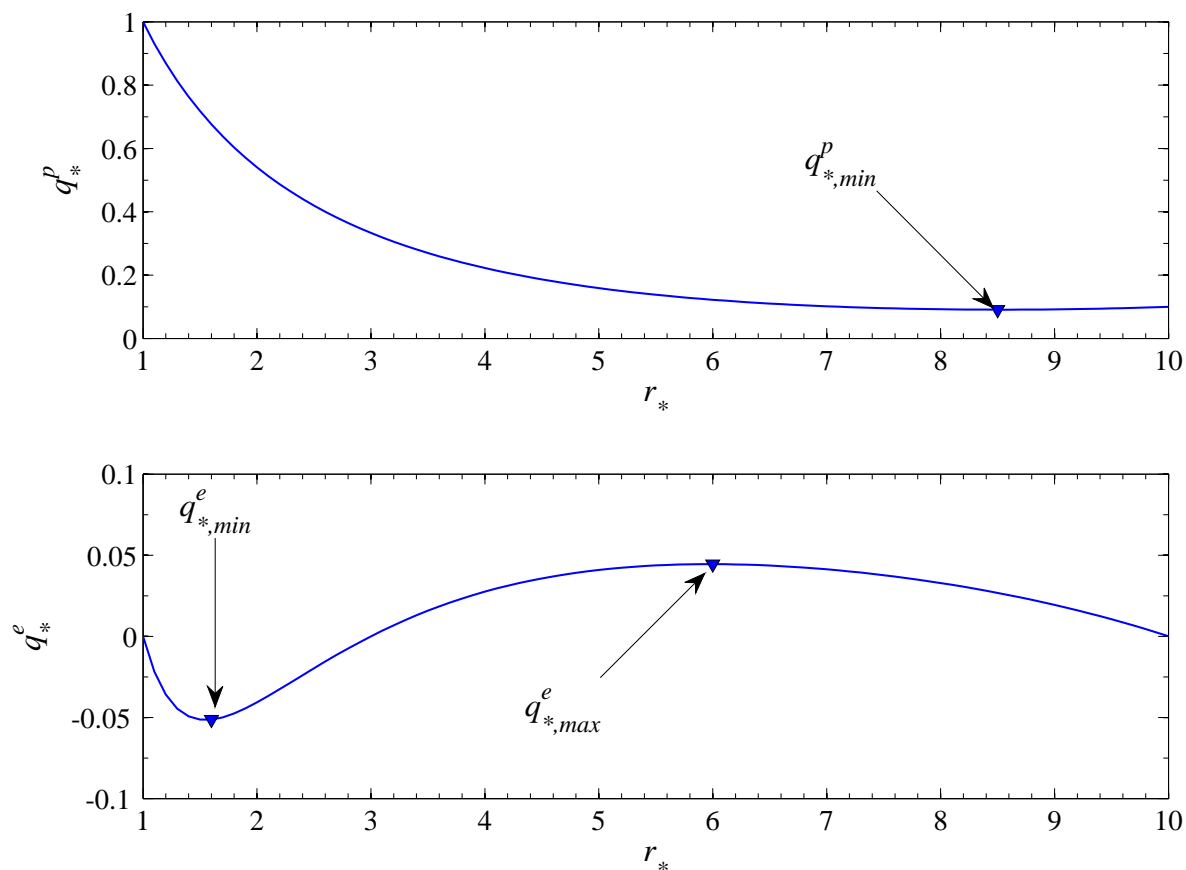


Figure 2. Behaviors of q_*^p and q_*^e versus the radial distance r_* in the circular layer that surrounds the hot device. The profile of q_*^p arises from the solution of the problem in Equation (12) with the boundary condition (13). The q_*^e profile, instead, follows from Equation (5). Refer to Table 1 for the different parameters involved in the problem.

