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## Mixed discrete least squares meshless method for solving the linear and non-linear propagation problems

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#### 1. Introduction

The interest in meshless methods to solve Partial Differential Equations (PDEs) has markedly grown over the last two decades. Approximating the unknown parameters by using some arbitrary distributed nodes, without needing the pre-defined connectivity of nodes, is the valuable advantage of meshless methods com-

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pared to mesh-based methods such as Finite Element Method (FEM) and Finite Volume Method (FVM). This advantage is rid of or at least moderates the difficulty of mesh generation.

Two major approaches are used in meshless methods to approximate the nodal parameters: i) kernel approximation, and ii) series representation. Kernel approximation is based on the theory of integral interpolation. The kernel approximation is represented using its information in a local influence domain via a weighted integral operation. The consistency is obtained by properly choosing the weight function. Series representation approach uses the polynomial basis functions to approximate the nodal parameters. The consistency is ensured by the completeness of

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the basic functions. Smoothed Particle Hydrodynamic (SPH), which uses the kernel approximation, is one of the earliest meshless methods successfully applied for simulation of different fluid mechanic problems, such as free surface Newtonian and non-Newtonian flows [1], sediment transport [2], and multiphase flows [3]. Koshizuka and Oka proposed Moving Particle Semiimplicit (MPS) method [4]. The MPS method is mainly similar to the SPH method [5], but in MPS method, the partial spatial derivatives are calculated without involving the gradient of a kernel function. This method has been applied to simulate dam break [6], 2D nonlinear fluid-structure interaction problems with free surface [7], and multiphase flow [8] problems. Kolahdoozan et al. also studied the effect of turbulence closure methods on the accuracy of MPS method for the viscous free surface flows [9]. However, it is shown that the kernel function used in the original MPS formula does not ensure continuity of the first derivatives [7,10]. Therefore, some modifications to the MPS method are proposed [11,12]. The computational cost of kernel approximations is less than that of series representation approximations. By contrast, the meshless methods using the series representation (Pascal polynomial) approximations provide higher order consistency leading to higher accuracy. Furthermore, the desired order of consistency can easily be achieved by raising the order of basis function, of course by extra effort, a capability that is absent in kernel approximation methods. This ability is highly useful when dealing with complex problems with high gradient solutions.

The series representation methods use the basic functions based on the Pascal polynomial triangle. The polynomial expansion can be directly used as a trial function to solve the linear problems. However, the main drawback of such approach is that the resulting linear algebraic equations are highly ill-conditioned. Some studies are performed to reduce the number of conditions of the linear system equations. Liu and Young used the multiple-scale in the Pascal polynomial to access accurate and stable solutions to 2D Stokes and inverse Cauchy-Stokes problems [13]. Liu introduced homogenized functions and differencing equations to recover time/space-dependent heat source in the heat conduction equation [14]. Since the number of conditions of the inverse heat source recovery problem was reduced using the method, the method was accurate and stable against large noise.

Element Free Galerkin (EFG) as a series representation method was proposed by Belytshko et al. [15]. The Galerkin procedure was utilized to discretize the governing PDEs, leading to symmetric coefficient matrices. This method was successfully used to investigate various engineering problems such as static and dynamic analyses of shell structures [16], temperature field problems [17], 2D fracture problems [18], and unsteady non-linear heat transfer [19]. The rate of convergence of the EFG method was shown to be higher than that of FEM [19]. Furthermore, the irregular nodal distribution was shown not to affect the efficiency of the EFG method [20]. However, use of a background mesh is unavoidable for numerical integrations due to using the weak-form of the governing equations. Therefore, some researchers do not consider the EFG method as a truly meshless method. The Meshless Local Petrov-Galerkin (MLPG) was proposed by Alturi and Zhu [21]. In this method, the shape functions were constructed using a series representation method, such as Moving Least Squares (MLS) approximation. The MLPG has been extensively used to solve a wide range of solid mechanic [22] and fluid flow and heat transfer [23] problems. A local weak-form is used in this method to avoid the use of background mesh. However, the method suffers from a major drawback of asymmetric coefficient matrix and difficulties arise in numerical integration procedure on and around the boundary nodes.

The least squares concept has been widely applied for solving PDEs. For example, Luan and Sun used the least squares method to solve a scattering problem in near field optics [24]. Recently, Discrete Least Squares Meshless (DLSM) method was proposed by Afshar et al. to solve the elliptic and convection-dominated problems [25-28]. The MLS approximation, a common type of the series representation, was used for interpolation (approximation) purposes. DLSM method has been successfully employed to solve solid [29,30] and fluid mechanic problems [31,32]. The DLSM as a truly meshless method uses the strong form of the governing equations. More recently, mixed formulation was applied in DLSM method leading to the Mixed Discrete Least Squares Meshless (MDLSM) method. The mixed formulation was employed in least squares finite element method [33,34]. In MDLSM method, the order of involved derivatives decreases by one unit by using the mixed formulation and, therefore, the complex and costly second derivatives of the shape functions are not required. However, the dimension of the coefficient matrix is larger. The larger coefficient matrix imposes an extra computational cost to solve the resulting linear algebraic system of equations. On the other hand, eliminating the cumbersome second order derivatives reduces the required computational effort of the MDLSM method. Mixed formulation also provides the possibility of simultaneously calculating both unknown parameters and their gradients without any post-processing procedure that is essential in the DLSM method for computing the gradients. Since the post-processing procedure involves less accurate derivatives of shape functions, the gradients are accurately computed in the MDLSM method compared to the DLSM method [30,35,36]. The MDLSM method is

