



Research article

Evaluation of a radial basis function node refinement algorithm applied to bioheat transfer modeling

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Abstract: Many bioheat transfer problems involve linear/non-linear equations with non-linear or time-dependent boundary conditions. For heat transfer problems, the presence of time and space-dependent functions under Neumann and Mixed type boundary conditions characterize trivial applications in bioengineering, such as thermotherapies, laser surgeries, and burn studies. This greatly increases the complexity of the numerical solution in several problems, requiring fast and accurate numerical solutions. This paper has a main objective evaluate an adaptive mesh refinement radial basis function method strategy for the classical Pennes's bioheat transfer modeling. Our numerical results had errors of $\sim 0.1\%$ compared to analytical solutions. Thus, the proposed methodology is accurate and has a low computational cost. For step function heating, two RBF shape parameters were applied, again achieving excellent results. The distributions of the nodes in the solution domain show that the primary source of error in the numerical solutions came from the boundary conditions. This finding should arouse the interest of engineers and scientists in the development of new strategies for problems involving boundary conditions with periodic functions.

Keywords: Radial Basis Function; Spatial Adaptive Algorithms; bioheat transfer; thermotherapies; Meshless Methods

Nomenclature

Symbol	Description	Unit
θ_c	Coarse node parameter	-
θ_r	Refine node parameter	-
ω_b	Blood perfusion	$\text{m}^3/\text{s}/\text{m}^3$
ϕ	Radial basis function	-
Δt	Numerical time step	s
c	Specific heat of tissue	$\text{J}/(\text{kg K})$
c_b	Specific heat of blood	$\text{J}/(\text{kg K})$
h_0	Heat convection coefficient	$\text{W}/(\text{m}^2 \text{ K})$
k	Thermal conductivity of tissue	$\text{W}/(\text{m K})$
L	Distance between skin surface and body core	m
Q_m	Metabolic rate of tissue	W/m^3
Q_r	Spatial heating	W/m^3
r	Euclidean distance	-
t	Time	s
T	Tissue temperature	$^{\circ}\text{C}$
T_0	Steady state temperature	$^{\circ}\text{C}$
T_a	Arterial temperature	$^{\circ}\text{C}$
T_c	Blood temperature	$^{\circ}\text{C}$
T_f	Fluid temperature	$^{\circ}\text{C}$
x	Spatial coordinate	m
α	Thermal diffusivity	m^2/s
ε	Multiquadric shape parameter	-
λ	Radial basis function interpolator	-
ρ	Density of tissue	Kg/m^3
ρ_b	Density of blood	Kg/m^3

1. Introduction

Penne [1] presented the classical model for governing heat transfer in tissues. The formulation is simplistic and is composed of an energy conservation balance with a source term. Penne's equation is the most used among the various models proposed to study heat transfer in biological tissues [2,3]. This is because of its mathematical simplicity and its ability to predict the temperature field reasonably well in various applications [4].

The simplicity of bioheat transfer modeling is associated with the absence of convective terms in the mathematical formulation. This feature is usually positive, making computational approaches more stable. However, bioheat transference problems can present high numerical complexity due to the presence of source terms, non-trivial boundary conditions [5], and complex geometries [6].

Analytical solutions to specific bioheat transfer problems are available in the literature [7–9]. However, [10] noted that there is not yet a general analytical form for the solution of Penne's equation. In this way, accurate numerical solutions remain essential for the resolution of the model in question.

The Radial Basis Function (RBF) method was applied by [11] to solve partial differential equations and thus garnered the attention of engineers and scientists because of its simplicity, even with complex geometries and non-linear problems of different dimensions [12]. However, the application of the RBF method for heat bioheat transfer problems is still in development [13]. It is currently restricted to applications and evaluations of the RBF method to classical problems (burns and thermotherapy in different dimensions and multi-domains) [14,15].

One of the main characteristics of the RBF method is that it does not make use of a mesh, which facilitates the construction of complex geometries in complex dimensions and enables the development of spatial adaptive algorithms. The literature on the RBF method presents a series of spatial adaptive algorithms for classical convection-diffusion equations [16,17]. The mentioned authors present different algorithms for choosing the best node location. They presented enthusiastic results involving convective problems and functions with high gradient regions. However, they have limitations concerning the domain's dimension. On the other hand, [18] presented the multi-domain Residual Subsampling Method (RSM), an adaptive node refinement strategy that enables the user to set parameters allowing the addition and removal of nodes based on residuals evaluated at a finer point set. This fact might be of interest to bioheat numerical problems since RSM can be useful to refine the boundary domain's (regions with high gradients in bioheat problems) and coarse-smooth gradient regions (usually on the internal nodes in bioheat problems), reducing computational time, and lowering the integration errors. However, the mentioned algorithms were not evaluated for the classical bioheat applications, and their stability and convergence in transient problems are still being researched. To our knowledge, for conventional bioheat modeling, no work involving the RBF technique has presented a spatial adaptive numerical solution. The presence of non-linear source terms and Neuman and Mixed boundary conditions suggests the implementation of such algorithms aiming at solutions of high accuracy and high computational performance as performed in applications concerning mesh-based methods [19], neural network modeling [20] and boundary value methods [21,22].

Recently [23] proposed a bioheat transfer study implementing the Radial Basis Function method. In this study, constant and sinusoidal boundary conditions were evaluated in three types of heat transfer models, i.e., Penne's bioheat model, single-phase lag model. The authors contributed with a physical analysis restricted to sinusoidal and constant heating boundary conditions and the physical parameters. However, the quality of numerical RBF solutions was not discussed or compared with benchmark solutions. The bioheat model and its respective boundary conditions and physical parameters directly impact the propagation of the tissue's thermal signals. Additionally, the precision of the numerical methodology may be a source of solution errors and numerical instability. This investigation topic becomes essential when the modeler proposes adopting numerical methods with empirical parameters and mesh refinement strategies. Moreover, the numerical method should overcome the complexity inherent to transient models, source terms, and step/sinusoidal boundary conditions. The importance of a performance analysis concerning RMS as a methodology for solving transient problems is emphasized by [18].

