

An evaluation of k-means as a local search operator in hybrid memetic group search optimization for data clustering

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Abstract

Cluster analysis is one important field in pattern recognition and machine learning, consisting in an attempt to distribute a set of data patterns into groups, considering only the inner properties of those data. One of the most popular techniques for data clustering is the K-Means algorithm, due to its simplicity and easy implementation. But K-Means is strongly dependent on the initial point of the search, what may lead to suboptima (local optima) solutions. In the past few decades, Evolutionary Algorithms (EAs), like Group Search Optimization (GSO), have been adapted to the context of cluster analysis, given their global search capabilities and flexibility to deal with hard optimization problems. However, given their stochastic nature, EAs may be slower to converge in comparison to traditional clustering models (like K-Means). In this work, three hybrid memetic approaches between K-Means and GSO are presented, named FMKGSO, MKGSO and TMKGSO, in such a way that the global search capabilities of GSO are combined with the fast local search performances of K-Means. The degree of influence of K-Means on the behavior of GSO method is evaluated by a set of experiments considering both real-world problems and synthetic data sets, using five clustering metrics to access how good and robust the proposed hybrid memetic models are.

Keywords Data clustering · Evolutionary algorithms · Group search optimization · K-means

1 Introduction

In the past few decades, the amount of daily produced data in electronic devices, such as smartphones, tablets, computers, cars, GPS, smart TVs, Internet of Things applications, and so on, has increased exponentially, in such a way that automatic and scalable computational systems are even more required. To extract useful information from the large data sets such systems are based on, it is impossible to rely in human analysis only, once the need for precise and reliable information in a short period of time has become mandatory (Naldi and Campello [2014](#page-24-0)).

Data clustering is one of the most important and primitive activities in pattern recognition, consisting in an important mechanism for exploratory data analysis. Clustering is characterized by an unsupervised attempt to categorize a set of data patterns in clusters, in such a way that observations belonging in a same cluster are more close related (according to their feature set) than observations from different clusters. In clustering, no prior knowledge about the data set at hand is required, so the clustering models take decisions about how to group the observations taking into consideration only the inner properties of such observations. Clustering algorithms have been successfully employed in many real-world applications, given their flexibility and adaptability, in fields such as: Agriculture (Rahamathunnisa et al. [2020](#page-24-0)), Data Mining (Sapkota et al. [2019](#page-25-0)), Ecology (Sreepathi et al. [2017](#page-25-0)), Image Understanding (Saraswathi and Allirani [2013](#page-25-0); Wan et al. [2017](#page-25-0); Diderot et al. [2019;](#page-23-0) Wei et al. [2019](#page-25-0)), Medicine (Li et al. [2016](#page-24-0); Bruse et al. [2017;](#page-22-0) Premalatha and Subasree [2017](#page-24-0); BEDDAD et al. [2019\)](#page-22-0), Text Mining (Liu and Xiong [2011](#page-24-0)), Web Services (Parimalam and Sundaram [2017\)](#page-24-0), Wireless

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Sensors Networks (WSN) (Misra and Kumar [2016;](#page-24-0) Dhiviya et al. [2017](#page-23-0); Masoud et al. [2019](#page-24-0)), and so on. An interesting survey on data clustering algorithms can be found in (Xu and Tian [2015](#page-25-0)).

From an optimization perspective, clustering is considered as a particular kind of NP-hard grouping problem. The most popular clustering approaches are the partitional clustering algorithms, which provide a partition of the data set into a prefixed number of clusters (an input parameter for the models). Each cluster is represented by its centroid vector, and the clustering process is driven in an effort to optimize a criterion function iteratively, by updating these cluster centroids, seeking out to improve the quality of the final solution (final partition) provided by the algorithm.

One of the most popular partitional clustering methods is the K-Means algorithm (MacQueen et al. [1967](#page-24-0)). Although traditional partitional clustering models are able to promote fast convergence speeds, these methods are known for their sensibility to the initial centroid position, what may lead to weak solutions (i.e., the model may be trapped in a local minimum point) if the algorithm starts in a poor region of the problem space. Standard partitional clustering approaches lack the mechanisms to escape from local optima points, promoting local searches on the problem space only.

Evolutionary Algorithms (EAs) have been increasingly applied to solve a great variety of difficult problems, such as hard functions optimization (He et al. [2009;](#page-23-0) Oliveira et al. [2013;](#page-24-0) Pacifico and Ludermir [2014;](#page-24-0) Jain et al. [2019](#page-23-0)), weights and architecture optimization in Neural Networks (Silva et al. [2011](#page-25-0); Idrissi et al. [2016;](#page-23-0) Darwish et al. [2020](#page-23-0)), feature weighting and selection (Ramos and Vellasco [2018](#page-24-0); Tassci et al. [2018](#page-25-0); Xu et al. [2020\)](#page-25-0), and so on, given their global search capability and their mechanisms to avoid local optima points. In EAs, a set (*population*) of candidate solutions (individuals) for the problem at hand is kept and evolved according to a generational process, seeking out the optimization of a criterion function (the fitness function). In EAs such as Genetic Algorithm (GA) (Holland [1992\)](#page-23-0), Evolutionary Programming (EP) (Fogel et al. [1966](#page-23-0); Fogel [2009\)](#page-23-0), Evolution Strategies (ES) (Rechenberg [1973](#page-25-0); Schwefel [1993\)](#page-25-0) and Differential Evolution (DE) (Storn and Price [1995](#page-25-0), [1997\)](#page-25-0), the searching scheme is driven by operators that simulate biological processes like mutation, recombination and selection. In this context, Swarm Intelligence (SI) methods are extensions of EAs which execute their search in an attempt to simulate self-organizing collective behavior of social animals, like swarming, flocking and herding (Bonabeau et al. [1999\)](#page-22-0). Examples of SI algorithms are the Ant Colony Optimization (ACO) (Dorigo et al. [1996\)](#page-23-0), Particle Swarm Optimization (PSO) (Kennedy and Eberhart [1995](#page-23-0); Kennedy et al. [2001\)](#page-23-0) and Group Search Optimization (GSO) (He et al. [2006,](#page-23-0) [2009](#page-23-0)).

Table [1](#page-2-0) presents some of the most recent and popular algorithms from Evolutionary Computing literature. Interesting texts in Evolutionary Computing can be found in (Kennedy [2006;](#page-23-0) Simon [2013](#page-25-0); Eiben and Smith [2015](#page-23-0)).

EAs have been increasingly applied in clustering applications (Hruschka et al. [2009;](#page-23-0) Inkaya et al. [2016](#page-23-0); Canuto et al. [2018;](#page-22-0) Figueiredo et al. [2019\)](#page-23-0), but, given their stochastic nature, such techniques may be too slow to converge in comparison to standard partitional clustering algorithms. It is also known that, in general, evolutionary algorithms local search mechanisms are not very much effective (Ren et al. [2014](#page-25-0)).

To overcome this drawback, Hybrid Intelligence Systems (HISs) and Memetic Algorithms (MA) combining evolutionary approaches and K-Means have been proposed in literature, such that global searches performed by EAs are complemented by local searches performed by K-Means. Some examples of hybrid evolutionary systems that use K-Means as a local searcher are found in (Ahmadyfard and Modares [2008](#page-22-0); Abdel-Kader [2010;](#page-22-0) Bhavani et al. [2011;](#page-22-0) Pacifico and Ludermir [2018](#page-24-0), [2019](#page-24-0)). But most works in data clustering literature only employ K-Means after the generational process of the selected EA is concluded, which means that the best solution found by the EA may already be trapped in a local minimum point, making the application of K-Means sub-optimal (it will only exploit the problem region that contains a local minimum). To avoid such problem, in this work three hybrid memetic algorithms are proposed, named FMKGSO, MKGSO and TMKGSO, which combine the global search capabilities of GSO with local search speed provided by K-Means, which use K-Means (in different ways) to improve the population of GSO during the generational process. GSO was proposed as a method for continuous optimization problems based on the behavior of social animals attempting to find food resources, and, in the past few years, GSO has been successfully applied in many real world applications (Chen et al. [2014b](#page-23-0); Krishnaprabha and Aloor [2014](#page-24-0); Li et al. [2015](#page-24-0); Pacifico et al. [2018\)](#page-24-0), presenting competitive optimization performances in comparison to other evolutionary algorithms, such as GA, PSO, EP and ES.

The proposed approaches are implemented in an attempt to comprehend the influence of K-Means on the behavior of GSO when dealing with data clustering task, and, for that, a testing bed consisting of 20 (twenty) real-world problems (obtained from UCI Machine Learning Repository - Asuncion and Newman [2007\)](#page-22-0) and 10 synthetic data sets is proposed. Five evaluation metrics from clustering literature are employed to validate the quality of the solutions achieved by the proposed models, and five partitional evolutionary algorithms are selected for comparison purposes.