based on minimizing a least squares functional with respect to the nodal parameters. The least squares functional is calculated as a summation of the squared residuals of the governing differential equation and its boundary conditions at the nodal points. Usually, non-satisfaction of LBB condition can produce some difficulties when solving incompressible Navier-Stokes equations, requiring special shape functions. The coefficient matrix of MDLSM method is always symmetric positive-definite so that it is not subject to difficult LBB condition unlike the least squares mixed finite element method [33,34,37]. Hence, the method overcomes the critical difficulties that arise as a result of the non-satisfaction of the LBB. This property is more useful to solve the Navier-Stokes equations [37] and will be investigated in future studies. MDLSM method was successfully employed to solve equilibrium problems such as linear elasticity problem [30,36] and linear quadratic differential equations [35]. The obtained results indicated high accuracy and efficiency of MDLSM method compared to the DLSM method.

The propagation PDEs are dominant in a variety of physical problems such as pollution transport, unsteady fluid, and heat transfer. Exploiting the mentioned merits for the MDLSM method, the MDLSM method is developed in this paper for solving both linear and non-linear propagation problems. The performance of the proposed method is evaluated by several benchmark problems.

#### 2. Moving Least Squares (MLS) approximation method

Several methods were developed to approximate or interpolate the nodal parameters such as Partition of Unity (PU) [38], Radial Basis Function (RBF) [39], Moving Kriging (MK) [40], Maximum Entropy (Max-Ent) [41], Local Maximum Entropy (LME) [42], and Radial Point Interpolation Method (RPIM) [43,44]. These methods are classified in two major groups: i) interpolation, and ii) approximation. The methods based on interpolation satisfy the Kronecker delta function property, which facilitates imposing the essential boundary conditions. However, most of these methods involve some free parameters. The MLS method as an approximation is not subject to this drawback as it only requires the radial of support domain as a free parameter. The MLS approximation is the most popular method for calculating the shape functions in meshless methods [45-47]. However, this approximation does not satisfy the Kronecker delta function property. The penalty approach is an appropriate method to impose the essential boundary conditions when using MLS method [45].

In the MLS method, the unknown function, u, is defined as follows:

$$\mathbf{u}(\mathbf{X}) = \sum_{i=1}^{k} p_i(\mathbf{X}) \mathbf{a}_i(\mathbf{X}) = \mathbf{P}^{\mathrm{T}}(\mathbf{X}) \mathbf{a}(\mathbf{X}), \qquad (1)$$

where,  $\mathbf{a}(\mathbf{X})$  is the coefficient vector;  $\mathbf{X}$  is the vector of the nodal coordinates;  $p_i$  is the component of  $\mathbf{P}$  matrix (the polynomial basis function) defined as Eq. (2); and k is the number of terms in the basis function that guarantees the required consistency:

$$\mathbf{P}^{\mathrm{T}}(\mathbf{X}) = [1, \langle \mathbf{x} \rangle_{1}, \langle \mathbf{x} \rangle_{2}, ..., \langle \mathbf{x} \rangle_{\mathrm{r}}, ..., \langle \mathbf{x} \rangle_{\mathrm{n}}]_{1 \times k}, \qquad (2)$$

in which:

$$\langle \mathbf{x} \rangle_{1} = [x_{1}, x_{2}, \dots, x_{i}, \dots, x_{nd}],$$

$$\langle \mathbf{x} \rangle_{r} = [x_{1}^{r}, \dots, x_{1}^{\gamma_{1}} x_{2}^{\gamma_{2}} \dots x_{i}^{\gamma_{i}} \dots x_{nd}^{\gamma_{nd}}, \dots, x_{nd}^{r}], \sum_{i=1}^{nd} \gamma_{i} = r,$$

$$(3)$$

where, n is the order of basis function;  $x_1$ ,  $x_2$ , and  $x_i$  are the elements of vector  $\mathbf{x}$ ; and nd is the size of  $\mathbf{x}$ .