The results obtained for the heat-flux profiles can be used to pursue our goal, namely, to obtain the temperature profile in the neighborhood of the hot device. In particular, we can insert the previous result for q_*^p in the following nondimensional version of Equation (10):

$$\gamma \frac{dT_*}{dr_*} = \text{Kn}_p^2 \frac{d^2q_*^p}{dr_*^2} + \text{Kn}_p^2 \left(\frac{1}{r_*} \right) \frac{dq_*^p}{dr_*} - q_*^p \tag{15}$$

wherein $T_* = T/T_0$ and $\gamma = \lambda_p T_0 / (r_0 q_0)$. Recalling that in the inner hot device, only the phonons may contribute to the heat transfer, it is easy to understand that q_0 should be proportional to T_0 , λ_p and r_0 . These considerations allow us to claim that:

$$\gamma = 1 \tag{16}$$

in Equation (15). Then, its coupling to the nondimensional version of the boundary condition (11), *i.e.*,

$$T_*(1) = 1 \tag{17}$$

produces the result of Figure 3.

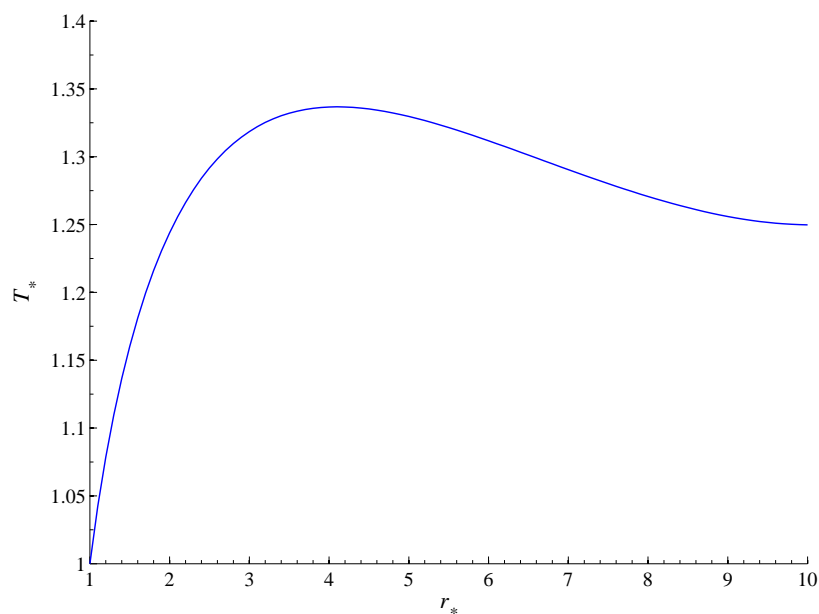


Figure 3. Behavior of T_* versus the radial distance r_* in the circular layer that surrounds the hot device. The theoretical profile for T_* arises from the solution of the problem in Equation (15) with the boundary condition (17). Refer to Table 1 for the different parameters involved in the problem.

As it can be seen from Figure 3, in the neighborhood of the hot device, the temperature profile increases for increasing radial distances. The anomalous hump in the temperature profile that appears in Figure 3 is strictly related to the role played by nonlocal effects, as it is clearly shown by the presence of the coefficients Kn_p in the right-hand side of the heat-transport Equation (2). These effects should be taken into account in practical applications [22]. In fact, if the surrounding layer acts as an insulator among the different components of a same chip, which have to be kept at different temperatures, for example, then the results plotted in Figure 3 show that great attention has to be paid in shortening its size. In particular, Equation (15) points out that the temperature at a radial distance $r_* = 10$ is about 1.25-times higher than that of the hot device at $r_* = 1$, and this value might be unacceptable.

3. Thermodynamic Compatibility

The behavior shown in Figure 3 is anomalous, since the temperature profile in the layer is not everywhere decreasing with the radial distance, as it would be expected from the classical Fourier law [23]. Since in the present problem, the only heat source is the hot device at $r_* \equiv 1$, that behavior (arising from the model Equation (7)) points out that in the circular layer, there are regions wherein the heat is flowing from colder points to hotter ones.

