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Solving time-dependent engineering problems with multiquadrics

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Abstract

This paper deals with the solution of time-dependent problems. The multiquadric radial basis function method is formulated, with a new approach for transient problems. One- and two-dimensional problems are considered. The forward difference and the Crank–Nicolson time-marching schemes for parabolic cases are considered. The central difference integration method of the Newmark family is considered for hyperbolic problems. The method proves its accuracy in four numerical examples. © 2004 Elsevier Ltd. All rights reserved.

1. Introduction

Time-dependent problems are of considerable relevance in engineering and science. This paper deals with the solution of time-dependent problems with radial basis functions. The multiquadric radial basis function method is formulated, with a new approach for transient problems. One-(1D) and two-dimensional (2D) problems are considered. The method proves its accuracy in some numerical 1D and 2D examples.

In this paper, a recent meshless approximation technique is used, based on radial basis functions (RBFs). This truly meshless technique is insensitive to spatial dimension, and considers

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only a cloud of nodes (centers) for the spatial discretization of both the problem domain and the boundary.

Other meshless methods have also been proposed. They may be classified as smooth particle hydrodynamics [1–3], diffuse element method [4], element free Galerkin [5–7], reproducing kernel particle method [8–13] and HP clouds [14]. In recent years, Liu's group gave significant contributions, namely the local point interpolation method (LPIM) and the local radial point interpolation method (LR-PIM) [15,16], the radial point interpolation method (RPIM) [17] and the boundary radial point interpolation method (BRPIM) [18].

The radial basis function method was first used by Hardy [19,20] for the interpolation of geographical scattered data, and later used by Kansa [21,22] for the solution of partial differential equations (PDEs). Many other radial basis functions can be used as reviewed in the recent book of Liu [23], namely Powell [24], Coleman [25], Sharan et al. [26], Wendland [27], among others.

The use of RBFs for 2-D solids has been proposed by Liu et al. [28–30] and by Ferreira [31,32] for composite plates and beams. The method has also been applied to other engineering problems such as in Refs. [33–35].

This paper concentrates on the solutions of 1D and 2D engineering problems, such as heat conduction and beams in bending, using both the forward difference and Crank–Nicolson time-marching schemes with interpolation by the unsymmetrical multiquadric method.

2. The multiquadric method

The multiquadric method relies on the Euclidian distance between nodes and in some cases on a shape parameter (c), user-defined and object of various discussions. The influence of such parameters not only defines the RBF, but may also provide ill-conditioned problems with inadequate solutions.

The numerical solution of PDEs is traditionally dominated by finite element methods, finite volume methods or finite difference methods. All of these methods are based on local interpolation strategies and depend on a mesh for local approximation. In these methods, although the function is continuous across meshes, its partial derivatives are not [36–38].

A new approach for solving partial differential equations is based on RBFs. An RBF depends only on the distance to a center point \mathbf{x}_j and is of the form $g(||\mathbf{x} - \mathbf{x}_j||)$. The RBF may also depend on a shape parameter c, in which case $g(||\mathbf{x} - \mathbf{x}_j||)$ is replaced by $g(||\mathbf{x} - \mathbf{x}_j||, c)$ [21,22,39–42].

Consider a set of nodes $x_1, x_2, \ldots, x_N \in \Omega \subset \mathbb{R}^n$. The radial basis functions centered at \mathbf{x}_j are defined as

$$g_i(\mathbf{x}) \equiv g(\|\mathbf{x} - \mathbf{x}_i\|) \in \mathbb{R}^n, \quad j = 1, \dots, N,$$
(1)

where $\|\mathbf{x} - \mathbf{x}_i\|$ is the Euclidian norm.