A weighted discrete  $L_2$  norm, Z, is defined by Eq. (4) as follows:

$$Z = \sum_{j=1}^{n_s} w_j (\mathbf{X} - \mathbf{X}_j) (\mathbf{P}^T (\mathbf{X}_j) \mathbf{a} (\mathbf{X}) - \hat{u}_j)^2, \qquad (4)$$

where,  $\mathbf{X}_{j}$  represents the nodal points with influence domains covering the point  $\mathbf{X}$  (center of support domain);  $\hat{u}_{j}$  is the nodal parameter of  $\hat{u}$  at *j*th node and defines the number of nodes in the support domain; and  $w_{j}$  is the weight function. In this paper, a cubic spline weight function is used as follows:

$$w_{j}(d) = \begin{cases} \frac{2}{3} - 4d^{2} + 4d^{3} & d \leq \frac{1}{2} \\ \frac{4}{3} - 4d + 4d^{2} - \frac{4}{3}d^{3} & \frac{1}{2} \leq d \leq 1 \\ 0 & d \geq 1 \end{cases}$$
$$d = \|\mathbf{X} - \mathbf{X}_{j}\| / d_{wj}. \tag{5}$$

Here,  $d_{wj}$  is the radius of the support domain at *j*th node defined as follows:

$$d_{wj} = \lambda D_k,\tag{6}$$

where  $D_k$  is the distance of the *j*th node from the *k*th nearest point and  $\lambda$  is a constant parameter.

The approximate nodal values can be calculated by minimizing Z with respect to  $\mathbf{a}(\mathbf{X})$  as follows:

$$u(\mathbf{X}) = \mathbf{N}(\mathbf{X})\hat{\mathbf{u}},\tag{7}$$

where,  $\hat{\mathbf{u}}$  is the vector of nodal parameters defined by:

$$\hat{\mathbf{u}}^{T} = [\hat{u}_{1}, \hat{u}_{2}, ..., \hat{u}_{ns}], \tag{8}$$

and  $\mathbf{N}(\mathbf{X})$  is the MLS shape function defined as:

$$\mathbf{N}(\mathbf{X}) = \mathbf{P}^{T}(\mathbf{X})\mathbf{E}^{-1}(\mathbf{X})\mathbf{G}(\mathbf{X}),$$
(9)

 $\mathbf{E}(\mathbf{X})$  and  $\mathbf{G}(\mathbf{X})$  are defined as follows:

$$\mathbf{E}(\mathbf{X}) = \sum_{j=1}^{n_s} w_j (\mathbf{X} - \mathbf{X}_j) \mathbf{P}(\mathbf{X}_j) \mathbf{P}^T(\mathbf{X}_j), \qquad (10)$$

$$\mathbf{G}(\mathbf{X}) = [w_1(\mathbf{X} - \mathbf{X}_1)\mathbf{P}(\mathbf{X}_1), w_2(\mathbf{X} - \mathbf{X}_2)\mathbf{P}(\mathbf{X}_2),$$
$$\dots, w_{n_s}(\mathbf{X} - \mathbf{X}_{n_s})\mathbf{P}(\mathbf{X}_{n_s})].$$
(11)

A necessary condition for the moment matrix (E(X))to be invertible is that there are at least k nodes covered in the support domain of every collocated point [48,49]. Hence,  $\lambda$  should be greater than or equal to 1 for satisfying the minimum number of nodes within the influence domain. For ensuring a well-posed momentum matrix and ensuring its invertibility, it is usually let  $(n_s \gg k)$  [45,50]. However, the shape functions tend to be more and more linearly dependent in the local area with increasing number of nodes covered in the support domain [50,51]. Determining the best value for the number of nodes covered in the support domain is an open problem [49].

The first order of derivatives can be calculated by:

$$\frac{\partial \mathbf{N}}{\partial x_i} = \frac{\partial \mathbf{P}^T}{\partial x_i} \mathbf{E}^{-1} \mathbf{G} + \mathbf{P}^T \frac{\partial \mathbf{E}^{-1}}{\partial x_i} \mathbf{G} + \mathbf{P}^T \mathbf{E}^{-1} \frac{\partial \mathbf{G}}{\partial x_i}.$$
 (12)

#### 3. Proposed Mixed Discrete Least Squares Meshless (MDLSM) method

Consider the following PDE governing a typical propagation problem:

$$\frac{\partial C(\mathbf{x},t)}{\partial t} + \mathbf{V}(\mathbf{x},t) \cdot \nabla C(\mathbf{x},t) = \nu \nabla^2 C(\mathbf{x},t) + S(\mathbf{x},t).$$
(13)

Subject to the following Dirichlet and Neumann boundary conditions:

$$\begin{cases} C(\mathbf{x},t) = \bar{C}(\mathbf{x},t), \\ \nabla C(\mathbf{x},t) = \nabla \bar{C}(\mathbf{x},t), \end{cases}$$
(14)

where C denotes the unknown variable of the problem; **V** is the velocity vector; **x** is the position vector; t is time;  $\nu$  and S are the diffusion coefficient and the source term, respectively; and  $\nabla$  and  $\nabla^2$  are gradient and Laplace operators, respectively.

A semi-discretization is first carried out using the relaxation method in time as follows:

$$C^{n+1} + \Delta t \theta (\mathbf{V}^n \cdot \nabla C^{n+1} - \nu \nabla^2 C^{n+1}) = g^n,$$
  
$$g^n = C^n - \Delta t (1 - \theta) (\mathbf{V}^n \cdot \nabla C^n - \nu \nabla^2 C^n) + S(\mathbf{x}, t),$$
  
(15)

here,  $\theta$ , is the relaxation parameter with a value between zero and one. The superscripts denote the time steps.

