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OPTIMAL DECOMPOSITION OF FINITE ELEMENT MESHES VIA K-MEDIAN METHODOLOGY AND DIFFERENT METAHEURISTICS

A. Kaveh^{*,†} and A. Dadras

Centre of Excellence for Fundamental Studies in Structural Engineering, Iran University of Science and Technology, Narmak, Tehran-16, Iran

ABSTRACT

In this paper the performance of four well-known metaheuristics consisting of Artificial Bee Colony (ABC), Biogeographic Based Optimization (BBO), Harmony Search (HS) and Teaching Learning Based Optimization (TLBO) are investigated on optimal domain decomposition for parallel computing. A clique graph is used for transforming the connectivity of a finite element model (FEM) into that of the corresponding graph, and *k*-median approach is employed. The performance of these methods is investigated through four FE models with different topology and number of meshes. A comparison of the numerical results using different algorithms indicates, in most cases the BBO is capable of performing better or identical using less time with equal computational effort.

Keywords: domain decomposition; Finite elements meshes; graph theory; optimization; metaheuristic algorithms; k-median; k-means++.

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1. INTRODUCTION

Parallel processing was proposed to perform computational tasks in shorter time. In this mode of operating, a process is split into parts and each of which are executed by different processors simultaneously. In engineering, sometimes heavy computational works are required to solve the problems, where this approach is necessary to reduce the computing time. Finite element analysis as an important case for the time consuming problems can be performed through this approach. In this, given a number of available processors q, the finite meshes of the main model are decomposed into q subdomains and each subdomain is

^{*}Corresponding author: Centre of Excellence for Fundamental Studies in Structural Engineering, Iran University of Science and Technology, Narmak, Tehran-16, Iran alikaveh@iust.ac.ir (A. Kaveh)

analyzed by one processor. This means, all formulation of the equations, assembling and solution of the system of equations for each one of q subdomains are carried out in parallel. The total time required for this mode is slightly more than that of the longest subtask, hence an optimum decomposition that yields equal execution times is desired. Finding this optimum decomposition is the main objective of this paper.

In mathematics the pointed problem belongs to a family of problems in graph theory, called 'finding medians of a graph' which is an NP-hard problem [1]. The exact solution of this problem, itself is highly time consuming, hence the approximate methods are suggested. Researchers have developed various methods to deal with domain decomposition especially for FEM [2-11]. A simple and powerful approach to tackle them is *k*-median method which can be simplified as follows. A graph is associated with the considered finite element. Then, the optimal medians of the graph are selected through the sum of the distances of nodes to optimum medians. Metaheuristics are suitable tools to solve this problem approximately. Recently a significant number of metaheuristics are proposed and applied to different engineering problem [12-15]. Of these, most well known Genetic algorithm [16,17], bionomic [18, 19], ant colony [20] and colliding bodies optimization [21] are developed in order to obtain solutions for *k*-median problem.

In this paper, four well-known algorithms; Artificial Bee Colony (ABC) [22], Biogeographic Based Optimization (BBO) [23], Harmony Search (HS) [13] and Teaching Learning Based Optimization (TLBO) [24] are implemented for optimal domain decomposition of finite element meshes. Since most of the referred papers used random initialization, authors also studied the effects of a special initialization technique so called k-means++ [25]. The developed programs are applied to four benchmark problems with different numbers of subdomains. The efficiency of the algorithms is compared in terms of statistical results, convergence rate and consumed running times.

The rest of this paper is organized as follows. The domain decomposition using *k*-median methodology and utilized initialization methods are comprehensively outlined in section 2. In Section 3 the reviews of different algorithms are provided. The experiment results are presented in section 4, and finally the conclusions are derived in section 5.