This work is organized as follows. Firstly, some introductory concepts are introduced (Sect. 2), like K-Means (Sect. 2.1), partitional evolutionary algorithms (Sect. [2.2\)](#page-3-0) and GSO (Sect. [2.3](#page-4-0)). After that, the proposed hybrid memetic GSO approaches are explained (Sect. [3](#page-5-0)), followed by the experimental evaluation (Sect. [4](#page-7-0)). Finally, some conclusions and interesting trends for future works are presented (Sect. [5](#page-15-0)).

2 Preliminaries

The baseline models adopted in this work are presented in the following sections (Sects. 2.1, [2.2](#page-3-0) and [2.3\)](#page-4-0).

2.1 K-Means

K-Means (MacQueen et al. [1967](#page-24-0)) has been proposed as a partitional clustering algorithm for continuous data, which groups real-valued data vectors into a predefined number of clusters. Consider a partition P_C of a data set with N data patterns (each data pattern is represented by a vector $\mathbf{x}_i \in \mathbb{R}^m$, where $j = 1, 2, \dots, N$ in C clusters, where C is a required input parameter for the algorithm. Each cluster is represented by its centroid vector $\mathbf{g}_c \in \mathbb{R}^m$ (where $c = 1, 2, ..., C$.

In K-Means, clusters are formed based on a dissimilarity measure, the Euclidean distance $[Eq. (1)]$. For each iteration (until a maximum number of iterations $t_{maxkmeans}$ is reached or another stopping criterion is satisfied), a new cluster centroid vector is calculated, for each cluster, as the mean of its current data vectors (i.e., the data patterns currently assigned to the cluster). After that, the new partition is formed, and each pattern is associated to the cluster with the nearest centroid.

$$
d(\mathbf{x}_j, \mathbf{g}_c) = \sqrt{\sum_{k=1}^m (x_{jk} - g_{ck})^2}
$$
 (1)

where

$$
\mathbf{g}_c = \frac{1}{N_c} \sum_{\forall \mathbf{x}_l \in c} \mathbf{x}_l \tag{2}
$$

where N_c is the number of patterns associated to cluster c. The criterion function for K-Means is the Within-Cluster Sum of Squares, given in Eq. (3).

$$
J(P_C) = \sum_{c=1}^{C} \sum_{\forall \mathbf{x}_j \in c} d(\mathbf{x}_j, \mathbf{g}_c)
$$
 (3)

K-Means algorithm is presented in Algorithm 1.

Algorithm 1 K-Means

 $t \leftarrow 0$

Initialization: Pick C patterns randomly as the initial cluster centroids \mathbf{g}_c $(c = 1, \ldots, C)$. After that, assign each pattern x_i to its closest cluster;

Calculate the initial criterion value $J(P_C^0)$ (eq.(3)).

while $(t < t_{maxkmeans})$ do

New Centroids Determination: for each cluster c , update its centroid \mathbf{g}_c using eq.(2). New Partition Determination: for each data pattern

 \mathbf{x}_i , assign it to the cluster with the nearest centroid \mathbf{g}_c . **Calculate** the new criterion value $J(P_C^t)$ (eq.(3)). $t \leftarrow t + 1$

end while

Return the final partition $P_C^{t_{maxkmeans}}$

2.2 Evolutionary algorithms for data clustering

In this section, we explain the most commonly adopted representation schema for evolutionary algorithms, when adapted as partitional clustering methods (Chen and Ye [2004\)](#page-22-0).

Once more, consider a partition P_C of a data set with N patterns $\mathbf{x}_j \in \mathbb{R}^m$ $(j = 1, 2, ..., N)$ in C clusters. Each cluster is represented by its centroid vector $\mathbf{g}_c \in \mathbb{R}^m$ $(c = 1, 2, \ldots, C)$. Each population individual $X_i \in \mathbb{R}^n$ (where $n = m \times C$) in population G represents C cluster centroids at the same time, one for each cluster (Chen and Ye [2004](#page-22-0)). For instance, if $m = 3$ and $C = 5$, each individual will be a vector $X_i \in \mathcal{R}$, where the first three features will represent the centroid \mathbf{g}_1 , features 4-th to 6-th will represent the centroid g_2 , features 7-th to 9-th will represent the centroid g_4 , features 10-th to 12-th will represent the centroid \mathbf{g}_4 , and the last three features (features 13-th to 15-th) will represent the centroid g_5 , just like in Fig. 1.

The population of evolutionary algorithms is generally initialized by a random process, but in the context of partitional data clustering, an initialization by the random choice of C patterns from the data set in analysis to compose the initial cluster centroids (just like in K-Means - see Sect. [2.1](#page-2-0)), for each individual, leads to a faster exploration of the problem search space.

As the fitness function, many works adopt the Within-Cluster Sum of Squares [Eq. ([3\)](#page-2-0)] or some alternative

function that takes such criterion as its main component, just like in (Chen and Ye [2004](#page-22-0); Ahmadyfard and Modares [2008](#page-22-0); Hruschka et al. [2009;](#page-23-0) Prabha and Visalakshi [2014](#page-24-0); Pacifico and Ludermir [2019,](#page-24-0) [2019b\)](#page-24-0). But sometimes, different functions are adopted as the fitness function (Das et al. [2007](#page-23-0); Liu et al. [2011;](#page-24-0) Wong et al. [2011](#page-25-0); He and Tan [2012](#page-23-0); José-García and Gómez-Flores [2016;](#page-23-0) Pacifico and Ludermir [2016](#page-24-0)).

Once the initial population is obtained and the fitness value for each individual **X** $_{i}^{(0)}$ in population G is computed, the evolutionary operators for the selected evolutionary algorithm are applied to the population, evolving the cluster centroids represented by each individual through a generational process, until a termination condition is reached. The global best individual found by the EA is provided as the clustering solution. A generic evolutionary algorithm for partitional data clustering is presented in Algorithm 2.

Algorithm 2 Generic Partitional Evolutionary Algorithm

$t \leftarrow 0$.
Initialize each individual $\mathbf{X}_{i}^{(0)} \in G^{(0)}$ by randomly pick-
ing C patterns from the current data set as its initial cluster
centroids.
Generate the initial partition $\mathbf{X}_{i}^{(0)} \cdot P_C^{(0)}$, assigning each
pattern \mathbf{x}_j to its closest cluster, for each individual $\mathbf{X}_i^{(0)}$.
Calculate the fitness function for each individual $\mathbf{X}_{i}^{(0)}$.
while (termination conditions are not met) do
Execute all evolutionary operators, according to the se-
lected evolutionary algorithm, on current population Gt .
Assign each pattern x_j to its closest cluster in X_i^t . P_C^t ,
for each X_i^t .
Calculate the new fitness value for each population in-
dividual $\mathbf{X}_i^t \in G^t$.
$t \leftarrow t + 1.$
end while
Return $\mathbf{X}_{best}^{t_{max}}$.

Although EAs are able to find global solutions even when dealing with complex optimization problems, the searching process performed by such strategies may be to slow, and in many applications (just like in Cui et al. [2005](#page-23-0); Abdel-Kader [2010](#page-22-0); Ahmadi et al. [2010;](#page-22-0) Chen et al. [2014a](#page-22-0); Pacifico and Ludermir [2018](#page-24-0), [2019\)](#page-24-0), the EAs are hybridized with K-Means in such a way that the EAs are used to find a

better set of starting points (i.e., cluster centroids) to K-Means. Algorithm 3 illustrates such process.

Algorithm 3 Hybrid Generic EA-k-means Algorithm
$$
t \leftarrow 0
$$
.

Initialize each individual $\mathbf{X}_{i}^{(0)} \in G^{(0)}$ by randomly picking C patterns from the current data set as its initial cluster centroids.

Generate the initial partition $\mathbf{X}_{i}^{(0)}$, $P_{C}^{(0)}$, assigning each pattern \mathbf{x}_j to its closest cluster, for each individual $\mathbf{X}_i^{(0)}$.

Calculate the fitness function for each individual $\mathbf{X}_{i}^{(0)}$. while (termination conditions are not met) do

Execute all evolutionary operators, according to the selected evolutionary algorithm, on current population G^t . **Assign** each pattern x_j to its closest cluster in $X_i^t.P_C^t$, for each X^t .

Calculate the new fitness value for each population individual $\mathbf{X}_i^t \in G^t$.

 $t \leftarrow t + 1.$ end while

Use the best individual found by the generational process $\mathbf{X}_{best}^{t_{max}}$ as the input set o cluster centroids to K-Means algorithm.

Execute K-Means on $\mathbf{X}_{best}^{t_{max}}$ for $t_{maxkmeans}$ iterations. Return the best set of cluster centroids found by K-Means \mathbf{X}_{best}^{new}

2.3 Group search optimization

Group search optimization is inspired by animal social searching behavior and group living theory. GSO employs the Producer-Scrounger (PS) model as a framework. The PS model was firstly proposed by Barnard and Sibly ([1981\)](#page-22-0) to analyze social foraging strategies of group living animals. PS model assumes that there are two foraging strategies within groups: producing (e.g., searching for food); and joining (scrounging, e.g., joining resources uncovered by others). Foragers are assumed to use producing or joining strategies exclusively. Under this framework, concepts of resource searching from animal visual scanning mechanism are used to design optimum searching strategies in GSO algorithm (He et al. [2009](#page-23-0)).