The following definition of gradients is used in mixed formulation:

$$\nabla C^{n+1} = \mathbf{q}^{n+1}.\tag{16}$$

Using  $\mathbf{q}^{n+1}$  to represent the gradient term  $\nabla C^{n+1}$  in Eq. (15) leads to the following system of equations to be solved for  $C^{n+1}$  and  $\mathbf{q}^{n+1}$ :

$$C^{n+1} + \Delta t \theta (\mathbf{V}^n \cdot \mathbf{q}^{n+1} - \nu \nabla \cdot \mathbf{q}^{n+1}) = g^n,$$
  

$$\nabla C^{n+1} - \mathbf{q}^{n+1} = \mathbf{0}.$$
(17)

The set of Eq. (17) can be written in a compact form as:

$$\mathbf{A} \cdot \nabla \boldsymbol{\varphi}^{n+1} + \mathbf{B} \boldsymbol{\varphi}^{n+1} = \mathbf{G}^n; \quad \mathbf{G}^n = [\mathbf{0}, g^n]^T, \quad (18)$$

subject to the following Dirichlet type boundary condition:

$$\boldsymbol{\varphi} = \bar{\boldsymbol{\varphi}}; \quad \bar{\boldsymbol{\varphi}} = [\bar{C}, \bar{\mathbf{q}}]^T,$$
(19)

where,  $\varphi$  is the vector of unknown nodal values defined as:

$$\boldsymbol{\varphi} = [C, \mathbf{q}]^T. \tag{20}$$

A and B are defined by the following matrices:

Γ.

$$\mathbf{A} = [\mathbf{A}_{1}, \mathbf{A}_{2}, \dots, \mathbf{A}_{i}, \dots, \mathbf{A}_{nd}],$$

$$\mathbf{B} = \begin{bmatrix} \mathbf{0} & -\mathbf{I} \\ 1 & \Delta t \theta. \mathbf{V}^{n} \end{bmatrix},$$

$$\mathbf{A}_{i} = \begin{bmatrix} \mathbf{e}_{i} & \mathbf{0} \\ 0 & -\nu \Delta t \theta \mathbf{e}_{i}^{T} \end{bmatrix},$$

$$\mathbf{e}_{i}^{T} = [e_{1}, e_{2}, \dots, e_{j}, \dots, e_{nd}] \quad e_{j} = \begin{cases} 1 & i = j \\ 0 & i \neq j \end{cases}$$
(21)

where **I** is identity matrix of size  $nd \times nd$ . nd is equal to the dimension of the problem. Eq. (7) is used for approximating the unknown nodal values in terms of the unknown nodal parameters as follows:

$$\boldsymbol{\varphi} = \mathbf{N}(\mathbf{X})\hat{\boldsymbol{\varphi}}, \quad \hat{\boldsymbol{\varphi}} = [\hat{C}_1, \hat{\mathbf{q}}_1, ..., \hat{C}_{nt}, \hat{\mathbf{q}}_{nt}]^T,$$
$$\mathbf{N}_{m,l}(X) = \begin{cases} N_j(\mathbf{X}), & l = (nd+1)(j-1) + m, \\ 0, & \text{otherwise}, \end{cases}$$
$$m = 1, 2, ..., nd + 1, \quad j = 1, 2, ..., nt, \qquad (22)$$

where nt is the total number of nodes used to discretize the problem domain and  $N_j(\mathbf{X})$  is the shape function of the *j*th node at  $\mathbf{X}$ . Similarly, the gradient of the nodal values can be approximated as follows:

$$\nabla \boldsymbol{\varphi} = \nabla \mathbf{N}(\mathbf{X}) \hat{\boldsymbol{\varphi}}.$$
(23)

Substituting Eqs. (22) and (23) into Eqs. (18) and (19)

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leads to the residuals of differential equation  $(R_{\Omega})$  and Dirichlet boundary condition  $(R_{\Gamma})$  defined as:

$$\begin{cases} R_{\Omega}^{n+1}(\mathbf{X}) = (\mathbf{A} \cdot \nabla \mathbf{N}(\mathbf{X}) + \mathbf{B} \mathbf{N}(\mathbf{X}))\hat{\boldsymbol{\varphi}}^{n+1} - \mathbf{G}^{n}, \\ R_{\Gamma}^{n+1}(\mathbf{X}) = \mathbf{N}(\mathbf{X})\hat{\boldsymbol{\varphi}}^{n+1} - \bar{\boldsymbol{\varphi}}^{n+1}. \end{cases}$$
(24)

The least squares functional of residuals  $R^{n+1}$  at all collocation points in time step n+1 is defined as follows:

$$R^{n+1} = \frac{1}{2} \left( \sum_{j=1}^{nt} \left( R_{\Omega}^{n+1}(\mathbf{X}_{j}) \right)^{2} + \alpha \sum_{j=1}^{nb} \left( R_{\Gamma}^{n+1}(\mathbf{X}_{j}) \right)^{2} \right),$$
(25)

where nb is the number of nodes on the boundaries. Since the MLS does not enjoy the Kronecker delta function property, the boundary conditions cannot be imposed directly.