2. DOMAIN DECOMPOSITION USING K-MEDIAN METHODOLOGY

2.1 Preliminaries from graph theory

A simple graph G is defined as a set N(G) of nodes and a set E(G) of edges together with a relation of incidence which associate two distinct nodes with each edge. Some of distinct nodes of graph are adjacent i.e. they are connected together with an edge. The edge is also called incident with both of its nodes. A simple graph with 9 nodes and 10 edges are drawn in Fig. 1(a). Graphs have several types, in overall the paper, the simple type is considered. A subgraph G_i of a graph is a graph for which $N(G_i) \subseteq N(G)$, $E(G_i) \subseteq E(G)$ and each edge of $E(G_i)$ has the same ends as in G. A path is a finite sequence of nodes and edges like $\{n_0, e_1, n_1, \dots, e_p, n_p\}$, where the nodes (n_{i-1}, n_i) are the ends of the edge (e_i) , while neither edges nor nodes are repeated. Cycle is a closed path with the last node and the first node being the same. A subgraph which has no cycle is known as a tree. A spanning tree of

a graph is a tree containing all the nodes of the graph. The distance between two nodes is the number of edges in the shortest path between these nodes. Shortest route tree (SRT) rooted from the node n is a spanning tree of connected graph G, such that the path from the root n to any other node u of the SRT is the shortest path from n to u in G. The schematic of a sample graph and its SRT rooted from node '1' is illustrated in Fig. 1(b). A clique graph has the same nodes as those of the corresponding finite element model, and the nodes of each element are cliqued, avoiding the multiple edges for the entire graph [26, 27].



Figure 1. Examples form basic concepts from graph theory: (a) A simple graph. (b) An SRT rooted from node '1'

2.2 Domain decomposition method for finite element analysis

As mention in section 2.1, all the FE analysis steps must be performed for each subdomain i.e. each of them can be considered as a super element. A finite element equation system can be assembled for each subdomain as Eq. (1):

$$[k]\{u\} = \{f\}$$
(1)

where [k] is a subdomain stiffness matrix, $\{u\}$ is a subdomain displacement vector, and $\{f\}$ is a subdomain load vector. As illustrated in Fig. 2, the subdomain nodes are grouped into interior nodes and interface boundary nodes designated by the subscripts *i* and *b*, respectively. If the interior nodes are numbered first and the interface boundary nodes are numbered last, then the subdomain equation system can be written in the following matrix form:

$$\begin{bmatrix} k_{ii} & k_{ib} \\ k_{bi} & k_{bb} \end{bmatrix} \begin{pmatrix} u_i \\ u_b \end{bmatrix} = \begin{cases} f_i \\ f_b \end{cases}$$
 (2)

where matrices $[k_{ii}]$ and $[k_{bb}]$ correspond to interior and interface boundary nodes, respectively, and matrix $[k_{ib}]$ reflects the interaction between the interior and boundary nodes. By disassembling the matrix, Eqs. (3 and 4) are obtained as:

$$[k_{ii}] \cdot \{u_i\} + [k_{ib}] \cdot \{u_b\} = \{f_i\}$$
(3)

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$$[k_{bi}]. \{u_i\} + [k_{bb}]. \{u_b\} = \{f_b\}$$
(4)

The determination of the displacements for interior subdomain is:

$$\{u_i\} = [k_{ii}]^{-1}\{\{f_i\} - [k_{ib}], \{u_b\}\}$$
(5)

By substituting Eq. (5) in Eq. (4), the following equation can be obtained:

$$[k_{bi}] \cdot [k_{ii}]^{-1} \{\{f_i\} - [k_{ib}] \cdot \{u_b\}\} + [k_{bb}] \cdot \{u_b\} = \{f_b\}$$
(6)

Denoting super-element stiffness matrix and corresponding modified load vector by $[k^*]$ and $\{f^*\}$ leads to:

$$[k^*] = [[k_{bb}] - [k_{bi}] \cdot [k_{ii}]^{-1} \cdot [k_{ib}]]$$

$$\{f^*\} = \{\{f_b\} - [k_{bi}] \cdot [k_{ii}]^{-1} \cdot \{f_i\}\}$$

$$(8)$$

According to these definitions, Eq. (6) can be rearranged as:

$$[k^*]\{u_b\} = \{f^*\} \tag{9}$$

Thus, structural stiffness matrix and nodal load vector for boundary nodes are assembled using Eq. (7) and Eq. (8), then the displacements of the nodes are evaluated by Eq. (9).