Some of the most common RBFs are [21,22,39–41]:

Multiquadrics:
$$g_i(\mathbf{x}) = (\|\mathbf{x} - \mathbf{x}_i\| + c^2)^{1/2},$$
 (2)

Inverse Multiquadrics :
$$g_i(\mathbf{x}) = (\|\mathbf{x} - \mathbf{x}_i\| + c^2)^{-1/2},$$
 (3)

Gaussians :
$$g_i(\mathbf{x}) = e^{-c^2 \|\mathbf{x} - \mathbf{x}_j\|^2}$$
, (4)

Thin Plate Splines :
$$g_j(\mathbf{x}) = \|\mathbf{x} - \mathbf{x}_j\|^2 \log \|\mathbf{x} - \mathbf{x}_j\|,$$
 (5)

where c is a shape (user-defined) parameter. In this paper, only the multiquadric method is used.

RBFs are insensitive to spatial dimension, making the implementation of this method much easier than, e.g., finite elements [21,22].

Time-dependent problems have been treated with multiquadrics since their reappearance due to Kansa in 1990 [21,22]. Kansa applied his nonsymmetric multiquadric method to solve two parabolic and hyperbolic PDEs [22], the linear advection–diffusion equation and the dynamic 1D von Neumann blast wave. To solve the first problem, the author used a standard implicit approximation scheme; as for the second equation, a fourth-order Runge–Kutta scheme is used to perform the integration in time.

Another interesting approach to solve the 1D parabolic problem using RBFs came from Fasshauer [43], who made use of the Newton method and Nash iteration with an explicit time-stepping method to interpolate the classical Heat Equation.

Further incursions on the fluid dynamics domain have been made by authors like Hon and Wong [34]. They used the RBF meshless method to solve a multilayer computational model for simulating 3D tidal flows in coastal waters [44]. In this paper, the authors also prove the applicability of the RBF method in the solution of large-scale systems using a domain decomposition technique. They continued the exploration of hydrodynamic problems using RBF-based techniques, solving the Shallow water equations [34]. A different base of interpolating functions was used, the compactly supported RBFs or CSRBFs, which according to these authors allow application of RBF interpolation schemes in large-scale problems. Although simple, the forward differences scheme allowed authors to obtain an excellent match between experimental and observational data.

Multiquadrics are dependent only on space coordinates. This characteristic imposes the use of mixed algorithms to treat time-dependent problems. From Newton iteration to the forward differences scheme, it is possible to combine a wide variety of time-dependent solvers with the basic RBF spatial treatment, used for example in an homogeneous Poisson equation.

In this paper, it is proposed to use Kansa's unsymmetric collocation method [21,22]. For the purpose of completeness, a brief explanation of the method follows.

Consider a boundary-valued problem with a domain $\Omega \subset \mathbb{R}^n$ and a linear elliptic partial differential equation of the form

$$Lu(x) = s(x) \subset \mathbb{R}^n,\tag{6}$$

$$Bu(x)_{|\partial\Omega} = f(x) \in \mathbb{R}^n,\tag{7}$$

where $\partial \Omega$ represents the boundary of the problem. We use points along the boundary $(\mathbf{x}_j, j = 1, ..., N_B)$ and in the interior $(\mathbf{x}_j, j = N_B + 1, ..., N)$.

Let the RBF interpolant to the solution $u(\mathbf{x})$ be

$$s(\mathbf{x},c) = \sum_{j=1}^{N} \phi_j g(\|\mathbf{x} - \mathbf{x}_j\|, c).$$
(8)

Collocation with the boundary data at the boundary points and with PDE at the interior points leads to equations

$$s_B(\mathbf{x},c) \equiv \sum_{j=1}^N \phi_j Bg(\|\mathbf{x} - \mathbf{x}_j\|, c) = \lambda(\mathbf{x}_i), \quad i = 1, \dots, N_B,$$
(9)

$$s_L(\mathbf{x},c) \equiv \sum_{j=1}^N \phi_j Lg(\|\mathbf{x} - \mathbf{x}_j\|, c) = \Phi(\mathbf{x}_i), \quad i = N_B + 1, \dots, N,$$
(10)

where $\lambda(\mathbf{x}_i)$, $\Phi(\mathbf{x}_i)$ are the prescribed values at the boundary nodes and the function values at the interior nodes, respectively.