Two approaches have been used to impose the boundary conditions: The Lagrange multipliers method and the penalty method. The penalty method keeps the coefficient matrix still symmetric, positively defined, and banded. Therefore, the penalty method is used here as a convenient alternative approach to impose the boundary conditions in which the penalty coefficient,  $\alpha$ , should be large enough to satisfy the boundary conditions [50]. However, choosing the appropriate value for the penalty parameter is not straightforward. Using too small or too big penalty value can lead to poor accuracy [51]. Some studies proposed an algorithm for determining the penalty value and investigated the performance of EFG method for different penalty parameter values [49,51,52]. In the practical computations, however, use of a fairly large penalty is usually sufficient to obtain accurate solutions [49,50,51].

Minimizing the least squares functional of residuals (Eq. (25)) with respect to the unknown nodal parameters leads to:

$$\sum_{j=1}^{nt} \frac{\partial (R_{\Omega}^{n+1}(\mathbf{X}_j))}{\partial \hat{\varphi}^{n+1}} (R_{\Omega}^{n+1}(\mathbf{X}_j)) + \alpha \sum_{j=1}^{nb} \frac{\partial (R_{\Gamma}^{n+1}(\mathbf{X}_j))}{\partial \hat{\varphi}^{n+1}} (R_{\Gamma}^{n+1}(\mathbf{X}_j)) = 0.$$
(26)

Eq. (26) yields the symmetric positive-definite system of algebraic equations as follows, which can be solved by iterative procedures such as the conjugate gradient:

$$\mathbf{K}\hat{\boldsymbol{\varphi}}^{n+1} = \mathbf{F}.\tag{27}$$

The coefficient matrix  $(\mathbf{K})$  and right-hand-side vector  $(\mathbf{F})$  are defined as follows:

$$\mathbf{K} = \sum_{j=1}^{nt} \mathbf{L}^{T}(\mathbf{X}_{j}) \mathbf{L}(\mathbf{X}_{j}) + \alpha \sum_{j=1}^{nb} \mathbf{N}^{T}(\mathbf{X}_{j}) \mathbf{N}(\mathbf{X}_{j}),$$
$$\mathbf{L}(\mathbf{X}_{j}) = \mathbf{A} \cdot \nabla \mathbf{N}(\mathbf{X}_{j}) + \mathbf{B} \mathbf{N}(\mathbf{X}_{j}), \qquad (28)$$

$$\mathbf{F} = \sum_{j=1}^{nt} \mathbf{L}^{T}(\mathbf{X}_{j}) \mathbf{G}^{n} + \alpha \sum_{j=1}^{nb} \mathbf{N}^{T}(\mathbf{X}_{j}) \bar{\boldsymbol{\varphi}}^{n+1}.$$
 (29)

Solution to Eq. (27) yields the value of the nodal parameters. Since the MLS shape function is not interpolant, the nodal values of the problem unknown must be retrieved by using Eq. (22) [50].

In the proposed MDLSM method, the complex and costly calculation of the second derivatives of MLS shape functions is not required. Furthermore, both unknown parameters and their gradients are simultaneously computed circumventing the need for the post-processing procedure in the DLSM method for the calculation of derivatives, which involves less accurate second derivatives of the shape functions. Therefore, the gradients of unknown parameters are computed more accurately in the MDLSM method than in the DLSM method.

#### 4. Numerical examples

In this section, several numerical examples are solved with the MDLSM method to investigate its efficiency and accuracy in solving linear and non-linear propagation problems. The value of the penalty coefficient is taken as  $\alpha = 10^8$  for the first four examples.

The constant parameter of support domain,  $\lambda$ , is usually determined by carrying out numerical experiments for a class of benchmark problems. It is generally recommended that a number between two and three will lead to satisfactory results [45,50]. For all examples studied in the current study, the relaxation parameter is chosen as  $\theta = 1$  and the source term is defined as  $S(\mathbf{x}, t) = 0$ . The following error norm is used as error indicator:

$$\operatorname{error} = \frac{\left\| C^{\operatorname{exact}} - C^{\operatorname{MDLSM}} \right\|_{2}}{\left\| C^{\operatorname{exact}} \right\|_{2}},\tag{30}$$

where  $C^{\text{exact}}$  and  $C^{\text{MDLSM}}$  are the vectors of exact and MDLSM solutions, respectively, and  $||.||_2$  is the  $l^2$ -norm.