In order to check the physical admissibility of the results plotted in Figure 3, in the present section, we analyze them in view of the second law of thermodynamics. Therefore, let us consider the local balance of the entropy

$$\dot{s} = -\nabla \cdot \mathbf{J}^{(s)} + \sigma^{(s)} \tag{18}$$

with s as the entropy per unit volume,

$$\mathbf{J}^{(s)} = \frac{\mathbf{q}}{T} + \left(\frac{\ell_p^2}{\lambda_p T^2} \right) \nabla \mathbf{q}_p \cdot \mathbf{q}_p + \left(\frac{\ell_e^2}{\lambda_e T^2} \right) \nabla \mathbf{q}_e \cdot \mathbf{q}_e \tag{19}$$

as the entropy flux in the absence of the electric-current density [19] and $\sigma^{(s)}$ as the entropy production per unit volume.

The second law of thermodynamics states that $\sigma^{(s)}$ has to be always non-negative along arbitrary thermodynamic processes [24]. Since in steady states from Equation (18), it follows $\nabla \cdot \mathbf{J}^{(s)} \equiv \sigma^{(s)}$, then the results of Figure 3 are admissible on physical grounds if, and only if, the divergence of the entropy flux is non-negative everywhere in the layer. Indeed, by using the previous results for q_*^p , q_*^e and T_* , the occurrence of that condition may be easily checked once the following nondimensional version of the constitutive relation in Equation (19) is introduced:

$$J_*^{(s)}(r_*) = \frac{\beta}{r_* T_*} + \left(\frac{\text{Kn}_p^2}{\gamma} \right) \frac{dq_*^p}{dr_*} q_*^p + \left(\frac{\text{Kn}_e^2}{\gamma} \right) \frac{dq_*^e}{dr_*} q_*^e \tag{20}$$

with $J^{(s)} = J_*^{(s)} q_0 / T_0$.

Figure 4 plots the behavior of the nondimensional entropy production (*i.e.*, the divergence of $J_*^{(s)}$) as a function of the radial distance from the hot device.

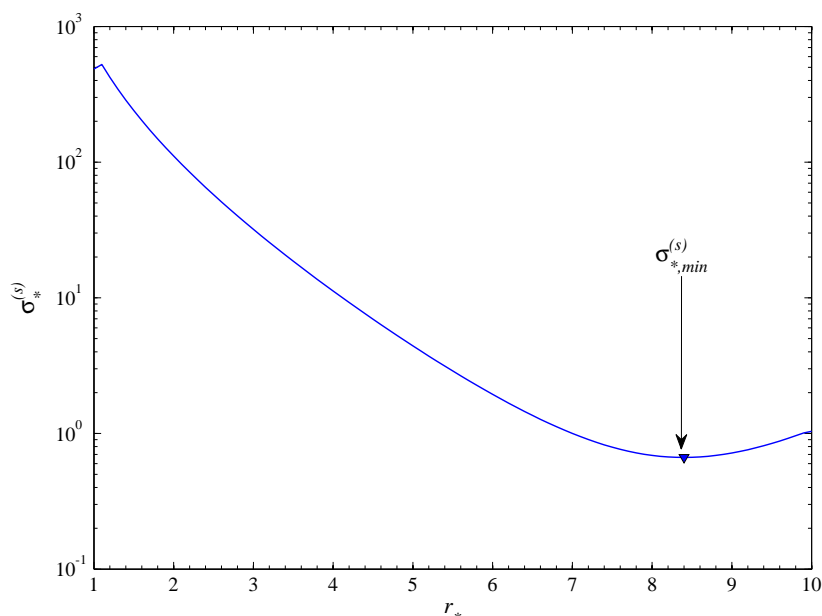


Figure 4. Behavior of $\sigma_*^{(s)}$ versus the radial distance r_* in the circular layer that surrounds the hot device: theoretical results arising from the spatial derivatives of Equation (20). The y-axis in the figure is in a logarithmic scale.