In GSO, the population G of S individuals is called group, and each individual is called a member. In a ndimensional search space, the i -th member at the t -th searching iteration (*generation*) has a current position \mathbf{X}_{i}^{t} $\in \mathbb{R}^n$ and a head angle $\alpha_i^t \in \mathbb{R}^{n-}$. The search direction of the *i*-th member, which is a vector **D** ${}_{i}^{t}(\alpha_{i}^{t}) = (d_{i1}^{t}, \ldots, d_{in}^{t})$ can be calculated from α_i^t via a polar to Cartesian coordinate transformation:

$$
d'_{i1} = \prod_{q=1}^{n-1} \cos(\alpha_{iq}^t),
$$

\n
$$
d'_{ij} = \sin(\alpha_{i(j-1)}^t) \prod_{q=1}^{n-1} \cos(\alpha_{iq}^t) \quad (j = 1, ..., n-1),
$$

\n
$$
d'_{in} = \sin(\alpha_{i(n-1)}^t)
$$
\n(4)

A group in GSO consists of three types of members: producers, scroungers and dispersed members (or rangers) (He et al. [2009\)](#page-23-0). The rangers are introduced by GSO model, extending standard PS framework.

During each GSO search iteration, a group member which has found the best fitness value so far (most promising area form the problem search space) is chosen as the producer (X_p) (Couzin et al. [2005](#page-23-0)), and the remaining members are scroungers or rangers. Standard GSO admits only one producer in each iteration, but there are some GSO variants that use multiple producers at the same time (Junaed et al. [2013](#page-23-0); Pacifico and Ludermir [2013\)](#page-24-0).

The producer employs a scanning strategy (producing) based on its vision field, generalized to a n-dimensional space, which is characterized by maximum pursuit angle $\theta_{max} \in \mathbb{R}^{n-}$ and maximum pursuit distance $l_{max} \in \mathbb{R}$, given by Eq. (5) .

$$
l_{max} = \|\mathbf{U} - \mathbf{L}\| = \sqrt{\sum_{k=1}^{n} (U_k - L_k)^2}
$$
 (5)

where U_k and L_k denote the upper bound and lower bound of the k-th dimension from the problem space, respectively.

In GSO, at the *t*-th iteration the producer X_p^t will scan laterally by randomly sampling three points in the scanning field: one at zero degree [Eq. (6)], one in the right hand side hypercube [Eq. (7)] and one in the left hand side hypercube [Eq. (8)].

$$
\mathbf{X}_z = \mathbf{X}_p^t + r_1 l_{max} \mathbf{D}_p^t (\alpha_p^t)
$$
\n(6)

$$
\mathbf{X}_r = \mathbf{X}_p^t + r_1 l_{max} \mathbf{D}_p^t (\alpha_p^t + \frac{\mathbf{r}_2 \theta_{max}}{2})
$$
 (7)

$$
\mathbf{X}_{l} = \mathbf{X}_{p}^{t} + r_{1}l_{max}\mathbf{D}_{p}^{t}(\alpha_{p}^{t} - \frac{\mathbf{r}_{2}\theta_{max}}{2})
$$
\n(8)

where $r_1 \in \Re$ is a normally distributed random number (mean 0 and standard deviation 1) and $\mathbf{r} \cdot z \in \mathbb{R}^{n-1}$ is a uniformly distributed random sequence in the range (0, 1).

If the producer is able to find a better resource than its current position, it will fly to this point; if no better point is found, the producer will stay in its current position, then it will turn its head to a new generated angle [Eq. (9)].

$$
\alpha_p^{t+1} = \alpha_p^t + \mathbf{r}_2 \alpha_{max} \tag{9}
$$

where $\alpha_{max} \in \Re$ is the maximum turning angle.

If after $a \in \mathcal{R}$ iterations the producer cannot find a better area, it will turn its head back to zero degree [Eq. (10)].

$$
\alpha_p^{k+a} = \alpha_p^k \tag{10}
$$

All scroungers will join the resource found by the producer, performing scrounging strategy according to Eq. (11).

$$
\mathbf{X}_{i}^{t+1} = \mathbf{X}_{i}^{t} + \mathbf{r}_{3} \circ (\mathbf{X}_{p}^{t} - \mathbf{X}_{i}^{t})
$$
\n(11)

where $\mathbf{r}_3 \in \mathbb{R}^n$ is a uniform random sequence in the range $(0, 1)$ and \circ is the Hadamard product or the Schur product, which calculates the entrywise product of two vectors.

The rangers will perform random walks through the problem space (Higgins and Strauss [2004](#page-23-0)), according to Eq. (12).

$$
\mathbf{X}_{i}^{t+1} = \mathbf{X}_{i}^{t} + l_{i} \mathbf{D}_{i}^{t} (\alpha_{i}^{t+1})
$$
\n(12)

where

$$
l_i = ar_1 l_{max} \tag{13}
$$

In GSO, when a member escapes from the search space bounds, it will turn back to its previous position inside the search space (Dixon [1959\)](#page-23-0). Some studies considering alternative treatments to deal with out-bounded population individuals can be found in (Xu and Shu [2006](#page-25-0); Silva et al. [2011;](#page-25-0) Pacifico et al. [2018](#page-24-0)).

GSO algorithm is presented in Algorithm 4.

Algorithm 4 GSO

 $t \leftarrow 0.$ **Initialize** randomly position $\mathbf{X}_{i}^{(0)}$ and head angles $\alpha_{i}^{(0)}$ of all members $\mathbf{X}_i^{(0)} \in G$.

Calculate the fitness value $(fit(\mathbf{X}_{i}^{(0)}))$ for each member $\mathbf{X}^{(0)}$

while (termination conditions are not met) do

Pick the best group member as the X_n^t for the current iteration.

Execute producing (\mathbf{X}_p^t) only) by evaluating three random points in its visual scanning field, \mathbf{X}_{τ}^{t} (eq. (6)), \mathbf{X}_{τ}^{t} (eq. (7)) and X_l^t (eq. (8)).

Choose a percentage from the members (but the X_p^t) to perform scrounging (eq. (11)).

Ranging: The remaining members will perform ranging through random walks (eq. (12)).

Calculate the new fitness value $fit(X_i^t)$ for each group member \mathbf{X}_i^t .

GSO scrounging operator focuses the search performed by the group in the most promising areas from the problem space, corresponding to the main exploration/exploitation strategy employed by many EAs (like crossover strategy in Genetic Algorithms and particle movement in Particle Swarm Optimization).

Producing and ranging are the main mechanisms employed by GSO for escaping local minima points. When the producer of one generation is trapped in a local minimum point (situation in which all scroungers would be trapped by following the producer, resulting in a premature convergence of the group), producing operator may find a better point in the search space, escaping from that local minimum. Even if that situation does not happen so easily, rangers will keep performing random walks that do not depend on the results found by the producer, which may lead to most promising areas than the ones found so far by the whole group, evading local minima points.

Some works have already adapted GSO to the context of partitional clustering, just like in (Pacifico and Ludermir [2014a,](#page-24-0) [2016](#page-24-0), [2019b](#page-24-0)).

3 Proposed approaches

This section presents the proposed hybrid memetic GSO and K-Means models, elaborated to test the influence of K-Means when employed as a local searcher to improve the capabilities of partitional Group Search Optimizer. Three memetic algorithms are presented: FMKGSO, MKGSO and TMKGSO. In this work, we opt to combine GSO and K-Means due to some advantages promoted by these algorithms, such as:

- 1. GSO has been proven to be a good global optimizer in many real-world applications;
- 2. GSO presents good mechanisms to escape from local optima points, represented by the producing behavior of producers, and by the ranging behavior from dispersed members;
- 3. Ranging and producing behaviors are also employed to prevent the premature convergence of the group, that is a known and hard to deal problem that affects some global search natural-inspired metaheuristics, such as Particle Swarm Optimization;
- 4. Standard GSO implements an effective mechanism to prevent out-bounded members, avoiding individuals that would not codify an actual solution to the problem at hand, by extrapolating the problem search space boundaries;
- 5. K-Means is a good local searcher, performing fast exploitation on problem space regions.

The proposed models are described as follows.

Firstly, as in generic partitional framework for EAs (see Sect. [2.2\)](#page-3-0), for all three memetic algorithms, each group member is represented as a continuous vector $X_i \in \mathbb{R}^n$ (with $n = m \times C$), codifying C cluster centroids at the same time, and at the *t*-th generation, \mathbf{X}_i will determine its own partition \mathbf{X}_{i}^{t} . P_C .

 $t \leftarrow t + 1.$

end while

Return $\mathbf{X}_{p}^{t_{max}}$.