This corresponds to a system of equations with an unsymmetric coefficient matrix, structured in matrix form as

$$\begin{bmatrix} B\phi\\ L\phi \end{bmatrix} [\mathbf{a}] = \begin{bmatrix} \lambda\\ \mathbf{\Phi} \end{bmatrix}. \tag{11}$$

It has been shown that the unsymmetric coefficient matrix can become ill-conditioned or singular [40].

The use of globally supported RBFs for large problems can bring problems due to the full populated matrices. To solve this drawback, a localization scheme is advisable. Domain decomposition methods [39,45] and localization of the basis functions [39,46] claim to be able to deal with tens of thousands of nodes.

The present model does not issue such methods, as for problems dealt with in the paper the number of nodes to provide good quality solutions is usually small. For large applications involving analysis of plates and shells, such refined approaches are certainly needed.

3. Solving time-dependent problems

Consider the following general time-dependent problem:

$$\frac{\partial u}{\partial t} + Lu = f(x), \quad x \in \Omega, \tag{12}$$

where Ω is a domain in \mathbb{R}^d , d = 1, 2, ..., n with boundary $\partial \Omega$ and L is some differential operator. We approximate u by \tilde{u} and assume

$$\widetilde{u}(t,x) = \sum_{j=1}^{N} a_j(t)\phi(\|x - x_j\|, c),$$
(13)

where x_j are N distinct data points in Ω , and a_j 's are unknown coefficients to be determined at each time step.

The derivative of the approximating solution on time is

$$\frac{\partial \widetilde{u}}{\partial t} = \sum_{j=1}^{N} \frac{\mathrm{d}a_j}{\mathrm{d}t} \,\phi(\|x - x_j\|, c). \tag{14}$$

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Spatial derivatives are obtained as

$$\frac{\partial \widetilde{u}}{\partial x} = \sum_{j=1}^{N} a_j \frac{\partial \phi}{\partial x} \phi(\|x - x_j\|, c),$$
(15)

$$\frac{\partial^2 \widetilde{u}}{\partial x^2} = \sum_{j=1}^N a_j \frac{\partial^2 \phi}{\partial x^2} \phi(\|x - x_j\|, c) \cdots$$
(16)

Substituting Eqs. (13)–(16) into Eq. (12) and collocating at N points x_i , gives

$$\frac{\partial \widetilde{u}(x_i)}{\partial t} + L\widetilde{u}(x_i) = f(x_i), \tag{17}$$

which can be expressed as

$$\Phi \dot{\mathbf{a}} + \Phi_L \mathbf{a} = \mathbf{f},\tag{18}$$

where **a** is the vector of unknown coefficients (a_j) , **f** is the vector $f(x_i)$ and the matrices Φ and Φ_L are given by $\Phi = \phi(||x - x_j||, c)$ and $\Phi_L = L\phi(||x - x_j||, c)$, respectively.

Eq. (18) can be expressed as

$$\dot{\mathbf{a}} = -\mathbf{\Phi}^{-1}\mathbf{\Phi}_L \mathbf{a} + \mathbf{\Phi}^{-1}\mathbf{f}.$$
 (19)

This is a typical system of first-order linear differential equations. With a time difference scheme applied to Eq. (19), the unknown coefficients **a** can be determined at each time step t if and only if the coefficient matrix $\mathbf{\Phi}$ is solvable.

To illustrate this problem more clearly, the first-order time difference scheme is applied to Eq. (18) obtaining

$$\mathbf{\Phi}\mathbf{a}^{n+1} = \mathbf{\Phi}\mathbf{a}^n - \Delta t \mathbf{\Phi}_L \mathbf{a} + \Delta t \mathbf{f}.$$
 (20)

At each time step, *n*, the right-hand term of Eq. (19) is known, so Eq. (19) is similar to $\Phi \mathbf{a} = \mathbf{f}$, the typical interpolation problem.