As it can be seen from Figure 4, in the surrounding layer, $\sigma_*^{(s)}$ is everywhere positive. This proves that the hump in the temperature profile observed in Figure 3 is physically admissible, since it agrees with the second law of thermodynamics.

In Figure 4, it is also possible to see that $\sigma_*^{(s)}$ qualitatively behaves as q_*^p for increasing r_* , attaining its maximum and minimum at the same radial distances at which q_*^p also reaches its maximum and minimum values, respectively. Moreover, $\sigma_*^{(s)}$ is increasing for $r_* > 8.4$. We feel that this result is due to the principal role played by the phonon contribution to the local heat flux in the particular problem analyzed here, since it is logical to suppose that $\mathbf{J}^{(s)}$ is mainly influenced by \mathbf{q}^p . However, the behaviors observed in Figures 2 and 4 deserve deeper investigations, especially with respect to their possible connection to the boundary conditions.

4. Conclusions

In the present paper, we have applied the nonlocal model Equations (2) to describe heat transport in quasicrystalline materials [25], namely, in materials that should be properly located at the border line between metals and semiconductors [26]. For these materials, the heat transport is due to the combined flows of phonons and electrons [27], which may also have different temperatures [18,19]. Assuming that both the phonon temperature and the electron one coincide, in the case of steady-state radial heat transfer from an inner hot device surrounded by an external circular layer, we have analyzed the consequences on the temperature profile of the contribution of nonlocal terms entering the temperature equation. Due to the special role played by nonlocal effects at nanoscale, we observed that the temperature profile shows an anomalous hump in the neighborhood of the hot device. The presence of that hump (see Figure 3) suggests that there are some areas wherein the heat flows from colder points to the hotter ones. This result completely disregards that arising from the classical Fourier law [23], which predicts that heat can only flow from the hotter zones to the colder ones. However, to check whether the theoretical predictions arising from Equation (2) are physically admissible, or not, we have analyzed the temperature hump in view of the second law of thermodynamics. Figure 4 shows that the entropy production is everywhere positive and points out that the temperature hump is only an apparent violation of the basic principles of thermodynamics.

There are some further problems related to the present investigation, which deserve consideration and that will be analyzed in our future research.

As we noticed in the Section 1, it is also possible to consider more complex situations, with each of the heat carriers (phonons and electrons) endowed with its own temperature. Accounting for two different temperatures may be important, for instance, when a laser pulse hits the surface of a system. In such a case, initially, the electrons capture the main amount of the incoming energy with respect to the phonons. Subsequently, through electron-phonon collisions, they give a part of it to the phonons. As a further example, we mention those situations in which the electron mean-free path corresponding to electron-phonon collisions is long, so that one may have the so-called “hot electrons”, namely, a population of electrons whose average kinetic energy (*i.e.*, the kinetic temperature) is considerably higher than that of the phonons.

In the presence of two temperatures, the evolution of the partial heat fluxes \mathbf{q}_p and \mathbf{q}_e is governed by the following equations:

$$\tau_p \dot{\mathbf{q}}_p + \mathbf{q}_p = -\lambda_p \nabla T_p + \ell_p^2 (\nabla^2 \mathbf{q}_p + 2\nabla \nabla \cdot \mathbf{q}_p) \quad (21a)$$

$$\tau_e \dot{\mathbf{q}}_e + \mathbf{q}_e = -\lambda_e \nabla T_e + \ell_e^2 (\nabla^2 \mathbf{q}_e + 2\nabla \nabla \cdot \mathbf{q}_e) \quad (21b)$$

wherein T_p and T_e , respectively, mean the phonon and the electron contribution to the average temperature T of the system [28]. Both T_p and T_e may be related to the internal energy per unit volume of phonons u_p and to the internal energy per unit volume of electrons u_e , respectively, by means of the constitutive equations [28]:

$$c_v^p T_p = u_p \quad (22a)$$

$$c_v^e T_e = u_e \quad (22b)$$

with c_v^p being the specific heat at constant volume due to the phonons, and c_v^e being the specific heat at constant volume due to the electrons, whereas the specific heat at constant volume of the whole system c_v is given by $c_v = c_v^p + c_v^e$.