For each algorithm, the initial population will be obtained by the random selection of C patterns from the data set being analyzed to compose each member $\mathbf{X}^{(0)}_i$. After random initialization, each member $\mathbf{X}^{(0)}_i$ has its fitness value computed to determine the quality of the partition **X** $^{(0)}_i$. P_C it represents, by associating each data pattern **x** *j* to its closest centroid vector **X** $_i^{(0)}$ **.g** *c* in **X** $_i^{(0)}$ **.**

After that, GSO generational process is initialized. Each memetic algorithm will employ K-Means for some degree during its generational process to improve the quality of the group.

As in standard GSO, for all three memetic approaches, the generational process starts, in each generation t , by the choice of the best member found so far (according to the selected fitness function) as the producer \mathbf{X}_{p}^{t} . The producer will perform *producing* operator as in standard GSO, by the evaluation of three random points from its visual scanning field [see Eqs. (6) (6) , (7) (7) and (8) (8)]. If a better point is found, the X_p^t will migrate to that point in the next generation.

The scrounging and ranging for the memetic approaches will be executed following Eqs. (11) (11) and (12) (12) , respectively. After executing all standard GSO operators, each memetic model will adopt K-Means to improve GSO group by different ways, as described bellow.

In FMKGSO, the algorithm keeps track on the quality of each member separately. For each new member X_i^{t+1} in the t-th generation (after the execution of producing, scrounging and ranging operators), determine its partition X_i^{t+1} . P_C by associating each data pattern x_j to its closest centroid vector **X** $_i^{t+1}$.**g** $_c$. Compute **X** $_i^{t+1}$ fitness. If **X** $_i^{t+1}$ has not been able to improve (in terms of the fitness value) by a predefined number of consecutive generations $(t_{maxkmeans})$, its cluster centroids are refined by the application of $t_{maxkmeans}$ K-Means steps. FMKGSO will only apply K-Means to the members that are unable to improve their partition for a period of time. This combination seeks out to speed up the search performed by each member, once GSO operators could present slow convergence rates, just like in all EAs. K-Means execution can speed up considerably the convergence rates of GSO members, promoting a better exploration and exploitation of the problem search space. The generational process executes until a termination condition is met, and the final best improved solution $\mathbf{X}_{p}^{t_{max}}$ is returned. FMKGSO algorithm is presented in Algorithm 5.

Algorithm 5 FMKGSO

 $t \leftarrow 0$

Initialization: For each member $\mathbf{X}_i^{(0)}$, pick C patterns randomly as the initial cluster centroids \mathbf{g}_c . After that, assign each pattern x_i to its closest cluster.

Calculate the initial fitness function for each member $\mathbf{X}^{(0)}$.

- while (termination conditions are not met) do
- **Pick** the best group member as the X_n^t for the current iteration.

Execute *producing* (\mathbf{X}_p^t only). For each evaluated point $(\mathbf{X}_{z}^{t}, \mathbf{X}_{r}^{t}$ and \mathbf{X}_{l}^{t}), determine its partition by assigning each data pattern to the cluster with the nearest centroid.

Choose a percentage from the members (but the X_n^t) to perform scrounging.

Ranging: The remaining members will perform *ranging* through random walks.

Determine the new partitions represented by each member X_i^{t+1} , by assigning each data pattern to the cluster with the nearest centroid, for each member.

Calculate the new fitness value for each group member. for each X_i^{t+1} do

if X_i^{t+1} has not improved for the last $t_{maxkmeans}$ generations then

Execute K-Means for $t_{maxkmeans}$ steps to refine \mathbf{X}^{t+1}

end if end for $t := t + 1$

end while

Return improved $\mathbf{X}_{n}^{t_{max}}$.

In MKGSO, after the determination of the new group by the execution of producing, scrounging and ranging operators, each member X_i^{t+1} will be improved by the execution of exactly one K-Means step in each generation before its fitness evaluation. All members are refined that way, independently if GSO operators have been able to improve or not. This operation seeks out to speedup the exploitation performed by each GSO member, by aggregating local information into GSO generational process smoothly. After that, the partitions represented by each member will be determined, and their fitness value will be computed. The generational process executes until a termination condition is met, and the final best improved solution $X_{p}^{t_{max}}$ is returned. MKGSO algorithm is presented in Algorithm 6.

Algorithm 6 MKGSO

$t \leftarrow 0$

Initialization: For each member $\mathbf{X}_i^{(0)}$, pick C patterns $\mathbf I$ randomly as the initial cluster centroids \mathbf{g}_c . After that, \mathbf{r} assign each pattern x_i to its closest cluster. έ Calculate the initial fitness function for each member ϵ $\mathbf{X}_{i}^{(0)}$. þ while (termination conditions are not met) do Ă **Pick** the best group member as the X_n^t for the current iteration. **Execute** *producing* (\mathbf{X}_p^t only). For each evaluated point $(\mathbf{X}_{z}^{t}, \mathbf{X}_{r}^{t}$ and \mathbf{X}_{l}^{t} , determine its partition by assigning each data pattern to the cluster with the nearest centroid. **Choose** a percentage from the members (but the X_n^t) to perform scrounging. **Ranging:** The remaining members will perform ranging through random walks. Determine the new partitions represented by each member \mathbf{X}_{i}^{t} , by assigning each data pattern to the cluster with the nearest centroid, for each member. **Execute** one K-Means step to refine each group member. Calculate the new fitness value for each group member. $t := t + 1$ end while **Return** improved $\mathbf{X}_{p}^{t_{max}}$ ϵ $\mathbf F$

The last proposed approach, TMKGSO, will apply K-Means to each member after a period, for a limited number of steps. After a prefixed number of generations $(t_{maxkmeans},$ an input parameter for the algorithm), each new group member X_i^{t+1} (obtained by the application of *producing*, scrounging or ranging operator) is refined by the application of $t_{maxkmeans}$ K-Means steps. By alternating some executions of GSO and K-Means, the new groups are improved by the interchanging of global and local information, in such a way that when GSO is executing, its operators will prevent the group from getting trapped in local optma regions, promoting the global exploration of the problem space, while during K-Means steps, the group will be guided through fast local exploitations on the region of the problem space where each member is placed. Each algorithm will complement the search performed by the other. The generational process executes until a termination condition is met, and the final best improved solution $X \frac{t_{max}}{p}$ is returned. TMKGSO algorithm is presented in Algorithm 7.

Algorithm 7 TMKGSO

\overline{t}

By employing three different strategies to use K-Means during GSO generational process, we can evaluate the actual influence of that technique as a local searcher. In that sense, in this work the total number of intended clusters C is supposed to be known a priori (Niu et al. [2017](#page-24-0)), so no other factors would influence on the behavior of the proposed models. The automatic determination of the optimal number of cluster for each problem is ongoing research, and that issue will be addressed in a future work. For more information on automatic clustering issue, please refer to Das et al. ([2007\)](#page-23-0), José-García and Gómez-Flores [\(2016](#page-23-0)), Elaziz et al. ([2019\)](#page-23-0).

4 Experimental results

In this section, we test the clustering capabilities of the proposed approaches, in comparison to other partitional evolutionary algorithms from literature, by means of 20 (twenty) real-world and 10 synthetic data sets. All realworld data sets are benchmark classification and clustering problems acquired from UCI Machine Learning Repository (Asuncion and Newman [2007](#page-22-0)). The selected real data set features are shown in Table [2](#page-8-0), presenting different degrees of difficulties, such as unbalanced and overlapping classes, different number of classes and features, and so on.

The proposed synthetic data sets are split into to main groups: Disp group (containing five data sets), where all classes varies in relation to its degree of dispersion, from a well-separated scenario (Disp01) to a scenario where all classes present some degree of overlapping; Prox group (containing five data sets), where the overlapping is obtained by the approximation of class centers, from a scenario where all class centers are well-split into the problem search space (Prox01), to a scenario where all class centers are close to each other (Prox05). The configurations for the proposed synthetic data sets are presented in Table [3](#page-9-0) and illustrated in Fig. [2](#page-9-0).

For comparison purposes, five clustering measures are employed: the Within-Cluster Sum of Squares [Eq. ([3](#page-2-0))], the Intra-Cluster Distance $[D_{max}, Eq. (14)]$ and the Inter-Cluster Separation $[D_{min}, Eq. (15)]$ (Wong et al. [2011](#page-25-0)), the Weighted Quantization Error $(J_e, Eq. 16)$ (Esmin et al. [2008\)](#page-23-0) and the Corrected Rand Index [CR, Eq. (17)] (Hubert and Arabie [1985](#page-23-0)).

$$
D_{max}(P_C) = \max_{c=1,\dots,C} \left\{ \sum_{\forall \mathbf{x}_j \in C} d(\mathbf{x}_j, \mathbf{g}_c) / |N_c| \right\}
$$
(14)

$$
D_{min}(P_C) = \min_{\forall c_1, c_2, c_1 \neq c_2} \{d(\mathbf{g}_{c_1}, \mathbf{g}_{c_2})\}
$$
(15)

$$
J_{e_2}(P_C) = \sum_{c=1}^{C} \left[\left(\sum_{\forall \mathbf{x}_j \in C} d(\mathbf{p}_j, \mathbf{g}_c) / |N_c| \right) \times \left(|N_c| / N \right) \right] \tag{16}
$$

The CR assesses the degree of similarity between an a priori partition and a partition provided by the clustering algorithm. Given that all data sets adopted in this work are real-valued classification problems (i.e., all data patters are labeled in each data set), CR represents a robust comparison metric for clustering studies, since its analysis takes into consideration only the relationships among the data patterns from an *a priori* and an *a posteriori* partitions, without taking into consideration the categories themselves.