For an instance finite element meshes illustrated in Fig. 2 can be performed by the following steps: first the finite meshes must be partitioned into several submeshes corresponding to the number of available processors. In this example, four processors are considered. The condensed stiffness matrix and modified load vector for each part is calculated by Eq. (7) and Eq. (8), respectively. Then processors are involved by solving Eq. (9) including approximately 1/4 of equations of the global matrix, so the internal nodal displacements are obtained. This is followed by the solution of the equations corresponding to interface nodes followed by the superposition of calculated displacements.

As it be seen the steps are clear, except the first step that points to domain decomposition. An automatic finite element domain decomposer should meet the following requirements [11]:

a. It should be able to handle irregular geometry.

b. It should equally distribute the computational burden among processors.

c. It should yield to minimum interface nodes to reduce the cost of synchronization between the processors.

d. It should be fast and result in a good convergence rate of domain decomposition upon iterative method.



Figure 2. Partitioned finite element meshes

2.3 The k-median of a graph

In the aforementioned problem k is the number of node subsets $N_k \in N$, and there is a median node m_k for each subset. Usually different solutions can be suggested for the same graph, hence, the solution is desired whose subsets have minimum total distances from medians. According to the presented definition, the cost function of this optimization problem can be expressed by a double summation as follows:

$$Z = \sum_{i=1}^{k} \sum_{n_j \in N_k} D(m_k, n_j)$$
(10)

where $D(m_k, n_j)$ is the shortest distances between kth median (m_k) and a node (n_j) of kth domain as defined in section 2.1. As suggested in [11], the clique graph should be plotted in Cartesian coordinate system. If the found components for median are not the same as the plotted nodes, the nearest node is substituted. The medians are evaluated by calculating the graph theoretical distance from all nodes. Nodes with the least distance from a particular median form the domain of that median.

2.4 k-means++

k-means method is a widely used clustering technique that seeks to minimize the average squared distance between points in the same cluster [28]. The basic *k*-means method starts from random initialization. Arthur and Vassilvitskii [25] proposed a probabilistic approach for initialization, so called *k*-means++. The initialization steps of this approach is presented in following:

a. Take one median m, chosen uniformly at random from N.

b. Take a new median m_{new} , choosing $n \in N$ with probability $\frac{D(m,n)^2}{\sum_{n \in N} D(m,n)^2}$.

c. Repeat step b. until we have taken k medians altogether.

In this paper, the k-means++ is used as well as random initialization in order to guide

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metaheuristics. The effects of these approaches are also investigated in the results of experiments.

3. METAHEURISTICS

Metaheuristics are non-deterministic optimization approaches, which do not guarantee that globally optimal solution can be found. However, they find feasible solutions for optimization problems in vicinity of the computation limits [29, 30]. Here, utilized algorithms consisting of the Artificial Bee Colony (ABC), Biogeographic Based Optimization (BBO), Harmony Search (HS) and Teaching Learning Based Optimization (TLBO) are considered and their pseudo codes are provided. In all cases the number of search agents (population) and number of iterations are limited to 20 and 100, respectively.

3.1 Artificial bee colony algorithm

ABC as a swarm intelligence algorithm imitates the foraging behavior of the honey bees [31]. This optimization technique has been successfully applied to various practical problems [32]. In this algorithm there are three types of bees: employed bees, onlooker bees, and scout bees. The bee's aim is to discover the places of food sources with high nectar amount. In which, the employed bees search food around the food source and share the information of these food sources to the onlooker bees. The onlooker bees tend to select good food sources from those found by the employed bees. The selection probability of each source by onlookers is calculated by the following equation:

$$p_s = \frac{f_s}{\sum_j^{Ns} f_j} \tag{11}$$

where f_s is the fitness of the *s*th source and *Ns* is the number of all candidate sources. The scout bees search new food sources and replace with abandoned sources. The steps of this algorithm are represented in the following pseudo code.