The following algorithm for time-dependent problems can now be formulated as

- 1. Initialize at time t := 0, approximate the initial condition using Eq. (13) and then compute the partial differential operator $L = L(u_0, u_x, u_{xx}, ...)$.
- 2. Solve Eq. (19) with some time-marching scheme to obtain coefficients **a** and then compute the solution **u** using Eq. (13) and the operator L at time t := t.
- 3. Correct the boundary values using boundary conditions.
- 4. Put $t := t + \Delta t$ and go to Step 2.

In the present paper, the Crank-Nicolson scheme was also used for parabolic problems.

In this case, the algorithm is established as follows. Consider a parabolic case. In the Crank–Nicolson formulation, the following two expressions hold for function and first derivative:

$$u^{t+\Delta t/2} = \frac{1}{2}(u^t + u^{t+\Delta t}),$$
(21)

$$\dot{u}^{t+\Delta t/2} = \frac{u^{t+\Delta t} - u^t}{\Delta t}.$$
(22)

When applying this scheme, the following equation is obtained:

$$u^{t+1} = u^t + \frac{\Delta t}{2} \left(\left(\frac{\mathrm{d}u}{\mathrm{d}t} \right)^{t+1} + \left(\frac{\mathrm{d}u}{\mathrm{d}t} \right)^t \right)$$
(23)

or

$$u^{t+1} - \frac{\Delta t}{2} \left(\frac{\mathrm{d}u}{\mathrm{d}t}\right)^{t+1} = u^t + \frac{\Delta t}{2} \left(\frac{\mathrm{d}u}{\mathrm{d}t}\right)^t.$$
(24)

The multiquadrics interpolation results in

$$\mathbf{\Phi}\mathbf{a}^{n+1} - \frac{\Delta t}{2}\mathbf{\Phi}_L\mathbf{a}^{n+1} = \mathbf{\Phi}\mathbf{a}^n + \frac{\Delta t}{2}\mathbf{\Phi}_L\mathbf{a}^n.$$
 (25)

The remaining procedure follows the same scheme as in the forward difference approach.

The central difference integration method of the Newmark family was used for hyperbolic problems, such as beams in bending. The algorithm for this case follows a similar radial basis interpolation approach as in the parabolic formulations.

4. Examples and discussion

4.1. 1D heat-conduction problem

Consider a 1D heat-conduction problem, with equation

•

$$\frac{\partial u}{\partial t} - \frac{\partial^2 u}{\partial x^2} = 0, \quad 0 < x < 1$$

with boundary conditions

$$u(0,t) = 0, \quad \frac{\partial u}{\partial x}(1,t) = 0$$

and initial condition u(x, 0) = 1.0. The problem is solved for time $0 \le t \le 1$, using $\Delta t = 0.05$. The grid used for this example is a regular grid as shown in Fig. 1.

As can be seen from Fig. 2 and Table 1, results obtained from the present methodology are in excellent agreement with exact results. Results are less adequate for $N \leq 5$.

Fig. 1. Regular grid used for heat-conduction problem; o — boundary nodes; • — interior nodes.



Fig. 2. 1D heat equation results; $-\Delta$, N = 5; ---, N = 10; $-\Delta$, N = 15.