Considering a partition $U_R = \{u_1, u_2, \ldots, u_R\}$ provided by a clustering algorithm and an a priori partition defined by classification $V_C = \{v_1, v_2, \ldots, v_C\}$, CR is defined as by Eq. (17).

$$
CR = \frac{a - b}{c - d} \tag{17}
$$

where

$$
a = \sum_{i=1}^{R} \sum_{j=1}^{C} \binom{n_{ij}}{2}
$$
 (18)

$$
b = \binom{n}{2}^{-1} \sum_{i=1}^{R} \binom{n_i}{2} \tag{19}
$$

$$
c = 1/2 \left[\sum_{i=1}^{R} {n_i \choose 2} + \sum_{j=1}^{C} {n_j \choose 2} \right]
$$
 (20)

Data set	Classes	Instances per class	Total no. of instances	Attributes	μ_k (1 to 5)	Σ_k (1 to 5)
Disp01	5	200	1000	3	$(-10, -5, 0, 5, 10)$	(1, 1, 1, 1, 1)
Disp02	5	200	1000	3	$(-10, -5, 0, 5, 10)$	(4, 1, 1, 1, 4)
Disp03	5	200	1000	3	$(-10, -5, 0, 5, 10)$	(4, 1, 4, 1, 4)
Disp04	5	200	1000	3	$(-10, -5, 0, 5, 10)$	(4, 4, 1, 4, 4)
Disp05	5	200	1000	3	$(-10, -5, 0, 5, 10)$	(4, 4, 4, 4, 4)
Prox01	5	200	1000	3	$(-8, -4, 0, 4, 8)$	(1, 1, 1, 1, 1)
Prox02	5	200	1000	3	$(-8, -6, 0, 4, 8)$	(1, 1, 1, 1, 1)
Prox03	5	200	1000	3	$(-8, -6, 0, 6, 8)$	(1, 1, 1, 1, 1)
Prox04	5	200	1000	3	$(-8, -6, -4, 6, 8)$	(1, 1, 1, 1, 1)
Prox05	5	200	1000	3	$(-8, -6, -4, -2, 0)$	(1, 1, 1, 1, 1)

Table 3 Disp and Prox synthetic data sets description: each Disp and Prox data set consists of 1000 3-dimensional patterns, equally distributed among 5 classes

The mean μ and covariance matrix Σ are the same for each attribute in each class, but varies from class to class (the values are presented from class 1–5, respectively)

Fig. 2 Synthetic data sets representation: a–e Disp01 to Disp05, respectively, f–j Prox01-Prox05, respectively

$$
d = \binom{n}{2}^{-1} \sum_{i=1}^{R} \binom{n_i}{2} \sum_{j=1}^{C} \binom{n_j}{2} \tag{21}
$$

where n_{ij} represents the number of objects that are in clusters u_i and v_i ; n_i indicates the number of objects in cluster u_i ; n_j indicates the number of objects in cluster v_i ; and n is the total number of objects. CR takes its values from the interval $[-1,1]$, in which the value 1 indicates perfect agreement between partitions, whereas values near 0 (or negatives) correspond to cluster agreement found by chance (Arabie et al. [1996\)](#page-22-0).

The evaluation criterion includes a rank system employed through the application of Friedman hypothesis test (Friedman [1937,](#page-23-0) [1940\)](#page-23-0) for all the comparison clustering measures. The Friedman test is a non-parametric hypothesis test that ranks all algorithms for each data set separately. If the null-hypothesis (all ranks are not significantly different) is rejected, Nemenyi test (Nemenyi [1962](#page-24-0)) is adopted as the post-hoc test. According to Nemenyi test, the performance of two algorithms are considered significantly different if the corresponding average ranks differ by at least the critical difference

$$
CD = q_a \sqrt{\frac{n_{alg}(n_{alg} + 1)}{6n_{data}}}
$$
 (22)

where n_{data} represents the number of data sets, n_{alg} represents the number of compared algorithms and q_a are critical values based on a Studentized range statistic divided by $\sqrt{2}$ (Demšar [2006](#page-23-0)). Since J, J_{e_2} and D_{max} are *minimization metrics* (indicated by \downarrow), the best methods will obtain lower ranks for the Friedman-Nemenyi test, while for D_{min} and CR (maximization metrics, indicated by \uparrow), the best methods will find higher average ranks for the Friedman-Nemenyi hypothesis test.

The proposed hybrid memetic GSO approaches are compared to five other EAs: Genetic Algorithm (GA), Differential Evolution (DE), Particle Swarm Optimization (PSO), standard Group Search Optimization and Backtracking Search Optimization (BSA). The selected approaches are state-of-the-art models from evolutionary computing and data clustering literature, being successfully applied in many applications recently: GA (Shi and Xu [2018;](#page-25-0) Akbari et al. [2019](#page-22-0); Islam et al. [2019](#page-23-0); Mortezanezhad and Daneshifar [2019](#page-24-0); Toman et al. [2020\)](#page-25-0), DE (Wang [2018](#page-25-0); Li and Dong [2019;](#page-24-0) Cho and Nyunt [2020;](#page-23-0) Zhang and Cao [2020;](#page-25-0) Zhu et al. [2020](#page-25-0)), PSO (Souza et al. [2018](#page-25-0); Wang et al. [2018;](#page-25-0) Li et al. [2019](#page-24-0); Pacifico and Ludermir [2019](#page-24-0); Miranda and Prudêncio [208\)](#page-24-0), GSO (Pacifico and Ludermir [2018;](#page-24-0) Lin and Huang [2019](#page-24-0); Pacifico and Ludermir [2019b](#page-24-0); Taj and Basu [2019;](#page-25-0) Abualigah [2020](#page-22-0)), and BSA (Latiff et al. [2016](#page-24-0); Günen et al. [2017;](#page-23-0) Li et al. [2019](#page-24-0); Toz et al. [2019;](#page-25-0) Hassan and Rashid [2020](#page-23-0)).

All selected algorithms have been adapted as hybrid partitional clustering models according to the schema presented in Algorithm 3. In this work, all algorithms used the Within-Cluster Sum of Squares $[Eq. (3)]$ $[Eq. (3)]$ $[Eq. (3)]$ as the fitness function, for simplicity, once our main objective is to evaluate the impact of K-Means on the behavior of memetic GSO, and not the influence of the fitness function on the behavior of EAs. For that issue, please, refer to Wong et al. ([2011\)](#page-25-0), Pacifico and Ludermir ([2016\)](#page-24-0).

All algorithms run in a MATLAB 7.6 environment. Thirty independent tests have been executed for each data set, and all evolutionary methods have started with the same initial population in each test, obtained by a random process, as explained in Sect. [2](#page-2-0)), as a manner to guarantee a fair evaluation and comparison among the selected techniques. Also, once the computational costs have been evaluated in terms of the average execution time of each model in each data set, each algorithm has been run and tested in a computer with an i7-7700K CPU, NVIDIA GeForce GTX 1060 6GB GPU and 32 GB RAM, independently (one algorithm each time), and no other programs, but the Operating System, were executed during the tests, granting the same environmental conditions to each method. For all tests, the adopted number of clusters C is equal to the number of classes per data set.

The hyperparameters for each evolutionary algorithm are presented in Table 4, and have been obtained from the literature: GA (Abdel-Kader [2010\)](#page-22-0), DE (Das et al. [2007](#page-23-0)), PSO (Abdel-Kader [2010](#page-22-0); Pacifico and Ludermir [2019](#page-24-0)), GSO (He et al. [2009;](#page-23-0) Pacifico and Ludermir [2014a](#page-24-0), [2018](#page-24-0)), and BSA (Civicioglu [2013\)](#page-23-0). The population size (S) and the maximum number of generations (t_{max}) have been chosen after a trial-and-error process, using values from Pacifico and Ludermir ([2018\)](#page-24-0),Pacifico and Ludermir [\(2019](#page-24-0)) as the staring point. Once the selected approaches

did not improve significantly with higher population sizes nor higher number of generations, we kept the same parameter values as presented in Pacifico and Ludermir [\(2018](#page-24-0)), Pacifico and Ludermir ([2019\)](#page-24-0) to perform the current evaluation, once the computational costs associated with such parameters may grow very fast for high parameter values. The maximum number of K-Means executions $(t_{maxkmeans})$ has been obtained from Pacifico and Ludermir [\(2018](#page-24-0)).