The present work aims to evaluate a spatial adaptive methodology coupled with the RBF-meshless method to obtain numerical solutions of the classic bioheat model.

The analytical solutions developed by [8] by Green's function method will be applied to

evaluate the numerical solution in three scenarios: constant skin surface heating, heating the skin surface using a step function, and heating the skin surface using a periodic function.

2. Materials and method

2.1. The bioheat transfer model

Most theoretical analyses on heat transfer in living tissues are based on Penne's equation, which describes the influence of blood flow on the temperature distribution in tissue.

In this work, in line with [8], only one-dimensional models with constant thermal parameters will be used, which results in a good approximation when heat propagates in the direction perpendicular to the skin surface.

Penne's bioheat transfer general equation for one-dimension analysis can be described as;

$$\rho c \frac{\partial T}{\partial t} = k \frac{\partial^2 T}{\partial x^2} + \omega_b \rho_b c_b (T_a - T) + Q_m + Q_r(x, t) \quad (1)$$

where ρ, c and k are the density, specific heat, and thermal conductivity of the tissue, respectively; ρ_b and c_b denote the blood's specific density and heat; ω_b blood perfusion; T_a the arterial temperature that is treated as a constant; T the temperature of the tissue; Q_m the heat generated by metabolism; and $Q_r(x, t)$ is the source of heat due to external applied heat.

According to [8], the initial temperature profile for the rest state of biological bodies can be obtained by solving the following equation and its boundary conditions.

$$k \frac{d^2 T_0(x)}{dx^2} + \omega_b \rho_b c_b [T_a - T_0(x)] + Q_m = 0 \quad (2)$$

$$T_0(x) = T_c, x = L \quad (3)$$

$$-k \frac{dT_0(x)}{dx} = h_0 [T_f - T_0(x)], x = 0 \quad (4)$$

where, $T(x, 0) = T_0(x)$ is the steady-state temperature profile before heating, T_c is the blood temperature and often considered as constant, h_0 is the apparent heat convection coefficient between the surface of the skin and the surrounding air physiologically in the basal state, and is a general contribution of natural convection and radiation, and T_f is the air temperature. The surface's skin is set on $x = 0$ while the center of the body in $x = L$.

During a practical thermal process, the boundary condition presented by Eq 4 is always dependent on time, which can be generalized to:

$$-k \frac{\partial T}{\partial x} = f_1(t), x = 0 \quad (5)$$

Or

$$-k \frac{\partial T}{\partial x} = h_f [f_2(t) - T(x)](t), x = 0 \quad (6)$$

where, $f_1(t)$ represents the surface heat flux dependent on time, and $f_2(t)$ is the medium cooling temperature on the time and the coefficient of heat convection between the medium and the surface of the skin. In this work, the body temperature was considered constant (T_c) considering that the biological body tends to maintain a stable central temperature, i.e.,

$$T(x) = T_c, x = L \quad (7)$$

2.2. The Radial Basis Function Method

A radial base function $\phi(r)$ is a radially symmetrical n -dimensional function: $R^n \rightarrow R$ which depends only on the Euclidean distance ($r = \|x - x_j\|$) between a center x and a point evaluated in x_j . In this work, multiquadric radial base function was applied to interpolate an $f(x)$ function into R^n assuming that

$$f(x) = \sum_{j=1}^N \lambda_j \phi(r), x \in R^n \quad (8)$$

where $\phi_j(r) = \sqrt{\varepsilon^2 + (x - x_j)^2}$ and ε is a non-zero empirical parameter.

The coefficients λ are calculated by solving the linear system presented in the form $\lambda = \Phi^{-1}f(x)$.

2.3. Discretization of bioheat transfer equation by Radial Basis Function Method

Beginning from Eq 8, the partial derivatives can be easily obtained from the following Eqs 9 and 10:

$$\frac{\partial f(x)}{\partial x} = \sum_{j=1}^N \lambda_j \frac{\partial \phi}{\partial x} = \sum_{j=1}^N \lambda_j \frac{x - x_j}{\phi}, x \in R^n \quad (9)$$

And

$$\frac{\partial^2 f(x)}{\partial x^2} = \sum_{j=1}^N \lambda_j \frac{\partial^2 \phi}{\partial x^2} = \sum_{j=1}^N \lambda_j \frac{\varepsilon^2}{[\varepsilon^2 + (x - x_j)^2]^{3/2}}, x \in R^n \quad (10)$$

Consider Eq 1 as a function temperature $T(x, t)$. The Crank-Nicolson [24] discretization scheme consists of a finite difference approach to approximate partial derivatives on $T(x, t)$ by the following equation.

$$\frac{T^{t+\Delta t} - T^t}{\Delta t} \approx \frac{1}{2} [T(x, t) + T(x, t + \Delta t)] \quad (11)$$