#### 4.1. One-dimensional Gaussian hill problem

Consider a linear one-dimensional convection-diffusion problem with  $\mathbf{V}(\mathbf{x}, t) = 1$  and  $\nu = 0.005$  in Eq. (13). The exact solution to this problem is defined as follows [53]:

$$C(x,t) = \frac{\xi(0)}{\xi(t)} e^{-\frac{(x-x_0 - \mathbf{v}_t)^2}{2\xi(t)^2}},$$
(31)

where  $\xi(t) = \sqrt{\xi(t) + 2\nu t}$ . The initial hill is centered around  $x_0 = 0.15$  and  $\xi(0) = 0.04$ . The size of the domain is 1.5. The Dirichlet boundary condition on right and left boundaries and initial conditions are defined according to the exact solution. The time step is chosen as 0.001. A quadratic basis function  $\mathbf{P}^{T}(\mathbf{X}) = [1, X, X^{2}]$  is used to produce the shape functions. The problem is solved using the three different uniform nodal distributions. The average error of solutions starts with 0.1784 for the coarse nodal distribution, which decreases to 0.0587 and 0.0425 for the finer distributions of 51 and 76 nodes, respectively, indicating the convergence of the method. Figures 1 and 2 compare the results of the MDLSM method and the exact solutions for two different regular nodal distributions. The problem is solved by the DLSM and MDLSM methods using the 51 nodes distributed uniformly. The gradients of the solutions obtained by the DLSM and MDLSM methods are compared in Figure 3. The results show that the gradients are slightly accurate in MDLSM method. Table 1 describes the CPU times of the DLSM and MDLSM methods. In the MDLSM method, the second order of the derivatives is not required so that the computational effort in approximation procedure (MLS) is less than that in the DLSM method. On the other hand, the computational cost required to solve the resulting system of equations is less in the DLSM method. Therefore, the total computational cost is dependent



Figure 1. Comparison of MDLSM results with the exact solutions for 26 nodes (first example).

on the type of problem being solved. When Eulerian type of simulation is considered, the computational cost of the MDLSM method is higher than that of the DLSM method since the MLS shape functions require to be constructed once. However, for the problems with the moving nodes, as encountered in Lagrangian simulation in which the MLS shape functions need to be constructed for each new nodal position, the computational effort of the MDLSM method would be lower than that of DLSM.



Figure 2. Comparison of MDLSM results with the exact solutions for 76 nodes (first example).



**Figure 3.** Comparison of the gradients of DLSM and MDLSM methods using 51 nodes (first example).

Table 1. Comparison of CPU times of the DLSM and MDLSM methods in one time step (first example).

Number of	MLS	MLS	Solving procedure	Solving procedure
nodes	$(\mathbf{DLSM})$	(MDLSM)	$(\mathbf{DLSM})$	$(\mathbf{MDLSM})$
26	0.005879	0.004556	0.000125	0.00235
51	0.012642	0.008764	0.000238	0.00541
76	0.017396	0.013887	0.000398	0.008247

#### 4.2. One-dimensional wave travel

Assuming the velocity field as  $\mathbf{V}(\mathbf{x}, t) = C$ , the onedimensional form of Eq. (13) can be used to model the one-dimensional nonlinear wave. The size of the domain is assumed to be equal to one. The quadratic basis function is used to solve this problem. The exact solution to this problem is available as follows:

$$C(x,t) = \frac{f_2 + f_1 e^{H(x-mt)}}{1 + e^{H(x-mt)}},$$
(32)

where H and m are defined as:

$$H = \frac{1}{2\nu}(f_2 - f_1), \tag{33}$$

$$m = \frac{1}{2}(f_2 + f_1). \tag{34}$$

 $f_1$  and  $f_2$  are assumed to be equal to zero and one, respectively. The exact solution is used to define the initial and Dirichlet-type boundary conditions. Different regular nodal distributions are used to solve this problem. A series of tests is also carried out to study the sensitivity of the results to different Peclet numbers defined as the dimensionless ratio of the advection term to the diffusion term. For this purpose, the problem is solved with a set of different diffusion coefficients leading to high gradient solutions and higher numerical errors.

Figures 4 and 5 compare the solutions of the MDLSM method with the exact solutions considering the diffusion coefficient as  $\nu = 0.01$ . The problem is also solved for  $\nu = 0.005$ . The results are shown in Figures 6 and 7. In all cases, the ratio of the time step size to the nodal spacing is equal to 0.1.

The time-averaged error of the results is presented in Table 2 to show the sensitivity of the results with respect to the number of nodes and diffusion coefficients. As expected, the results indicate that the error



Figure 4. Comparison of MDLSM results with the exact solutions for 26 nodes and  $\nu = 0.01$  (second example).

MDLSM on t = 0.1٥ 0.0 MDLSM on t = 0.2MDLSM on t = 0.3п 0.8MDLSM on t = 0.4÷ MDLSM on t = 0.50.7Exact solution 0.6 <sub>ლ</sub> 0.5 0.40.30.20.1 0.00.11.0 0.0 0.20.30.40.50.6 0.70.80.9X

Figure 5. Comparison of MDLSM results with the exact solutions for 101 nodes and  $\nu = 0.01$  (second example).