4.1 Discussion

The discussion is divided into three parts: first, we evaluate the experimental results for the real-world data sets (Sect. [4.1.1](#page-11-0)), followed by the evaluation on the experiments using the synthetic data sets (Sect. [4.1.2](#page-11-0)); An overall evaluation is also performed, based on the Friedman-Nemenyi hypothesis tests and the computational costs (in terms of the average execution times) obtained by each model (Sect. [4.1.3](#page-11-0)).

4.1.1 Real-world data sets

The experimental results for the real-world data sets are presented from Tables [5,](#page-12-0) [6](#page-13-0), [7](#page-14-0), [8](#page-15-0), [9.](#page-16-0) The best results for each metric in each data set are bold faced.

In an empirical analysis, we can observe that the proposed hybrid memetic GSO approaches are able to find better values in relation to the fitness function (J) in most cases than hybrid partitional models that only use K-Means to refine the best solution found by the generational process of the EA. As illustrated in Fig. [3,](#page-17-0) sometimes the comparison hybrid partitional approaches are too slow while performing the global search that the only significant improvement is obtained after the generational process, when their final solution is refined by K-Means method (which presents fast local search capabilities). Fig. [3](#page-17-0) also showed that even the use of just one K-Means step during the generational process for each member (MKGSO approach) is enough to improve significantly the convergence speed of GSO. The empirical analysis also showed that the hybrid memetic GSO models are more stable than the comparison hybrid partitional EA and K-Means algorithms.

The overall evaluation on the real-world problems obtained by the Friedman-Nemenyi hypothesis tests (Table [10\)](#page-17-0) showed that the proposed memetic approaches are able to find clusters that are, in average, better represented by their final centroids (according to J and J_{e_2}) indices), the final clusters are more similar to the actual classes (CR index), and the clusters are well-split in the problem space (according to D_{min}). The hypothesis tests also showed that, although MKGSO and TMKGSO use about the same number of K-Means steps for each member on current experimentation, MK-GSO presented more smooth convergence rates than TMKGSO, what may help MKGSO to avoid local minima points, but yet according to the current set of experiments, the performance of TMKGSO has been considered significantly better than the average results for MKGSO (in terms of the optimization of the fitness function $-J$).

According to the Friedman-Nemenyi tests, considering all five clustering metrics, the best results have been found by TMKGSO, followed by MKGSO and FMKGSO (second and third places, respectively).

When considering the average execution time obtained by each method in each data set (Table [11\)](#page-18-0), we can see that FMKGSO and TMKGSO (which use K-Means refinements only in determined conditions) presented average times compatible with the other state-of-the-art EAs from the literature. By the other hand, MKGSO presented average execution times about four times higher than the other evaluated approaches, what may limit its applications when

the data set is big (in terms of the number of data patterns). Such limitation may have occurred due to the fact that such method uses K-Means calls in each generation for each individual of the population, requiring many calls for allocation and deallocation of computational resources by the compiler.

4.1.2 Synthetic data sets

The results for both sets of synthetic data sets are presented from Tables [12,](#page-18-0) [13](#page-19-0), [14,](#page-19-0) [15,](#page-20-0) [16](#page-20-0). The best results for each metric in each data set are bold faced.

The empirical analysis considering both Disp and Prox sets showed that all algorithms have been significantly affected when the degree of class overlapping increased, specially when the class overlapping has been obtained by approximating the class centers of distribution (Prox data sets). In spite of that, the memetic approaches have been able, once again, to find the best average performances and degree of stability among all approaches. In particular, FMKGSO was able to find the best values for almost all tests, considering the five evaluation metrics, showing its capabilities and competitiveness. Fig. [4](#page-21-0) illustrates the obtained partitions for each algorithm.

Considering the ranks provided by Friedman-Nemenyi hypothesis test (Table [17](#page-21-0)), the memetic approaches have found the best values for all evaluated indices. Although, in an overall analysis on the synthetic data sets, FMKGSO has been able to present ranks slightly better then the ranks of MKGSO, the Friedman-Nemenyi tests showed that there is no statistical significantly differences between FMKGSO performance and the performances of MKGSO (their differences are not greater than the Critical Distance), so both memetic algorithms are considered to have achieved the best results together. TMKGSO has been capable of reaching the third place on the evaluation considering the synthetic approaches, but, for all five clustering metrics, this method has found better performances than all the selected state-of-the-art comparison approaches, reinforcing the advantages of memetic models.

The evaluation considering the average execution times for each model showed that (Table [18\)](#page-22-0) showed that, once more, FMKGSO and TMKGSO computational costs, in terms of time, are close to the computational costs of stateof-the-art algorithms, but MKGSO, although presenting better performances when considering the clustering metrics than the non-memetic comparison approaches, is still a little bit slower in execution.

4.1.3 Overall evaluation

At this point, an overall evaluation considering both the real-world and synthetic data sets is provided. For that, an

evaluation based on the Friedman-Nemenyi hypothesis tests considering all data sets (real and synthetic) is per-formed (Table [19\)](#page-22-0). By considering both sets of data sets, we can have a better understanding on the behavior of all selected algorithms (hybrid partitional approaches, and hybrid memetic partitional approaches) when dealing with either controlled and uncontrolled testing scenarios, showing how robust such approaches actually are.

The overall analysis ranked TMKGSO as the best approach among all the tested algorithms, considering the rank values for all metrics, followed by FMKGSO and MKGSO, respectively. The overall evaluation showed that including some K-Means during the generational process of EAs is quite advantageous (the memetic approach), and it is a preferred strategy than only use K-Means only to refine the best solution found so far by the EAs. Considering the selected five clustering metrics, we can concluded that, for the current testing bed, the memetric approaches are able to find final solutions with clusters that are, in average, better represented by their centroids vectors (according to J and J_{e_2} indices), more similar to the actual classes (CR index), more compact (according to D_{max} index), and well-split in the problem space in relation to each other (according to D_{min}). The proposed memetic

Table 7 Experimental results for real-world data sets: weighted quantization error $(J_{e_2}^{\downarrow})$