Considering the terms Q_m and Q_r as constant in the Eq 1 and expanding the Eq 11 will result in:

$$T^{t+\Delta t} - \frac{\Delta t}{2} \left[\alpha \frac{\partial^2 T^{t+\Delta t}}{\partial x^2} - w c_b T^{t+\Delta t} \right] + \Pi = T^t + \frac{\Delta t}{2} \left[\alpha \frac{\partial^2 T^t}{\partial x^2} - w c_b T^t \right] + \Pi \quad (12)$$

where $\alpha = k/\rho c$ and Π represents the sum of the constant terms contained in the model $\Pi = w c_b T_a + q_m + q_r$. From Eq 12, two new operators will be defined, H^+ and H^- , considering $T(x, t) \approx \phi(x, t)$ and $\frac{\partial^2 T}{\partial x^2} \approx \frac{\partial^2 \phi}{\partial x^2}$ as follows:

$$H^+ = \phi^{t+\Delta t} - \frac{\Delta t}{2} \left[\alpha \frac{\partial^2 \phi^{t+\Delta t}}{\partial x^2} - w c_b \phi^{t+\Delta t} \right] + \Pi \quad (13)$$

$$H^- = \phi^t + \frac{\Delta t}{2} \left[\alpha \frac{\partial^2 \phi^t}{\partial x^2} - w c_b \phi^t \right] + \Pi \quad (14)$$

Finally, the operators H^+ and H^- will be used to approximate the function $T(x, t)$ using the radial base function, shown below:

$$\sum_{j=1}^N \lambda_j^{t+\Delta t} H^+ = \sum_{j=1}^N \lambda_j^t H^- \quad (15)$$

Equation 15 generates a system of linear equations, which can be solved by LU decomposition

to obtain the unknowns, $\lambda^{t+\Delta t}$ from the known values of λ^t , at a previous time step. Then they give rise to $T(x, t)$ by means of Eq 8.

2.4. Meshfree node adaptive algorithms – Residual Subsampling Method

The residual subsampling method developed by [18] is a simple and easy method to add, reallocate, and remove the nodes based on an interpolation process.

This method consists of generating an initial discretization using equally spaced N points and finding the Radial Basis Function approximation. The Figure 1, illustrate this first RSM step for interpolate a step function.

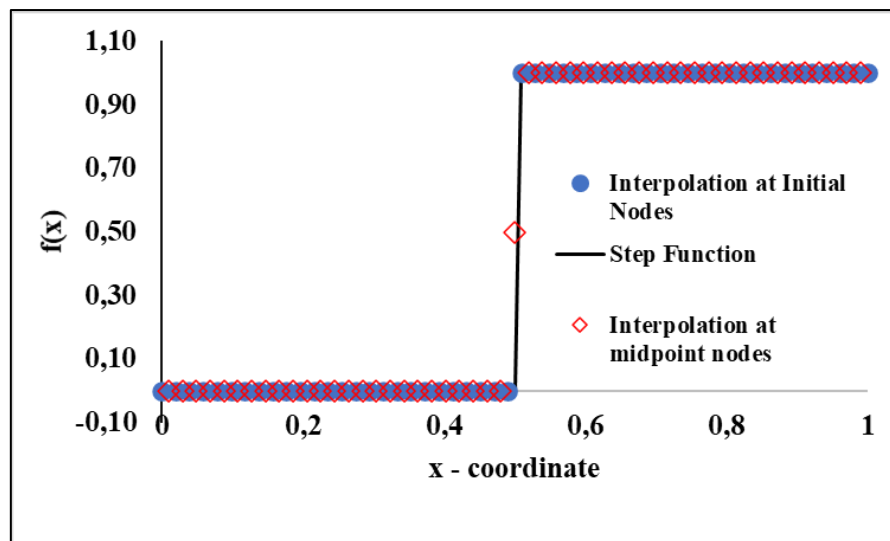


Figure 1. Interpolation of a step function on initial and midpoint nodes.

Next, the interpolation error of the midpoints between the nodes is calculated as illustrated in Figure 2.

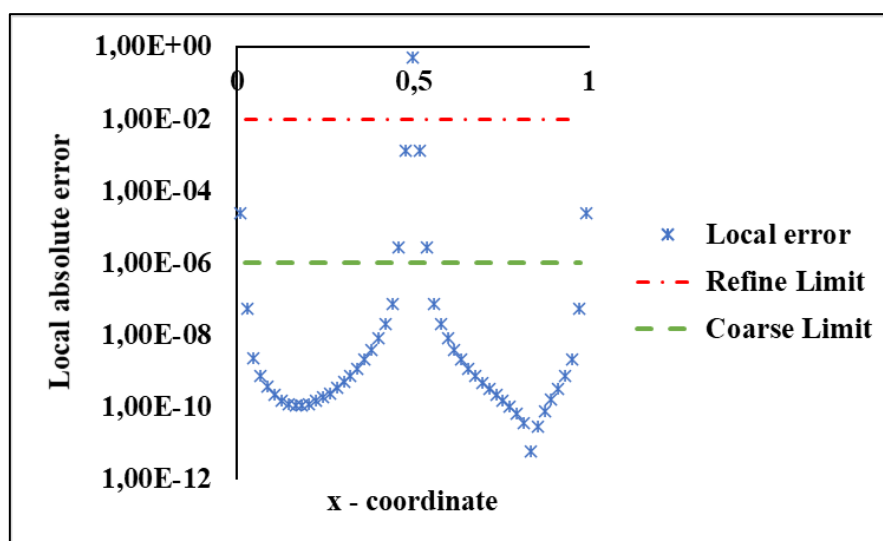


Figure 2. Local interpolation errors and limits to remove and add nodes.

Points at which the error exceeds a threshold become (refine limit) θ_r centers, and the centers that are between two points whose error is below a lower threshold (coarse limit) θ_c are removed. Contour nodes are always left intact. Figure 3 shows the final set node after an iteration of RSM for interpolating a step function.