Table 1

1D heat-conduction problem—a comparison of the present model (N = 5, 10 and 15) with the exact solution and finite element solution of Reddy [47, p. 236]

Time t	Reddy [47]	Exact solution	N = 5	N = 10	N = 15
0.00	1.0000	1.0000	1.1103	1.0000	1.0000
0.10	0.9549	0.9493	0.9607	0.9520	0.9502
0.20	0.7731	0.7723	0.7541	0.7749	0.7726
0.30	0.6006	0.6068	0.5786	0.6097	0.6069
0.40	0.4741	0.4745	0.4430	0.4775	0.4746
0.50	0.3701	0.3708	0.3392	0.3738	0.3709
0.60	0.2890	0.2897	0.2596	0.2926	0.2898
0.70	0.2258	0.2264	0.1987	0.2290	0.2265
0.80	0.1764	0.1769	0.1522	0.1793	0.1769
0.90	0.1378	0.1382	0.1165	0.1403	0.1383
1.00	0.1076	0.1080	0.0892	0.1098	0.1081

4.2. Transverse motion of Bernoulli and Timoshenko beams

Consider the transverse motion of an isotropic beam, clamped at both ends, according to the Euler–Bernoulli beam theory, with equation

$$\frac{\partial^2 w}{\partial t^2} - \frac{\partial^4 w}{\partial x^4} = 0, \quad 0 < x < 1,$$

with boundary conditions

$$w(0,t) = 0, \quad \frac{\partial w}{\partial x}(0,t) = 0, \quad w(1,t) = 0, \quad \frac{\partial w}{\partial x}(1,t) = 0$$

and initial condition

$$w(x,0) = \sin \pi x - \pi x(1-x), \quad \frac{\partial w}{\partial t}(x,0) = 0$$

The problem is solved for time $0 \le t \le 0.15$ using $\Delta t = 0.005$ and the present solution compared with a finite element solution by Reddy [47], using four beam elements and a Galerkin solution. The grid for Bernoulli beams is illustrated in Fig. 3.

The present model presents results in excellent agreement with the exact solution and with the finite element results of Reddy [47]. The agreement improves, as expected, with the increase of grid points, as seen in Table 2 and Fig. 4. The modelling of the Euler-Bernoulli beam needs the consideration of the following boundary conditions, imposed upon the transverse deflection (w) and x-direction bending moment (M_x) :

(a)
$$w = 0$$
 in $x = 0$ and $x = L(\Gamma_1)$,
(b) $M_x = 0$ in Γ_2 .
 $\underbrace{\delta}_{x}$ - Very small real number

Fig. 3. Irregular grids for Euler–Bernoulli beams, Neumann conditions imposed; • — w = 0 condition (Γ_1); * — $M_x = 0$ condition (Γ_2); • — interior nodes.

Table 2

Transient Bernoulli beam—a comparison of the present model (N = 9, 11 and 15) with the exact (Galerkin) solution and finite element solution of Reddy [47, p. 240]

Time <i>t</i>	Reddy [47]	Exact solution	N = 9	N = 11	N = 15
0.00	0.2146	0.2146	0.2146	0.2146	0.2146
0.01	0.2098	0.2157	0.2106	0.2095	0.2089
0.02	0.1951	0.1988	0.2005	0.2001	0.1985
0.03	0.1698	0.1716	0.1655	0.1674	0.1691
0.04	0.1350	0.1356	0.1186	0.1252	0.1304
0.05	0.0935	0.0925	0.0654	0.0738	0.0818
0.06	0.0483	0.0447	0.0214	0.0306	0.0384
0.07	0.0018	-0.0055	-0.0292	-0.0189	-0.0107
0.08	-0.0455	-0.0553	-0.0830	-0.0693	-0.0586
0.09	-0.0923	-0.1023	-0.1397	-0.1245	-0.1111
0.10	-0.1336	-0.1441	-0.1760	-0.1644	-0.1520
0.11	-0.1682	-0.1783	-0.1986	-0.1917	-0.1834
0.12	-0.1932	-0.2034	-0.2091	-0.2048	-0.1996
0.13	-0.2087	-0.2179	-0.2202	-0.2172	-0.2136
0.14	-0.2148	-0.2211	-0.2144	-0.2171	-0.2164
0.15	-0.2111	-0.2129	-0.1904	-0.2022	-0.2094



Fig. 4. Bernoulli beam results; *, Reddy [41]; \rightarrow , Exact; \rightarrow , N=9; \rightarrow , N=11; \rightarrow , N=15.