From the practical point of view, Equations (21) and (22) may lead to some perplexities, since one may naturally wonder whether T_p and T_e are measurable quantities, or not [29,30]. Here we notice that in [28] two possible strategies to measure those temperatures have been proposed.

Once Equations (22) have been assumed and both the specific heats are supposed to be constant, then the following partial energy balances can be postulated:

$$c_v^p \dot{T}_p + \nabla \cdot \mathbf{q}_p = 0 \quad (23a)$$

$$c_v^e \dot{T}_e + \nabla \cdot \mathbf{q}_e = 0 \quad (23b)$$

For temperature-dependent specific heats, Equations (23) become nonlinear. However, such a nonlinearity, which is due to the coefficients of the time derivatives of the partial temperatures, does not influence the temperature profile in steady states. Recalling that whenever both electrons and phonons contribute to the heat flow, the internal energy per unit volume u of the whole system is given by $u = u_p + u_e$, then the use of Equation (22) allows us to relate T_p and T_e to the average temperature T of the system as:

$$T = \frac{c_v^p T_p + c_v^e T_e}{c_v} \quad (24)$$

once the thermodynamic relation $u = c_v T$ is used. Note further that the summation of Equation (23) produces the local energy balance (3).

The combination of Equations (21) with Equations (23) allows us, in principle, to derive the temperatures behaviors analytically, which may be interesting in practical applications. However, the solution of that problem is much more complex with respect to the case considered here. In [19], the authors analyzed the consequences of accounting for Equations (21) and (23) in steady-state radial heat transfer from a point source when the partial contributions to the local heat flux are such that:

$$\mathbf{q}_p = \left(\frac{\lambda_p}{\lambda_p + \lambda_e} \right) \mathbf{q} \quad (25a)$$

$$\mathbf{q}_e = \left(\frac{\lambda_e}{\lambda_p + \lambda_e} \right) \mathbf{q} \quad (25b)$$

In this special case, both T_p and T_e display anomalous behaviors, which, however, are in accordance with the second law of thermodynamics [19].

Relaxing the constitutive assumptions (25), it would be interesting to obtain a numerical solution for both the temperature profiles $T_p(r)$ and $T_e(r)$, in order to infer the different role played by phonons and electrons in quasicrystalline materials.

Another interesting problem is related to the efficiency of the thermoelectric coupling to two temperatures. In the presence of thermoelectric effects, the system of Equations (21)–(23) needs to be completed by suitable balance equations for the electric charge and for the electric current (see [18,19] for more details). Such a problem, when the nonlocal effects may be neglected, has been analyzed in [28] with a uniform chemical potential due to the electric charge density and in [31] with a nonuniform chemical potential due to the electric charge density. From those references, it follows that for practical applications, in order to enhance the efficiency of the thermoelectric energy conversion, it would be useful to have:

- (1) the charge distribution as homogeneous as possible;
- (2) the ratio T_e/T as high as possible.

In our future research, we aim to extend the previous investigation to the case in which also nonlocal effects are taken into account, in order to check how these effects may influence the efficiency of thermoelectric devices.

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Author Contributions

In the preparation of this article, all three authors were fully involved in: substantial conception and design of the paper; drafting the article and revising it critically for important intellectual content; final approval of the version to be published. All authors have read and approved the final manuscript.

Conflicts of Interest

The authors declare no conflict of interest.

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