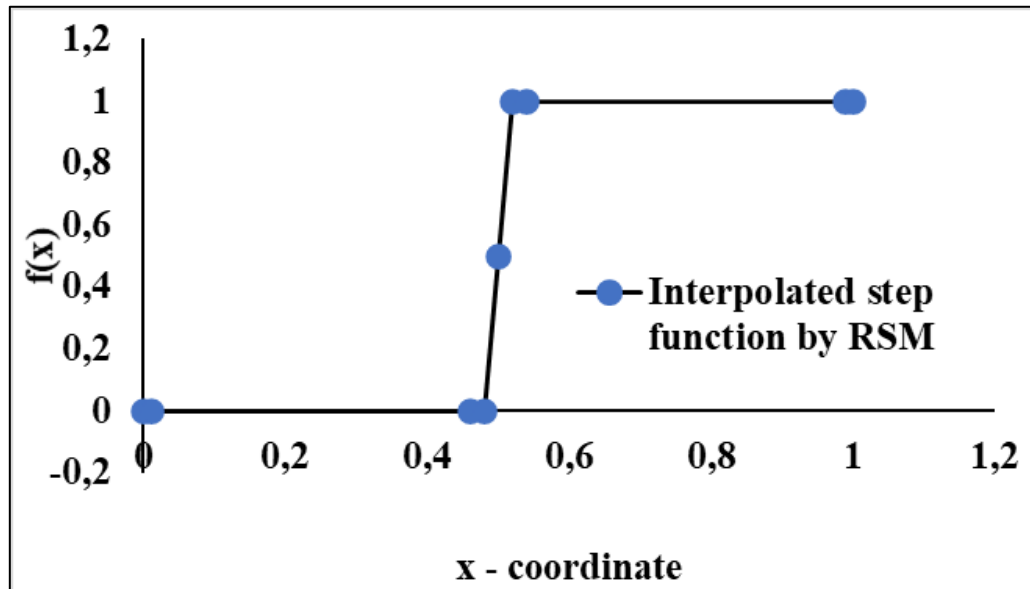


Figure 3. Interpolation result after only one iteration of RSM.

The shape parameter of each center is chosen based on spacing with the nearest neighbors, and the Radial Basis Function approximation is recalculated using the new center set. The original approach by the Residual Subsampling Method is made by Multiquadrics RBF. However, it should be easily adapted to other global support RBFs. The coarse and refinement process should be repeated until an error criterion be reached.

In short, the adaptation process follows the family paradigm of resolving estimating/approaching until a stopping criterion is met [18].

The Residual Subsampling Method can be easily modified to solve linear or non-linear problems of contour value. Figure 4 shows the flowchart of this method applied to the present study.

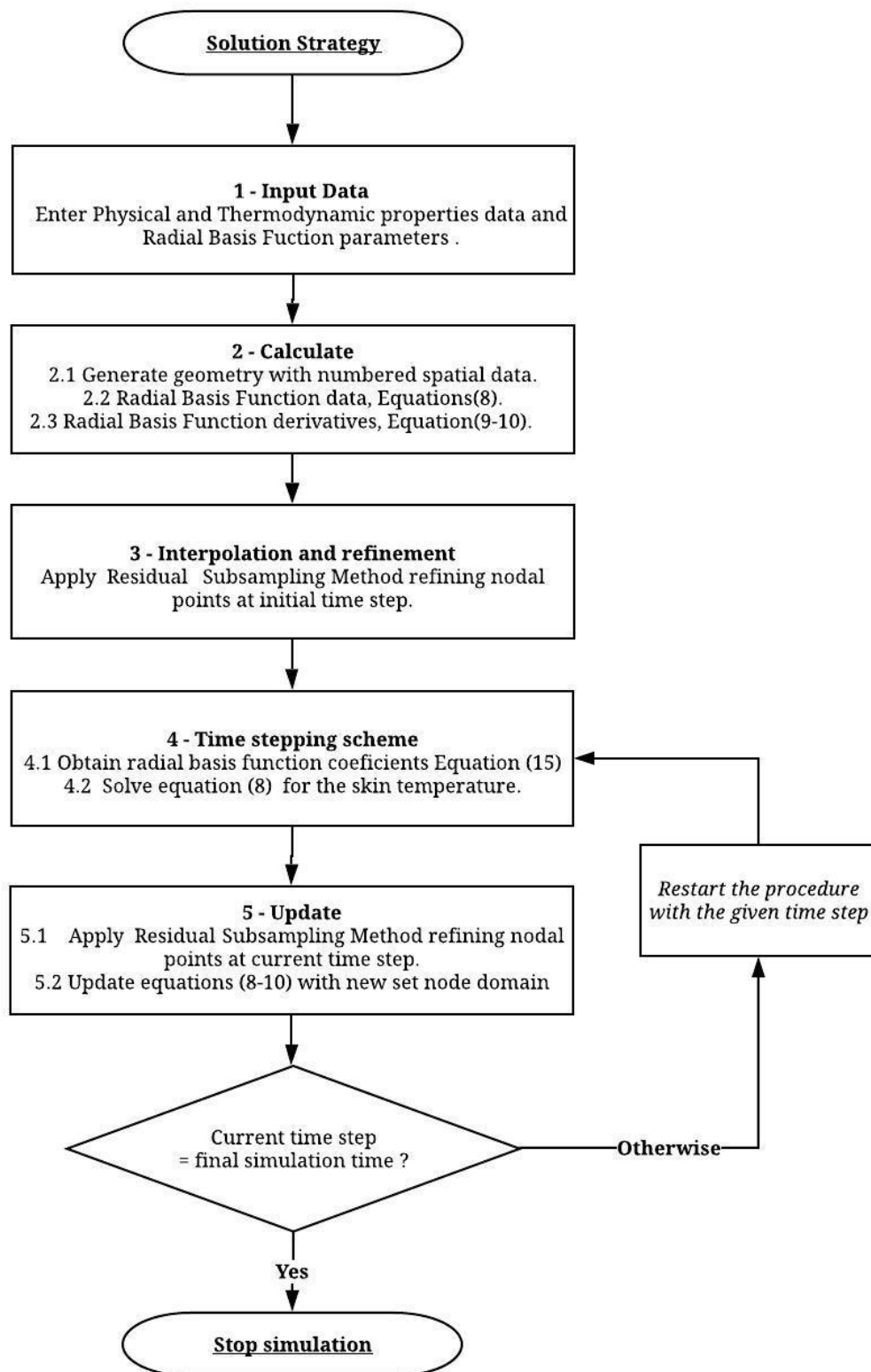


Figure 4. Flowchart of Radial Basis Function with mesh refinement strategy for bioheat transfer problems.