Figure 6. Comparison of MDLSM results with the exact solutions for 26 nodes and  $\nu = 0.005$  (second example).



Figure 7. Comparison of MDLSM results with the exact solutions for 101 nodes and  $\nu = 0.005$  (second example).

Table 2. Sensitivity analysis with respect to the number of nodes and diffusion coefficient (Peclet number) for one-dimensional wave travel problem.

$\mathbf{Number}$	Diffusion	Error	
of nodes	${\rm coefficient}(\nu)$		
25	0.01	0.0773	
50	0.01	0.0161	
100	0.01	0.0060	
25	0.005	0.2435	
50	0.005	0.0720	
100	0.005	0.0123	



Figure 8. Convergence rate for  $\nu = 0.005$  (second example).

is higher for the smaller diffusion coefficients (higher Peclet number) and the results are more accurate for the finer nodal distributions. A set of non-uniform nodal distributions is used to investigate the effect of non-uniform nodal distribution on the accuracy of results. Figure 8 compares the convergence rates of uniform and non-uniform nodal distributions. Using the 50 nodes for  $\nu = 0.005$ , the gradients are computed by the DLSM and MDLSM methods. The error norms of the gradients are 0.0912 and 0.0843 for DLSM and MDLSM methods, respectively, indicating the higher accuracy of the MDLSM.

# 4.3. Burger's one-dimensional equation with periodic boundary conditions

In this section, Burger's non-linear one-dimensional equation  $\mathbf{V}(\mathbf{x},t) = C$ , is considered with periodic boundary conditions. The size of the domain is considered as  $2\pi$ . The MLS shape functions are produced by the cubic basis function. To impose the periodic boundary conditions, the computational domain is extended at the periodic boundaries and the unknown values at the nodes on the right and left sides of the domain are considered equal [54]. Figure 9 illustrates the schematic view of the process. In this



Figure 9. Repetitive domain and the support domains to impose the periodic boundary conditions.



Figure 10. Comparison of MDLSM results with the analytical solutions for 100 regular nodal distributions and  $\nu = 1$  (third example).

figure, *I*-th node represents nodal points on the rightand left-hand sides of the boundaries. The function  $C(\mathbf{x}, t) = 10 \sin(x)$  is used as the initial condition.

Figure 10 compares the results of the proposed MDLSM method with the available analytical solutions for  $\nu = 1$  [55]. This problem is also solved for different values of the diffusion coefficients  $\nu = 0.01$  and 0.001 using different numbers of nodal points distributed uniformly in the domain. The results are shown in Figures 11 to 14. It can be seen that while the results are of high accuracy for the smallest diffusion coefficient, the accuracy of the results decreases with increase in the value of the diffusion coefficient. Furthermore, Figures 11 and 13 show that with coarse nodal distributions of 100 and 150 nodal points, the method is divergent for  $\nu = 0.01$  and  $\nu = 0.001$ , respectively. On the other hand, Figures 12 and 14 indicate that for diffusion coefficient values of 0.01 and 0.001, uniformed distributions of 150 and 512 nodes are enough to produce convergent and accurate results, respectively.

#### 4.4. Two-dimensional Gaussian hill problem

The diffusion of a Gaussian hill in a uniformly rotating flow field controlled by the velocity of  $\mathbf{V}(\mathbf{x},t) = (-\omega(x_2 - O_2), \omega(x_1 - O_1))$  is solved in this section. Here,  $(O_1, O_2)$  is the center of domain,  $x_1$  and  $x_2$  are the components of the coordinate vector, and  $\omega$  denotes



Figure 11. Results of the MDLSM for 100 regular nodal distributions and  $\nu = 0.01$  (third example).



Figure 12. Results of the MDLSM for 150 regular nodal distributions and  $\nu = 0.01$  (third example).



Figure 13. Results of the MDLSM for 150 regular nodal distributions and  $\nu = 0.001$  (third example).



Figure 14. Results of MDLSM for 512 regular nodal distributions and  $\nu = 0.001$  (third example).

the rotational velocity. The analytical solution to this problem on an infinite plate is defined as follows [56]:

$$C(\mathbf{x},t) = \frac{A_0}{1 + (\frac{2\nu t}{\sigma^2})} \exp(-\frac{\tilde{x}_1^2 + \tilde{x}_2^2}{2(\sigma^2 + 2\nu t)}),$$
(35)

where:

$$\tilde{x_1} = (x_1 - O_1) - Q_1 \cos \omega t + Q_2 \sin \omega t$$

and:

$$\tilde{x_2} = (x_2 - O_2) - Q_1 \sin \omega t - Q_2 \cos \omega t$$

 $Q_1$  and  $Q_2$  define the initials of the Gaussian hill center. The constant values considered are  $A_0 = 1$ ,  $\sigma^2 = 2$ ,  $\nu = 10^{-5}$ , and  $\omega = 10^{-6}$ .