Data set	GA-k-means	DE-k-means	PSO-k- means	GSO-k- means	BSA-k- means	FMKGSO	MKGSO	TMKGSO
Abalone	2.4191 \pm 0.071	2.3979 \pm 0.024	2.4191 \pm 0.071	2.3847 \pm 0.025	2.4076 \pm 0.020	2.3677 ± 0	2.3677 ± 0	2.3677 ± 0
Banknote authentication	32.106 ± 0	32.106 ± 0	32.106 ± 0	32.106 ± 0	32.106 ± 0	32.106 ± 0	32.106 ± 0	32.106 ± 0
Blood transfusion	$9.08\times$ 10^5 \pm 5.83×10^{3}	9.06×10^5 ± 0	$9.08\times$ 10^5 \pm 5.83×10^{3}	9.06×10^5 ± 0	$9.07\times$ 10^5 \pm 4.23×10^{3}	9.06×10^5 ± 0	$9 \cdot 06 \times 10^5$ ± 0	9.06×10^{5} ± 0
Cancer	0.3483 ± 0	0.3483 ± 0	0.3483 ± 0	0.3483 ± 0	0.3483 ± 0	0.3483 ± 0	0.3483 ± 0	0.3483 ± 0
Diabetes	6695.8 ± 0	6695.8 ± 0	6695.8 ± 0	6695.8 ± 0	6695.8 ± 0	6695.9 \pm 0.251	6695.8 ± 0	6695.8 ± 0
E. coli	$0.0449 \pm$ 0.002	$0.0446 \pm$ 0.002	$0.0447 \pm$ 0.002	$0.0441 \pm$ 0.002	$0.0445 \pm$ 0.002	$0.0412 \pm$ 0.0001	$0.0418 \pm$ 0.0009	$0.0413 \pm$ 0.0002
Glass	$1.7405 \pm$ 0.106	1.7044 \pm 0.135	$1.7405 \pm$ 0.106	1.6983 \pm 0.114	$1.7692 \pm$ 0.115	1.9592 \pm 0.204	1.7281 \pm 0.141	$1.6097 \pm$ 0.066
Heart	2034.5 \pm 0.047	2034.5 \pm 0.047	2034.5 \pm 0.047	2034.6 \pm 0.047	2034.5 \pm 0.043	2034.5 ± 0	2034.5 ± 0	2034.5 ± 0
Image segmentation	6435.9 \pm 867.5	6324.7 \pm 744.7	6435.9 \pm 867.5	6616.2 \pm 1074.9	6525.9 \pm 986.6	6567.5 \pm 455.7	6273.2 \pm 477.3	6063.9 \pm 222.2
Ionosphere	6.8928 \pm 0.0001	6.9711 \pm 0.428	6.8929 \pm 0.0001	6.8928 \pm 0.0001	6.8928 \pm 0.00006	6.8928 ± 0	6.8928 ± 0	6.8928 ± 0
Iris	$0.5257~\pm$ 0.00001	$0.5257 \pm$ 0.00001	$0.5266 \pm$ 0.0013	$0.5257 \pm$ 0.00001	$0.5256 \pm$ 0.00001	0.5257 ± 0	0.5257 ± 0	0.5257 ± 0
Landsat satellite images	2379.5 \pm 51.14	$2363.2 \pm$ 29.74	2383.4 \pm 57.65	2372.4 \pm 45.57	2396.8 \pm 65.87	2333.9 \pm 5.609	$2355.1 \pm$ 26.30	$2357.5 \pm$ 19.91
Letter recognition	31.022 \pm 0.213	31.049 \pm 0.179	$30.993 \pm$ 0.217	31.005 \pm 0.229	$31.002 \pm$ 0.200	$30.623 \pm$ 0.064	30.864 \pm 0.178	30.817 \pm 0.192
Optical recognition	666.47 \pm 9.538	665.05 \pm 13.13	666.47 \pm 9.538	667.58 \pm 11.20	$661.57 \pm$ 7.652	692.93 \pm 1.107	694.88 \pm 20.65	672.33 \pm 11.75
Page blocks classification	6.08×10^{6} ± 2.57×10^{6}	$3.72 \times 10^6 \pm$ 9.35×10^5	$4.34 \times 10^6 \pm$ 1.69×10^{6}	$4.22 \times 10^6 \pm$ 1.12×10^{6}	3.82×10^6 ± 7.16×10^5	3.69×10^{6} ± 8.95×10^{5}	$3.29 \times 10^6 \pm$ 1.02×10^{6}	2.96×10^6 \pm 8.77×10^5
Pen-digits	4619.3 \pm 77.58	4594.1 \pm 100.7	4610.2 \pm 75.39	4619.0 \pm 74.58	4617.5 \pm 98.40	4528.3 \pm 93.61	4566.0 \pm 80.69	4566.9 \pm 43.77
Seeds	2.7991 \pm 0.0034	2.7998 \pm 0.0035	2.7991 \pm 0.0034	2.7986 \pm 0.0031	2.7986 \pm 0.003	2.7968 ± 0	2.7968 ± 0	2.7968 ± 0
Waveform	26.624 ± 0	26.624 ± 0	26.624 ± 0	26.624 ± 0	26.624 ± 0	26.624 ± 0	26.624 ± 0	26.624 ± 0
Wine	13558.4 \pm 545.6	13318.5 ± 0	13558.4 \pm 545.6	13318.5 ± 0	13607.6 \pm 588.2	13318.5 ± 0	13318.5 ± 0	13318.5 ± 0
Yeast	$0.0320 \pm$ 0.0017	$0.0327 \pm$ 0.0022	$0.0322 \pm$ 0.0016	$0.0312 \pm$ 0.0003	$0.0322 \pm$ 0.0016	$0.0305 \pm$ 0.0001	$0.0306 \pm$ 0.0002	$0.0307 \pm$ 0.0003

approaches have also been able to find the best stability throughout the experimentation, reinforcing their reliability.

For the current testing bed, the best hybrid model has been the TMKGSO approach, which executes some K-Means steps after a prefixed number of generations for each group member, which either has found good performances, according to the clustering indices, and has been able to execute with a computational cost, in terms of average execution time, close to what is expected from any Evolutionary Algorithm from current state-of-the-art. By the other hand, although able to find good average performances, MKGSO approach has been quite slow in its execution, what may represent a limitation. But, if time is not the main concern when a new data clustering system is proposed, MKGSO would still be a better choice than nonmemetic approaches, once it still preserves GSO good global searching capabilities and K-Means fast convergence speed, so the designer of such system should have this trade-off in mind when making the decision.

Table 8 Experimental results for real-world data sets: intra-cluster distance (D_{max}^{\downarrow})

Data set	GA-k-means	DE-k-means	PSO-k- means	GSO-k- means	BSA-k- means	FMKGSO	MKGSO	TMKGSO
Abalone	6.8652 \pm 0.501	$6.6643 \pm$ 0.352	6.8652 \pm 0.501	6.8051 \pm 0.279	6.7650 \pm 0.484	6.7044 ± 0	6.7044 ± 0	6.7044 ± 0
Banknote authentication	35.978 ± 0	35.978 ± 0	35.978 ± 0	35.978 ± 0	35.978 ± 0	35.978 ± 0	35.978 ± 0	35.978 ± 0
Blood transfusion	$3.93\times$ 10^6 \pm 1.68×10^{5}	$3 \cdot 86 \times 10^6$ ± 0	$3.93\times$ 10^6 \pm 1.68×10^5	$3 \cdot 86 \times 10^6$ ± 0	$3.91 \times 10^6 \pm$ 1.33×10^5	$3 \cdot 86 \times 10^6$ ± 0	$3\cdot 86\times 10^6$ ± 0	$3\cdot86\times10^6$ ± 0
Cancer	0.8050 ± 0	$0.8050\,\pm\,0$	0.8050 ± 0	0.8050 ± 0	0.8050 ± 0	0.8050 ± 0	0.8050 ± 0	0.8050 ± 0
Diabetes	$1\cdot 68\times 10^4$ ± 0	$1\cdot 68\times 10^4$ ± 0	$1\cdot 68\times 10^4$ ± 0	$1\cdot 68\times 10^4$ ± 0	$1\cdot 68\times 10^4$ ± 0	$1\cdot 68\times 10^4$ ± 11.56	$1\cdot 68\times 10^4$ ± 0	$1\cdot 68\times 10^4$ ± 0
E. Coli	$0.1029 \pm$ 0.015	$0.1048 \pm$ 0.018	$0.1118 \pm$ 0.019	$0.1069 \pm$ 0.016	$0.1083 \pm$ 0.017	0.1330 ± 0	$0.1307 \pm$ 0.013	0.1330 ± 0
Glass	$9.0743 \pm$ 1.639	$9.2814 \pm$ 1.764	$9.0743 \pm$ 1.639	7.9628 \pm 2.163	$8.6093 \pm$ 1.730	8.4604 \pm 1.825	7.3165 \pm 0.912	7.4499 \pm 1.109
Heart	2718.0 \pm 2.542	2718.5 \pm 2.425	2718.0 \pm 2.542	2718.0 \pm 2.323	2718.9 \pm 2.423	2720.7 ± 0	2720.7 ± 0	2720.7 ± 0
Image segmentation	231386.2 \pm 58399.2	235047.8 ± 54517.2	231386.2 \pm 58399.2	213276.6 \pm 75049.5	222034.2 \pm 68870.5	183540.4 \pm 60964.4	206166.7 \pm 57345.6	209843.8 \pm 51638.3
Ionosphere	10.547 \pm 0.0127	10.505 \pm 0.2394	10.546 \pm 0.0126	$10.550 \pm$ 0.0127	$10.548 \pm$ 0.012	10.561 ± 0	10.561 ± 0	10.561 ± 0
Iris	$0.6488 \pm$ 0.004	$0.6485 \pm$ 0.004	$0.6488 \pm$ 0.004	$0.6504 \pm$ 0.003	$0.6497 \pm$ 0.003	0.6423 ± 0	0.6423 ± 0	0.6423 ± 0
Landsat satellite images	5769.5 \pm 1306.0	5729.4 \pm 1424.2	5738.7 \pm 1364.3	5678.1 \pm 1309.2	6343.7 \pm 1548.2	4726.1 \pm 6.984	5383.9 \pm 1105.2	5507.9 \pm 1033.7
Letter recognition	48.858 \pm 1.816	49.113 \pm 2.401	48.541 \pm 2.159	48.644 \pm 2.008	49.019 \pm 2.272	48.739 ± 0.639	49.035 ± 1.715	48.682 \pm 1.439
Optical recognition	800.70 \pm 51.17	781.49 \pm 46.10	800.70 \pm 51.17	795.23 \pm 49.98	772.15 \pm 20.07	829.63 \pm 66.73	827.20 \pm 64.67	798.93 ± 44.25
Page blocks classification	$1.20\times$ 10^9 \pm 4.42×10^{7}	6.28×10^8 \pm 3.31×10^8	$1.18\times$ 10^9 \pm $\overline{0}$	$1.16\times$ 10^9 \pm 1.42×10^8	$1.18\times$ 10^9 \pm $\mathbf{0}$	$1.01\times$ 10^9 \pm 3.27×10^8	8.30 \times 10 ⁸ \pm 3.84×10^{8}	$6.76 \times 10^8 \pm$ 3.65×10^8
Pen-digits	6534.3 \pm 427.0	6358.0 \pm 438.2	6455.9 \pm 463.5	$6560.1~\pm$ 427.8	6490.3 \pm 451.4	6411.8 \pm 262.2	6423.8 \pm 377.0	6525.4 \pm 484.0
Seeds	3.0182 ± 0	3.0182 ± 0	3.0182 ± 0	3.0182 ± 0	3.0182 ± 0	3.0182 ± 0	3.0182 ± 0	3.0182 ± 0
Waveform	27.140 ± 0	27.140 ± 0	27.140 ± 0	27.140 ± 0	27.140 ± 0	27.140 ± 0	27.140 ± 0	27.140 ± 0
Wine	27559.4 \pm 3196.2	28956.4 ± 0	27559.4 \pm 3196.2	28956.4±0	27364.0 \pm 3253.5	28956.4 ± 0	28956.4±0	28956.4 ± 0
Yeast	$0.0638 \pm$ 0.0107	$0.0598 \pm$ 0.0082	$0.0681 \pm$ 0.0161	$0.0662 \pm$ 0.0118	$0.0305 \pm$ 0.0043	$0.0931 \pm$ 0.0127	0.0980 ± 0	$0.0907 \pm$ 0.0150