3. Results and discussion

In the following calculations, the typical skin physical properties will be applied as proposed by [8]: $\rho = \rho_b = 1000 \text{ kg/m}^3$, $c = c_b = 4200 \text{ J/(kg K)}$, $T_a = T_c = 37^\circ\text{C}$, $k = 0.5 \text{ W/(m K)}$, $\omega_b = 0.0005 \text{ m}^3/\text{s/m}^3$, $Q_m = 33800 \text{ W/m}^3$. The convection coefficient of apparent heat due to natural convection and radiation is taken as $h_0 = 10 \text{ W/(m}^2 \text{ K)}$. The forced convection coefficient is applied as $h_f = 100 \text{ W/(m}^2 \text{ K)}$, while the temperature of the surrounding fluid was chosen as $T_f = 25^\circ\text{C}$. Also, as demonstrated in many works [6,25,26], the indoor temperature of the tissue usually tends to be constant at a short distance, such as 2 to 3 cm, so it will be used in $L = 3 \text{ cm}$ in this study.

The local (E^∞) and global (E^2) errors will be calculated using the following equations:

$$E^\infty = \max_{0 \leq i \leq N} |f(x_i) - s(x_i)| \quad (16)$$

$$E^2 = \sqrt{\frac{1}{N} \sum_{0 \leq i \leq N} |f(x_i) - s(x_i)|^2} \quad (17)$$

where $f(x)$ is the analytical solution, and $s(x)$ is the numerical solution.

3.1. Skin surface under constant heating

This method of heating is often used in hot plate tests. This test was introduced in pharmacology by [27] to test the efficiency of new analgesics in use at the time and has since been widely used by researchers and scientists in the field [28,29]. The hot plate test evaluates the reaction time to thermal stimuli of mice placed on hot plates at 55°C [27,29].

The numerical solution for surface temperature response is shown in Figure 5, while the numerical solution of the temperature response along the tissue compared to its analytical solution is shown in Figure 6. The curves presented are the transient temperatures of the tissues subject to three surface heating constants: $P_0(t) = 1000 \text{ W/m}^2$, $P_0(t) = 500 \text{ W/m}^2$, and $P_0(t) = 200 \text{ W/m}^2$.

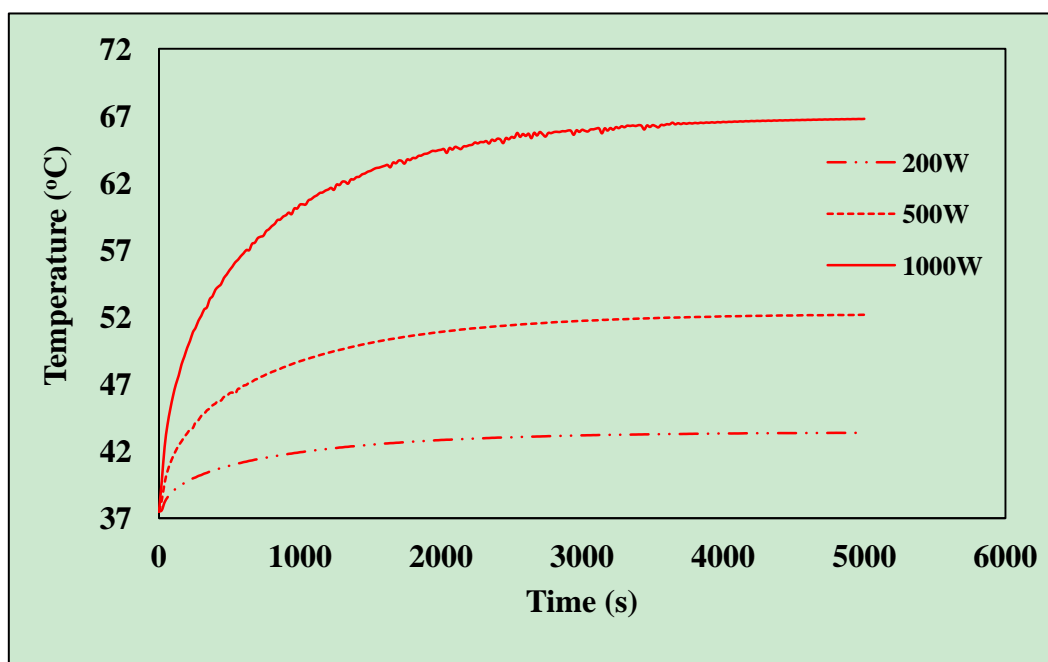


Figure 5. Skin surface temperatures for different surface heating constants.