Due to the collocation strategy, it is necessary to impose the boundary condition defined in (b) in a near point: $M_x = 0$ in $x = 0 + \delta$; $x = L - \delta(\Gamma_2)$, where δ is a very small number. This type of boundary condition produces significant improvement over boundary conditions imposed on regular grids (Fig. 3).

Another method for dealing with superposition of boundary conditions was proposed by Wu and his colleagues [48–54].

The evolution of the transverse displacement with time for larger times is illustrated in Fig. 5. It can be seen that a very regular pattern is obtained without divergence of the solution. It is now considered the same beam problem, but with a Timoshenko formulation with h = 0.01 and 0.1.

The governing equations of the Timoshenko beam with cross section A are given by

$$\rho A \frac{\partial^2 w}{\partial t^2} - \frac{\partial}{\partial x} \left[GAk \left(\frac{\partial w}{\partial x} + \phi \right) \right] = 0, \tag{26}$$

$$\rho I \frac{\partial^2 \phi}{\partial t^2} - \frac{\partial}{\partial x} \left[E I \frac{\partial \phi}{\partial x} \right] + G A k \left(\frac{\partial w}{\partial x} + \phi \right) = 0, \tag{27}$$

where G is the shear modulus, K is the shear correction coefficient, ρ is the material density and I the inertia moment.

In order to present a similar simulation as in the case of the Euler–Bernoulli beam, consider EI = 1.0, $\rho A = 1.0$.

In Fig. 6 and Table 3, results are presented and compared with a Galerkin formulation and finite element results of Reddy [47]. Again, $\Delta t = 0.005$ is used as in Reddy's book, with central



Fig. 6. Timoshenko beam results. \bullet , Reddy [41]; \bullet , Galerkin; $\cdot \bullet \cdot \cdot$, h = 0.01, N = 7; $\cdot + \cdot$, h = 0.01, N = 9; $\cdot \cdot \cdot$, h = 0.01, N = 11; \bullet , h = 0.1, N = 7; $\bullet + \cdot$, h = 0.1, N = 9; $- \cdot \cdot \cdot$, h = 0.1, N = 11.

difference (Newmark) scheme. Results with h = 0.01 or 0.1 agree quite well with both formulations. The grid used in this case is similar to the one illustrated in Fig. 1.

In general, the methodology presents quite good results.

Table 3

Time t	Reddy [47]	Galerkin	h = 0.01			h = 0.1		
			N = 7	N = 9	N = 11	N = 7	N = 9	N = 11
0.00	0.2146	0.2146	0.2146	0.2146	0.2146	0.2146	0.2146	0.2146
0.01	0.2098	0.2157	0.2185	0.2113	0.2096	0.2080	0.2077	0.2079
0.02	0.1951	0.1988	0.1922	0.1972	0.1974	0.1887	0.1883	0.1875
0.03	0.1698	0.1716	0.1436	0.1659	0.1684	0.1622	0.1636	0.1652
0.04	0.1350	0.1356	0.1070	0.1207	0.1294	0.1347	0.1375	0.1379
0.05	0.0935	0.0925	0.0440	0.0723	0.0824	0.1028	0.1059	0.1064
0.06	0.0483	0.0447	0.0236	0.0258	0.0379	0.0620	0.0663	0.0682
0.07	0.0018	-0.0055	-0.0699	-0.0235	-0.0102	0.0140	0.0202	0.0217
0.08	-0.0455	-0.0553	-0.1323	-0.0784	-0.0590	-0.0358	-0.0301	-0.0293
0.09	-0.0923	-0.1023	-0.1792	-0.1304	-0.1103	-0.0815	-0.0754	-0.0726
0.10	-0.1336	-0.1441	-0.1981	-0.1684	-0.1515	-0.1185	-0.1093	-0.1061
0.11	-0.1682	-0.1783	-0.2297	-0.1919	-0.1826	-0.1472	-0.1402	-0.1399
0.12	-0.1932	-0.2034	-0.2224	-0.2086	-0.2006	-0.1735	-0.1693	-0.1666
0.13	-0.2087	-0.2179	-0.1952	-0.2202	-0.2141	-0.1979	-0.1934	-0.1922
0.14	-0.2148	-0.2211	-0.1744	-0.2169	-0.2169	-0.2129	-0.2108	-0.2118
0.15	-0.2111	-0.2129	-0.1242	-0.1923	-0.2079	-0.2142	-0.2161	-0.2157