Initial food sources are produced for all employed bees
while terminating conditions are not met do
Employed bees:
for $i = 1$ to number of employed bees do
<i>i</i> th employed bee goes to a food source in her memory and determines a
neighbor source evaluates its nectar amount (fitness)
if nectar amount of neighbor is more than the previous food source then
Replace the new neighbor with the previous source
else
Mark it as a candidate to be abandoned
end if
end for
Onlooker bees:

3.2 Biogeographic Based Optimization algorithm

Simons proposed a new population based algorithm so called biogeography based optimization (BBO) algorithm [23]. In biogeography, the suitability of a habitat to live is called 'habitat suitability index' (HSI), according to this definition the amiable places have high HIS and the habitats unfriendly to live have a low HIS. This index pursues some features, which are called 'suitability index variables' (SIVs). Examples of this index are weather conditions, access to food, diversity, safety and so on. Habitats with low HSI have both a high emigration rate (μ) and a low immigration rate (λ); in contrary, habitats with high HSI have a low μ (Eqs. (12 and 13)).

$$\lambda_k = (1 - r_k) \tag{12}$$

$$\mu_S = r_k \tag{13}$$

where r_k is the rank of the *k*th habitat. Suppose a problem is presented with some candidate solutions. A good solution is analogous to a habitat with a high HSI, and a poor solution represents a habitat with a low HSI. High HSI solutions resist the change more than low HSI solutions and tend to share their features with low HSI solutions, while poor solutions accept many new features from the good solutions. The pseudo code of the BBO is presented in following.

Initialize habitats while terminating conditions are not met do *Apply migration* for j=1 to number of habitats do Select habitat H_j according to λ_j if rand $(0,1) < \lambda_j$ then

for $e=1$ to H do
Select habitat H_e according to μ_e
Replace the selected SIV of H_i by selected SIV of H_e
end for
end if
end for
Apply mutation
for $J = 1$ to number of nabitats do
if rand(0,1) < mutation probability then
Replace $H_i(SIV)$ with randomly generated SIV
end if
end for
Sort the population according to the increasing order of fitness
Keep the elite solution
Stop, if termination criterion satisfied
end while

3.3 Harmony search algorithm

Harmony Search (HS) algorithm simulates finding best harmony by musicians [13]. This algorithm is widely utilized in engineering problems [33]. In which, the candidate solutions are considered as harmonies and the best harmony is arranged by utilizing some mechanisms upon several parameters. These parameters are defined as follows:

hms : the size of the harmony memory.

hmcr : the rate of choosing a value from the harmony memory.

par : the rate of choosing a neighboring value.

fw (fret width): the amount of maximum change in pitch adjustment. The new harmony is calculated by the following equation:

$$x_n = x_p + fw \times r \tag{14}$$

where x_n, x_p, fw and r are harmony new position, harmony previous position, fret width and random number within (-1,1), respectively.

Initialize harmony memories as many as hms
Evaluate all memorized harmonies
while terminating conditions are not met do
Initialize new harmonies as many as <i>candidate solutions</i> (<i>nNew</i>)
for i=1: nNew do
for $j=1$ to number of variables do
if $rand(0,1) \leq hmcr$ then
Choose the <i>j</i> th variable from a randomly selected vector of harmony memory
Replace the chosen variable with the <i>j</i> th variable of the <i>i</i> th new harmony

end if
if $rand(0,1) \leq par$ then
Change the <i>j</i> th variable of the <i>i</i> th new harmony by Eq. (14)
end if
end for
Evaluate the ith harmony
end for
Consider all harmonies containing new created harmonies and harmony memories and
save <i>hms</i> best harmonies
end while