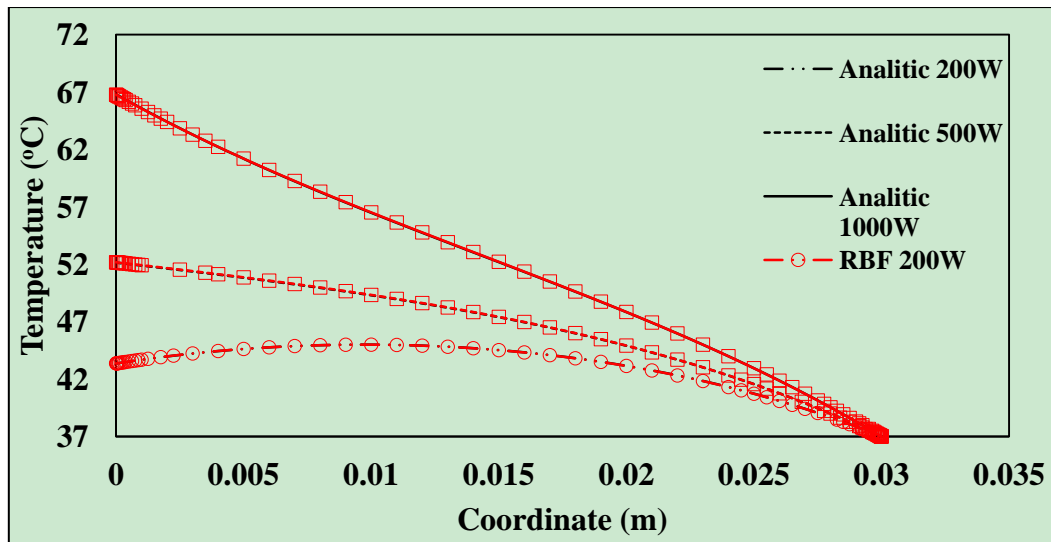


Figure 6. Comparison against numerical and analytical solutions for skin temperature profiles at simulation time = 5000 seconds.

Table 1 shows the values of the shape parameters used for model calibration and the values of errors E^2 and E^∞ resulting from the comparison of numerical results with the analytical solution for the constant heating condition of the surface.

Table 1. Numerical parameters and respective relative and absolute errors for adaptive strategy.

P_0 (W/m ²)	Δt	N	ε	θ_r	θ_c	E^2 (%)	E^∞ (%)
200	20	67	0.74	10^{-3}	10^{-6}	0.023	0.051
500	20	64	0.82	10^{-3}	10^{-6}	0.039	0.063
1000	20	70	1.09	10^{-3}	10^{-6}	0.031	0.058

Table 2. Numerical parameters and respective relative and absolute errors for an RBF method without node mesh refinement.

P_0 (W/m ²)	Δt	N	ε	E^2 (%)	E^∞ (%)
200	20	301	7668	0.023	0.056
500	20	301	8569	0.026	0.076
1000	20	301	11268	0.053	0.117

Analyzing Figures 5 and 6, and the results presented in Table 1, it was concluded that all error rates were below 0.1%. Also, the final values of nodes were extremely low for the three powers used, indicating an excellent performance of the proposed numerical methodology. It is important to focus on the uniformity of the errors presented by the technique, which can be observed more specifically in Table 1, which presents the maximum errors close to the global errors. Another phenomenon to be highlighted is the greater concentration of nodes on the boundaries, indicating that the contour regions present greater instability, caused by the formulation of the boundary condition or the non-linear terms of the equation. Table 2 corroborates the mentioned results when comparing the RBF adaptive node strategy in Table 1 against the RBF uniform node strategy in Table 2.

3.2. Surface heating by a step function

Practical examples of this case can be found in ophthalmologic surgeries through a single laser pulse [30] or skin burns due to an instant fire, hot plate, liquid, and gas for a short period. In the atomic explosion, the burn caused by the high-temperature shock wave also belongs to this problem [8].

Figure 7 shows the transient temperature response to a heating on the surface of the step function $P_0(t) = 1000 \text{ W/m}^2$ for $t < 1200 \text{ s}$ and $P_0(t) = 0$ for $t \geq 1200 \text{ s}$, while spatial heating $Q_r = 0$ and $\omega_b = 0.0005 \text{ m}^3/\text{s/m}^3$.

The numerical solution of the temperature response to heating by a step function compared to its analytical solution is presented in Figure 8, as well as the distribution of the resulting nodes in solving the problem.

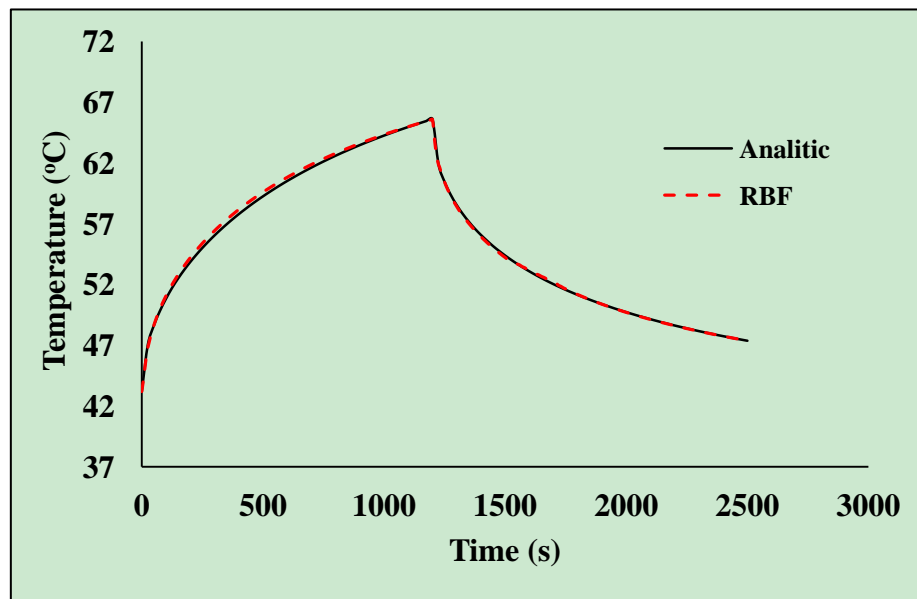


Figure 7. Temporal skin surface temperature profile under step heating boundary condition.