The domain of the problem is a square domain measuring  $20 \times 20$  and  $(Q_1, Q_2) = (15, 10)$ . The analytical solution is used to impose the initial and Dirichlet boundary conditions. Linear basis function is used and the results are presented after one rotation produced in 2000 time steps. The domain is discretized by two sets of regular nodal distributions of 441 and 1681 nodes. The errors of the MDLSM method for each set of coarse and fine nodal distributions are 0.1113 and 0.0635, respectively. The analytical and numerical results are compared in Figure 15.

#### 4.5. Burger's two-dimensional equation

Burger's two-dimensional equation can be represented by taking  $\mathbf{V}(\mathbf{x},t) = C$  in Eq. (13). The problem is solved in a unit square domain discredited by a regular nodal distribution of 441 nodes. The exact solution to this problem is available as:

$$C(\mathbf{x},t) = \frac{1}{1 + \exp(x + y - t - 0.25)/(2\nu)},$$
 (36)

where x and y define the coordinates and  $\nu$  is assumed to be 0.03 in this problem. The exact solution is



Figure 15. Comparison of the contour lines of exact and MDLSM solutions (forth example): (a) Exact solution, (b) MDLSM results obtained for regular distribution of 441 nodes, (c) MDLSM results obtained for regular distribution of 1681 nodes, and (d) all results.

used to impose the Dirichlet boundary conditions. The time step size,  $\Delta t$ , is equal to 0.001. The penalty coefficient is taken as  $\alpha = 10^4$ . Figure 16 compares the numerical and the exact results. This comparison shows the high accuracy of the proposed MDLSM method for solving non-linear convection-dominated problems. The problem is also solved with  $\Delta t = 0.01$ to investigate the influence of the penalty parameter, of which the results are demonstrated in Figure 17. The results show that accuracy generally improves by increasing penalty parameter. However, when the penalty exceeds a specified value, the coefficient matrix becomes ill-posed and the error sharply increases. A set of irregular nodal distributions, presented in Figure 18, are also used to study the effect of the irregular nodal configuration on the results. The convergence rate of the method is shown in Figure 19 for the results at t = 0.5 s for both uniform and irregular nodal distributions. The gradients of the solutions obtained on a uniform  $21 \times 21$  nodal distribution are also compared in Figure 20. The obtained results show that the gradients of the solutions are remarkably accurate in MDLSM method compared to DLSM method.



Figure 16. Comparison of MDLSM results with the exact solutions at t = 0.5 and y = 0.5 (fifth example).

In Table 3, the CPU times of the DLSM and MDLSM methods are compared. In the MDLSM method, the cumbersome second derivatives of the MLS are not required and, hence, the CPU time of the

Table 3. Comparison of CPU times of the DLSM and MDLSM methods in one time step (fifth example).

Number of	$\mathbf{MLS}$	$\mathbf{MLS}$	Solving procedure	Solving procedure
$\mathbf{nodes}$	(DLSM)	(MDLSM)	$(\mathbf{DLSM})$	(MDLSM)
36	0.02918	0.01549	0.00041	0.009195
121	0.1171	0.0591	0.01521	0.02476
441	2.3205	0.2417	0.02120	0.24771



**Figure 17.** Influence of the penalty factor  $\alpha$  (fifth example).

approximation procedure is less than that of DLSM method. However, the size of the coefficient matrix is larger in the MDLSM method and the CPU time of the solving procedure increases compared to the DLSM method. Therefore, the total computational cost is dependent on the type of the problem being solved. When Eulerian type of simulation is considered, the computational cost of the MDLSM method is higher than that of the DLSM method since the MLS shape functions require to be constructed once. However, for the problems with the moving nodes as those encountered in Lagrangian simulation, in which the MLS shape functions need to be constructed for each new nodal position, the computational effort of the MDLSM method would be lower than that of DLSM.



Figure 19. Convergence rate of MDLSM method at t = 0.5 (s) for uniform and irregular nodal distributions (fifth example).

#### 5. Conclusion

A truly meshless method, namely, MDLSM, was presented in this paper to solve propagation problems. The method was based on the minimization of the least squares functional with respect to the nodal parameters. The least squares functional was defined as the sum of the squared residuals of the differential equation and its boundary conditions. The MLS shape function was used to approximate the solutions. The coefficient matrix of the MDLSM method is always symmetric positive-definite, circumventing the LBB condition. With the MDLSM method, calculation of second derivatives of shape function was not required and the gradient of solutions was accurately computed



Figure 18. Irregular nodal distribution (fifth example).



Figure 20. Comparison of the gradients of the DLSM and MDLSM methods at t = 0.5 and y = 0.5 (fifth example).

compared with the DLSM method. The efficiency and accuracy of the method were evaluated via some numerical examples. The numerical results were presented and compared with the available analytical solutions. The results indicated high efficiency of the MDLSM method in solving both linear and nonlinear convection-dominant problems. Comparison of the CPU times required by the DLSM and MDLSM methods for the test examples showed that although MDLSM was more expensive than DLSM for problems with fixed nodal point positions, it could be cheaper for the problems with moving nodes, i.e. Lagrangian simulations.

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