3.4 Teaching learning based optimization algorithm

This population-based method is motivated by the influence of a teacher on learners. The population is considered as a group of learners and the process of TLBO is divided into two parts: the first part consists of the 'Teacher Phase' and the second part consists of the 'Learner Phase'. In 'Teacher Phase' teacher learns and in 'Learner Phase' learning is performed by the interaction between learners. In the first phase, the position of learners are updated by Eq. (15):

$$L_{new}^{i} = L_{old}^{i} + Random \times (T - TF \times M)$$
⁽¹⁵⁾

where L_{new}^i , L_{old}^i , Random, T, TF and M are new position of the *i*th learner, old position of the *i*th learner, random number within (0,1), teacher's position, teaching factor and mean position of class.

In the learner phase, the position of learners are updated by Eqs. (16.a, 16.b):

$$L_{new}^{i} = \begin{cases} L_{old}^{i} + Random \times (L_{old}^{i} - L^{rand.}) & \text{if } C(L_{old}^{i}) < C(L^{rand.}) \\ L_{old}^{i} - Random \times (L_{old}^{i} - L^{rand.}) & \text{if } C(L_{old}^{i}) > C(L^{rand.}) \end{cases}$$
(16a)
(16b)

where L^{rand} is a random learner who interacts with *i*th learner and C is the cost function.

The pseudo code of TLBO is provided here:

Initialize individuals randomly

while terminating conditions are not met do

Evaluate all individuals and determine the teacher with the minimum cost and mean position of learners

Teacher Phase for i=1: number of learners do Update the position of learners by Eq. (15) if $C(L_{new}^i) < C(L_{old}^i)$ then Replace L_{new}^i with L_{old}^i end if end for A. Kaveh and A. Dadras

Learner Phase	
for j=1: number of learners do	
Update the position of learners by Eqs. (16 a, 16 b)
if $C(L_{new}^j) < C(L_{old}^j)$ then	
Replace L_{new}^{j} with L_{old}^{j}	
end if	
end for	
end while	

4. NUMERICAL RESULTS

In this section, four well-known algorithms described in section 3 with random initialization and k-means++ initialization approaches, are examined by four FE models. The algorithms are ran 30 times and the statistical results and convergence histories are obtained. These results are obtained by Core TM i7-5500U CPU @ 2.4 GHz and the clock time of computations are calculated. Additional information about the models and results are presented in the following subsections.

4.1 Square plate

The first example is a rectangular plate with $2601(51\times51)$ meshes plotted in Fig. 3. This model with 3,4,5 and 6 medians was studied by Kaveh and Mahdavi [11] to decompose optimal domains using Colliding Bodies Optimization (CBO) algorithm [34]. The number of medians are set according to the referred article and the statistical results are presented in Table 1. Convergence histories are depicted for mean of 30 independent trial in Fig. 4.



Figure 3. FE models of Example 1. (a) FE meshes. (b) The corresponding clique graph

					1					
			ABC		BBO		HS		TLBO	
			Random	k-means++	Random	k-means++	Random	k-means++	Random	k-means++
	k=6	19958.8	20045.8	18936.6	18935.8	19520.7	19472.9	19031.9	19026.6	
0.00	ago	k=5	21272.7	21220.6	20509.7	20509.6	20814.4	20761.2	20537	20553.7
ct Avers		k=4	22681.5	22878.2	22113	22113	22148.1	22131.8	22153.5	22158.9
	7	k=3	28160.7	28151.7	28097.2	28098.3	28099.2	28102.7	28097.9	28100
		k=6	19607	19696	18921	18921	19270	19180	18944	18941
ŧ	7	k=5	20903	20750	20505	20505	20619	20600	20505	20505
Rect		k=4	22259	22405	22113	22113	22113	22113	22113	22113
		k=3	28100	28107	28097	28097	28097	28097	28097	28097
		k=6	238.1	193.1	25.6	24.7	131.2	143.4	67.6	58.9
lard	tion	k=5	264.7	282.4	3.9	4.6	131.4	115.7	34.9	48.3
tanc	evia	k=4	240.2	300.9	0	0	54.3	35.4	2.4	61.9
01	р	k=3	35.6	26.3	0.6	4.9	5.1	11.1	18944	7.2
		k=6	54.9	51.9	27.1	25	26	28	50	52.3
PU time Sta	$\hat{\mathbf{u}}$	k=5	52.2	51.3	26	24.7	27.3	26.7	50.2	49.5
	(sec	k=4	51.6	51	25.4	24.4	27.7	27.7	50.8	49
U		k=3	53.2	51.2	26.2	24.2	26.1	27.5	48.6	48.8