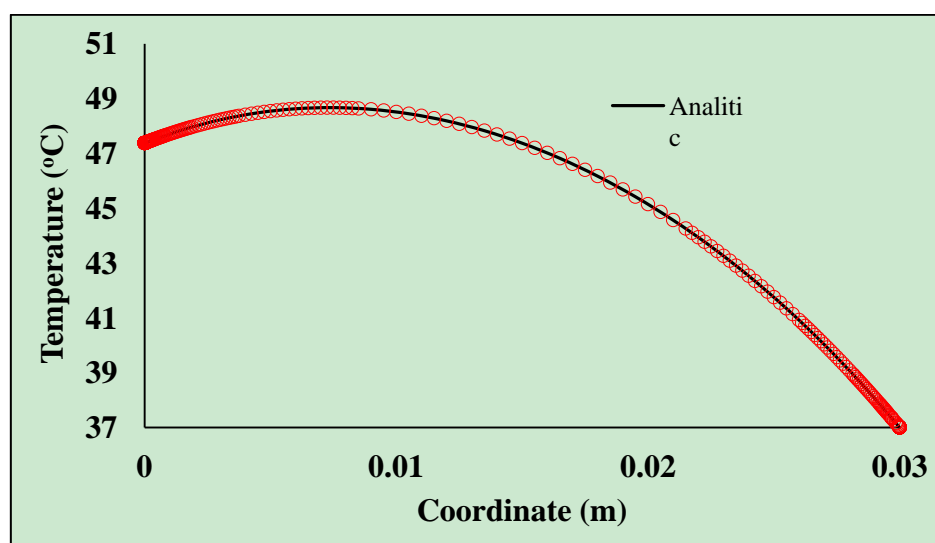


Figure 8. Temperature spatial profile for step heating boundary condition at simulation time = 2500 seconds.

Table 3 shows the respective values of the shape parameters used for RBF mesh refinement strategy and regular node distribution. The values of errors E^2 and E^∞ result from the comparison with the analytical solution for the condition of the heating of the surface by a step function.

Table 3. Shape parameters and errors for the step function scenario.

RBF strategy	P_0 (W/m ²)	Δt	N	ε_1	ε_2	θ_r	θ_c	E^2 (%)	E^∞ (%)
Node refinement	1000	10	288	0.673	0.448	10^{-6}	10^{-10}	0.025	0.049
Without node refinement	1000	10	501	11762	7300	-	-	0.028	0.051

To preserve good performance, the form factor of the multiquadric function had to be changed during the simulation, justified by the behavior of the step function in the contour. Therefore, for heating times below 1200 s, the applied value was $\varepsilon_1 = 0.673$, and for longer times $\varepsilon_2 = 0.448$.

Analyzing Figures 7 and 8, and the results presented in Table 3, it was concluded that the error rates presented values below 0.1%. The final number of nodes was very low, thus supporting the application of the proposed numerical methodology. It was also observed that the regions with the highest concentration of nodes were on the boundaries. The skin surface is a region of greater instability in the execution of the model, due to the sudden change of heat flux on the surface. This fact raises the difficulty of maintaining the desired stability and convergency criteria and makes it necessary to triple the number of nodes compared to the previous scenario.

3.3. Surface heating under a periodic function

Many researchers have investigated the effects of warming by a periodic function of biological tissues [32,25,33]. This type of heating is used in medicine and reflects a situation in which heating is caused by repeated laser irradiation in thermotherapy procedures [8], and is also used to measure blood perfusion in biological tissues [25].

Using the current solution in the finite domain, the temperature for periodic heating – both of the skin surface and inside biological bodies – can be easily obtained. Periodic surface heating can be expressed as [8].

$$P_0(t) = q_0 + q_w \cos(\omega_1 t) \quad (18)$$

Figure 9 shows the response of transient temperature in biological bodies subject to surface heating by a periodic function.