Transient Timoshenko beam—a comparison of the present model (N = 7, 9 and 11) with the exact solution and finite element solution of Reddy [47, p. 240]

4.3. 2D Poisson problem

Consider the 2D Poisson transient problem, with equation

$$\frac{\partial T}{\partial t} - \nabla^2 T = 1, \quad \text{in } \Omega = \{(x, y): 0 < (x, y) < 1\},\$$

with boundary conditions

$$\Gamma = 0$$
 on $\Gamma_1 = \{\text{Lines } x = 1 \text{ and } y = 1\}, t \ge 0,$

$$\frac{\partial T}{\partial n} = 0$$
 on $\Gamma_2 = \{\text{Lines } x = 0 \text{ and } y = 0\}, t \ge 0$

and initial condition

$$T=0, \text{ in } \Omega.$$

The square grid used in this example is shown in Fig. 7. The problem is solved for time $0 \le t \le 1.0$. In Table 4 and Fig. 8, the present method is compared with the finite element results of Reddy [55] and shows high accuracy.



Fig. 7. Grid for 2D problems.

Table 4

Transient 2D Poisson—a comparison of the present model (N = 5, 9 and 17) with the exact solution and finite element solution of Reddy [55, p. 240]

Time t	Formulation	Temperature	Temperature along the $y = 0$ line $(T(x, 0, t) \times 10)$				
		x = 0.0	x = 0.25	x = 0.50	x = 0.75		
	Reddy [55] (R2)	0.9945	0.9853	0.9264	0.6360		
0.1	5×5 grid	1.0580	1.0182	0.9195	0.6451		
	9×9 grid	0.9872	0.9673	0.8822	0.6258		
	17×17 grid	0.9841	0.9718	0.9020	0.6323		
	Reddy [55] (R2)	1.8115	1.7329	1.4997	0.9612		
0.2	5×5 grid	1.8272	1.7395	1.4881	0.9607		
	9×9 grid	1.7376	1.6699	1.4365	0.9347		
	17×17 grid	1.7257	1.6597	1.4275	0.9274		
	Reddy [55] (R2)	2.2479	2.1432	1.8018	1.1319		
0.3	5×5 grid	2.3058	2.1845	1.8309	1.1147		
	9×9 grid	2.2117	2.1090	1.7742	1.1183		
	17×17 grid	2.1998	2.0982	1.7644	1.1102		
	Reddy [55] (R2)	2.9621	2.8037	2.3065	1.4053		
1.0	5×5 grid	3.0131	2.8413	2.3352	1.4187		
	9×9 grid	2.9308	2.7740	2.2836	1.3941		
	17×17 grid	2.9237	2.7672	2.2767	1.3878		



Fig. 8. 2D Poisson results.

5. Conclusions

In this paper, the multiquadric radial basis function method was applied to the analysis of some time-dependent problems. A one- and two-dimensional conduction and a beam in bending were analyzed. Euler–Bernoulli and Timoshenko formulations were used for the transient analysis of isotropic beams. Results were compared with existing solutions showing excellent performance.

Results showed that the use of unsymmetric colocation strategy gives very good agreement with available theories or previous results for all cases.

This method, based on radial basis functions has very large potential for the solution of structural problems, as a real meshless method, insensible to spatial dimension.

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