Table 1: Comparison of the statistical results for Example 1







Figure 4. The average convergence curves obtained in 2601 meshes FE model: (a) k=3. (b) k=4. (c) k=5. (d) k=6

As presented in Table 1, the BBO outperformed other considered algorithms. The BBO outperformed CBO for k=5 and 6 cases and obtained the same results for k=3 and 4 presented in [13]. As written in bold font, in some cases using k-means++ improved the results of BBO, while this statement is not true about all cases. Fig. 4 shows higher convergence rate of the BBO and TLBO. Fig. 5 illustrates the optimal domains with distinct colors and medians positions with black filled circles.

4.2 H-shaped finite element model

This FEM is consider with 4949 meshes as illustrated in Fig. 6. The number of 5 and 10 subdomains are considered and the statistical results are presented in Table 2. Average convergence histories are illustrated in Fig. 7.







Figure 5. A FEM decomposed into k subdomains using the BBO algorithm: (a) k=3. (b) k=4. (c) k=5. (d) k=6



Figure 6. The FE model of Example 2

Г٤	ıb	le 2	2:	C	omparison	of	the	statistical	resul	ts f	for	Examp	le	2
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		ABC		B	BO]	HS	TLBO		
		Random	$n \frac{k}{\text{means}++} \text{Random} \frac{k}{\text{means}++} \text{H}$		Random <i>k</i> -means++		Random	k-means++		
A	k=5	60311.1	60030.4	55136.7	55008	56601.4	56637.7	55265	60311.1	
Average	k=10	45326.2	44938.7	38198.6	37896.1	41934.4	42256.3	38760.7	45326.2	
Deet	k=5	57719	58308	55008	55008	55186	55347	55013	57719	
Dest	k=10	43699	41932	37717	37717	40869	41341	37900	43699	
Standard	k=5	1031.9	1168.3	704.9	0	1039.1	899	292.4	1031.9	
deviation	k=10	853.4	991.9	578.6	195.6	581.7	446.8	888.9	853.4	
CPU time	k=5	96.6	93.3	50.3	47.7	51.1	50.7	107.3	96.6	
(sec)	k=10	104.3	97.3	50.6	51.2	53.2	53	100.9	104.3	



Figure 7. The average convergence curves obtained by different methods in Example 2: (a) k=5. (b) k=10

According to Table 2, the BBO obtained the better optimum results, also, using k-means++ improved the robustness of the method. As illustrated in Fig. 7, the BBO especially with k-means++ initialization reached the higher average convergence rate in both cases. Zero value for standard deviation of the BBO in the first case shows that has it led to 55008 for the cost function in all 30 trial, which is the best so far reported result. The optimal solutions are displayed in Fig. 8.



Figure 8. A FEM divided into k subdomains using the BBO algorithm: (a) k=3. (b) k=4. (c) k=5. (d) k=6

4.3 Rectangular plate with four openings

As shown in Fig. 9, this rectangular plate with 4 openings, has 760 meshes. The numbers of medians considered as $k=\{5 \text{ and } 10\}$. The results are presented in Table 3 and the average convergence histories are depicted in Fig. 10.