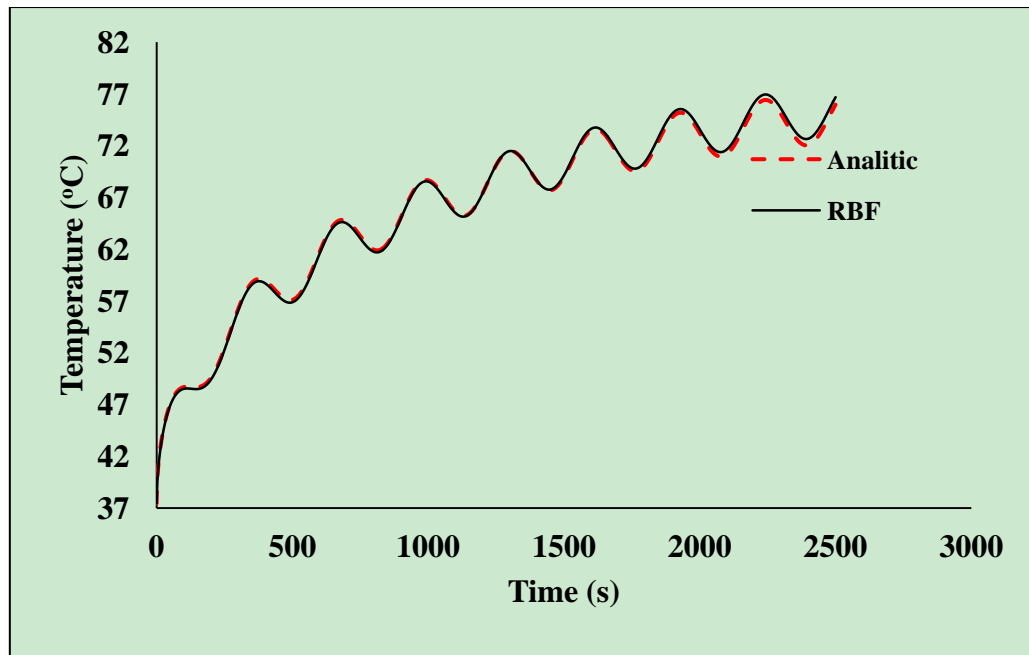


Figure 9. Transient temperature response to surface heating by a periodic function.

The numerical solution of the temperature response to heating by a periodic function compared to its analytical solution is presented in Figure 10, as well as the distribution of the resulting nodes in solving the problem.

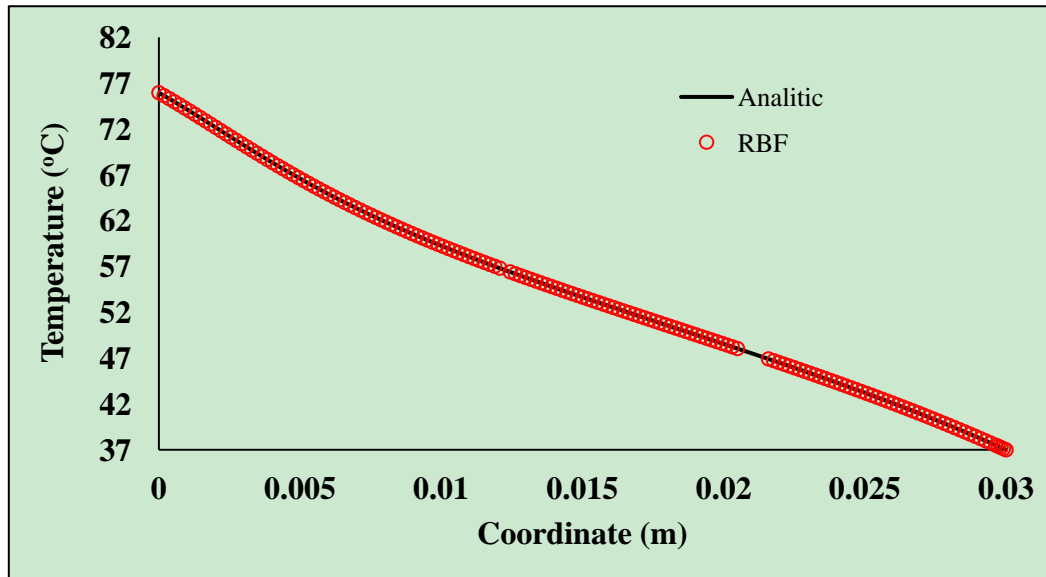


Figure 10. Node location and tissue temperature under a surface heating by a periodic function at simulation time = 2500 seconds.

Table 4 shows the values of the shape parameters used for model calibration and the values of errors E^2 and E^∞ resulting from the comparison with the analytical solution for the condition of the heating surface by a periodic function.

Table 4. Shape parameters and errors for surface heating condition by a periodic function.

RBF strategy	P_0 (W/m ²)	Δt	N	ε	θ_r	θ_c	E^2 (%)	E^∞ (%)
Node refinement	$1000 + 500 \cdot \cos(0.02 \cdot t)$	1.25	162	0.2	10^{-4}	10^{-10}	0.425	0.982
Without node refinement	$1000 + 500 \cdot \cos(0.02 \cdot t)$	1.0	151	5849	-	-	0.680	2.252

Analyzing Figures 9 and 10, and Table 4, it was concluded that the error rates were around 0.1%, which are acceptable considering that only 162 nodes were needed to solve the problem. The points were distributed almost evenly along the tissue, but in the central region, it can be noted that there was a slight absence of points because the tendency is for the points to focus on the contours. It is essential to highlight the error uniformity of adaptive strategy when compared against the regular node distribution. This is a desirable phenomenon for numerical approaches representing the numerical stability and absence of local oscillations.

4. Conclusion

Here we have presented a performance evaluation of the RBF node refinement strategy to solve bioheat transfer problems.

The numerical results had an excellent approximation relative to the analytical solutions, thus confirming that RBFs (coupled with a spatial node refinement strategy) are acceptable methods for solving bioheat transfer problems with low set of nodes. The novel RBF approximation for boundary condition involving step function presented excellent results. Additionally, the numerical solution showed uniformity on local and global error norms for all simulated scenarios.

However, it is essential to highlight that for surface heating by step function and periodic function, the transient terms in these respective boundary conditions directly affected the proposed methodology's performance. Additionally, the adaptive strategy may work excessively to maintain the error precision imposed by θ_r and θ_c parameters inducing instability in integration on time—however, the numerical approach still presenting good results. This fact should be considered in problems with irregular geometries. It should not be trivial to guarantee the presence of nodes only inside the domain coordinates. Additionally, remove a boundary node in 2D or 3D models within complex geometries may generate serious stability problems.

As a proposal for future work, we suggest expanding the adaptive technique proposed in this work for applications in 2- and 3-dimension bioheat transfer models involving multilayer problems, complex geometries, and the presence of convective terms. Practical adaptive radial basis function approach by bio-heat transfer modeling computed together with biomechanics may present complex modeling phenomena as soft tissue movement and thermal expansion/shrinkage effect [34,35] and should be investigated.

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