Figure 9. The FE meshes of example 3

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Algorithm Initialization Average Best deviation	CPU time (sec)
k=5 $k=10$ $k=5$ $k=10$ $k=5$ $k=1$	0 <i>k</i> =5 <i>k</i> =10
ADC Random 4812.5 3465.5 4670 3362 63.9 47	4 18.4 19.5
ABC k -means++ 4815.6 3475.6 4694 3372 67.6 48	2 17.8 18.9
Random 4488.2 2971.1 4467 2934 65.4 20.	1 9.1 9.8
k-means++ 4492.5 2962.9 4467 2934 50.2 31	9.1 9.8
Random 4582.3 3304.6 4502 3215 45 33.	3 8.9 9.5
k-means++ 4582.5 3296.3 4482 3202 62.4 51.	2 9 9.7
TLPO Random 4507.9 3055.4 4467 2968 39.7 41.	4 19.6 20.2
<i>k</i> -means++ 4513.5 3064.9 4467 2959 36.5 61	3 19.1 19.5

Table 3: Comparison of the statistical results for Example 3



Figure 10. The average convergence histories of Example 3 obtained by different methods: (a) k=5. (b) k=10

As presented in Table 3, the BBO outperformed other methods in terms of average cost and best cost, while HS completed the optimization process in lower CPU time. According to Fig. 10 the convergence speed of the BBO is more than other compared methods. The optimal domains and medians of both cases are plotted in Fig. 11.



Figure 11. A FEM divided into k subdomains using the BBO algorithm: (a) k=5. (b) k=10

4.4 Perforated circular plate

A circular plate is considered as shown in Fig. 12. This FEM contains 1152 meshes and the number of subdomains are set to $k = \{4 \text{ and } 5\}$. The statistical results are provided in Table 4 and convergence curves are plotted in Fig. 13.



Figure 12. The FE meshes of a circular plate

Algorithm	Initialization	Ave	Average		Best		Standard deviation		CPU time (sec)	
		<i>k</i> =5	<i>k</i> =10	<i>k</i> =5	<i>k</i> =10	<i>k</i> =5	<i>k</i> =10	<i>k</i> =5	<i>k</i> =10	
APC	Random	7562.6	5623.1	7283	5486	100.5	60.6	28.1	29.7	
ADC	k-means++	7589.7	5651.5	7325	5549	98.3	54.2	29.1	31.1	
PRO	Random	7271	5261.6	7230	5210	58.1	35.5	13.5	15.2	
bbo	k-means++	7312.1	5257.3	7230	5188	107.7	36.5	13.4	15.2	
ЦС	Random	7540.2	5605.6	7413	5419	71	55.1	13.2	14.8	
пз	k-means++	7521.6	5602.6	7296	5489	95.1	49.3	14.3	16.9	
	Random	7257.6	5281.1	7230	5210	26.1	41	26.8	30.6	
ILDU	k-means++	7273	5292.2	7230	5207	35.2	50.7	27.4	30	

 Table 4: Comparison of the statistical results for example 4



Figure 13. The average convergence histories of circular plate obtained by different methods: (a) k=5. (b) k=10

In this case, the best results are obtained by BBO and TLBO, while consumed time by HS is slightly lower than the BBO. Convergence curves also show better convergence of the BBO and TLBO. The optimal solutions are illustrated in Fig. 14.



Figure 14. A FEM divided into k subdomains using the BBO algorithm: (a) k=5. (b) k=10

5. CONCLUSION

Four well-known metaheuristics are applied to optimal domain decomposition of FE models with different topology and number of meshes. According to results, BBO algorithm has generally performed better in most cases. According to Fig. 15(a) BBO algorithm has better performance, when is initialized by k-means++, except in standard deviation item. However by considering all algorithms, both initializations are almost equal and the only advantage of above initialization approach is lower required time to optimize, as depicted in Fig. 15(b). After BBO, the TLBO has obtained good convergence rate, average and minimum results, also, HS algorithm has consumed little time for completing the optimization procedure. Finally, this study suggests the implementation of BBO for the domain decomposition of finite elements due to the superiority of the results obtained by this method.



Figure 15. Outperforming percent of different initialization approaches: (a) in BBO algorithm. (b) by considering all algorithm

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