5 Conclusions

This paper presented an evaluation on the behavior of partitional Group Search Optimization when employing K-Means algorithm as a local search operator, to deal with data clustering problem. In that sense, three hybrid memetic algorithms are proposed (FMKGSO, MKGSO and TMKGSO), which use different ways to combine GSO and K-Means in an attempt to improve GSO population during the generational process of the method. Although K-Means is easily trapped in local optima points, it is known that it would represent a good local searcher when combined to global search metaheuristics, like evolutionary algorithms. Also, GSO, which is known to present good global searching capabilities and mechanisms to escape from local optima points, would benefit considerably by adopting K-Means to speedup the algorithm convergence when dealing with data clustering problems.

To evaluate the proposed methods, five state-of-the-art partitional clustering evolutionary algorithms are tested:

GA, DE, PSO, standard GSO and BSA. As a current and usual practice in data clustering literature, the comparison approaches are hybridized with K-Means in such a way that their best final solutions are refined by K-Means after the generational process, so we could access the actual influence of such technique on the behavior of EAs.

The experimental testing bed adopted in this work included 20 real-world and 10 synthetic data sets. To access the potential of the proposed memetic models, five clustering metrics have been employed (within-cluster sum of squares, correct Rand index, weighted quantization error, intra-cluster distance and inter-cluster separation), and the experimental analysis has been obtained by means of both an empirical analysis, and by an overall evaluation obtained through the application of a rank system considering the Friedman-Nemenyi hypothesis tests on each clustering metric.

Fig. 3 Average convergence graph for the best solution found so far X_p^t for a Abalone, b E. coli, c Glass, d Image segmentation, e Page blocks classification, f Yeast

Table 10 Overall evaluation on real-world data sets: average Friedman–Nemenyi ranks (and position) for each metric. The best results are bold faced

The critical distance is $CD = 2.3477$

The experimental results showed that, when K-Means is executed within the generational process of GSO at least by a few steps, it is able to aggregate more local information than when it is adopted only to improve the final solution found so far by an EA, speeding up the exploration/exploitation promoted by the memetic system. EAs are known to present slow convergence speeds, so when K-Means is only employed after the generational process of the EA, given K-Means lack of recovering mechanisms from local optima points, the final solution found by the hybrid model may still be in a region of the problem space that do not contain the global optimum point, so K-Means will still be trapped in such region.

All memetic approaches have been able to improve the comparison approaches, and the best evaluated memetic combination between K-Means and GSO (according to the Friedman-Nemenyi ranks and the evaluation on the average execution times) has been achieved when K-Means executes for a limited number of steps, for each population individual, after a prefixed number of GSO generations (TMKGSO algorithm), followed by FMKGSO (where individuals are improved by K-Means only if they fail to improve for a prefixed number of consecutive generations).

Data set	GA-k-means	DE-k-means	PSO-k-means	GSO-k-means	BSA-k-means	FMKGSO	MKGSO	TMKGSO
Abalone	3.2324	3.1665	3.1848	3.7053	3.2208	3.2161	11.993	2.2860
Banknote authentication	0.6849	0.6814	0.6726	0.8054	0.6885	0.7939	2.5665	0.6990
Blood transfusion	0.3938	0.3792	0.3808	0.4633	0.3883	0.3616	1.4789	0.2945
Cancer	0.4844	0.4633	0.4654	0.5878	0.4986	0.4299	1.8070	0.4041
Diabetes	0.5166	0.4940	0.4947	0.6159	0.5155	0.5596	1.8619	0.5105
E. coli	0.5914	0.5820	0.5820	0.8138	0.7540	1.4727	3.0436	0.9399
Glass	0.3681	0.3524	0.3547	0.5540	0.5159	0.6873	1.8540	0.5448
Heart	0.2361	0.2202	0.2214	0.3206	0.2607	0.2866	0.9268	0.2626
Image segmentation	6.2777	6.3355	6.2147	8.2467	7.3594	10.891	21.707	7.0738
Ionosphere	0.4540	0.4443	0.4522	0.7103	0.6920	0.5505	1.8175	0.5118
Iris	0.1545	0.1413	0.1473	0.2002	0.1546	0.2267	0.6869	0.1886
Landsat satellite images	47.173	47.168	47.064	55.600	50.703	85.043	152.46	53.855
Letter recognition	254.40	254.61	253.51	294.70	263.82	553.67	809.80	334.92
Optical recognition	105.40	106.37	106.07	129.77	129.52	169.31	343.48	100.96
Page blocks classification	7.2430	7.8097	6.9642	8.0242	8.0102	11.203	24.978	7.1443
Pen-digits	52.908	52.925	52.971	61.555	54.205	87.005	174.22	55.801
Seeds	0.2066	0.1930	0.1937	0.2777	0.2206	0.2795	0.8892	0.2312
Waveform	9.8666	9.8409	9.8524	11.496	10.095	11.378	33.455	9.7679
Wine	0.2291	0.2132	0.2155	0.3475	0.3063	0.3398	0.9966	0.2940
Yeast	3.1202	3.0142	3.0257	3.6732	3.3673	8.0542	11.848	4.5985

Table 11 Average execution times for the real-world data sets (in seconds)

Table 12 Experimental results for synthetic data sets: within-cluster sum of squares (J^{\downarrow})

Data set	GA-k-means	DE-k-means	PSO-k-means	GSO-k-means	BSA-k-means	FMKGSO	MKGSO	TMKGSO
Disp01	$3232.8 \pm$ 1320.6	$3468.9 \pm$ 1815.9	2991.7 ± 0	2991.7 ± 0	$3241.3 \pm$ 1367.3	2991.7 ± 0	2991.7 ± 0	2991.7 ± 0
Disp02	7138.9 \pm 1831.8	$7820.9 \pm$ 2646.7	$7125.8 \pm$ 1781.2	$6905.3 \pm$ 1355.5	$7631.2 \pm$ 2525.1		6657.8 ± 0 6657.8 ± 0	6657.8 ± 0
Disp03	8756.5 \pm 1791.9	$8602.0 \pm$ 1646.6	8402.1 \pm 1275.0	8363.8 \pm 1065.8	8386.2 \pm 1187.7		8169.1 ± 0 8169.1 ± 0	8169.1 ± 0
Disp04	9809.4 \pm 1395.1	$10904.6 \pm$ 2466.6	9442.8 ± 0	$10173.1 \pm$ 1893.8	$9825.9 \pm$ 1458.2		9442.8 ± 0 9442.8 ± 0	9442.8 ± 0
Disp05	11942.3 \pm 1042.3	$12706.0 \pm$ 2175.1	11752.0 ± 0	$11937.5 \pm$ 1015.7	11923.7 \pm 940.7	$11752.0 \pm$ $\mathbf{0}$	11752.0 ± 0	11752.0 ± 0
Prox01	$3407.4 \pm$ 1186.0	$3700.9 \pm$ 1569.1	3095.8 ± 0	3095.8 ± 0	$3251.8 \pm$ 854.7	3095.8 ± 0	3095.8 ± 0	3095.8 ± 0
Prox ₀₂	3226.4 ± 526.8	3266.2 ± 545.1	3188.9 \pm 508.8	3227.1 ± 528.0	3382.9 \pm 577.1	2887.3 ± 0	2887.3 ± 0	2924.7 \pm 205.2
Prox03	3304.6 ± 594.8	3530.1 ± 544.2	3266.1 \pm 498.6	3267.0 ± 500.0	3393.3 \pm 553.7	2919.8 ± 0	$2987.5 \pm$ 257.8	2956.1 \pm 199.0
Prox04	3078.7 ± 446.8	3156.4 ± 460.2	$3060.6 \pm$ 411.5	3066.8 ± 422.3	$2920.8 \pm$ 240.9	2841.9 ± 0	$2868.0 \pm$ 143.1	2983.6 \pm 327.1
Prox05	$2725.4 \pm$ 0.2862	2773.5 ± 182.9	$2725.4 \pm$ 0.2564	$2725.3 \pm$ 0.2300	$2725.4 \pm$ 0.2785	2725.2 ± 0	2725.2 ± 0	2725.2 ± 0

It means that interchange between the global search offered by EA operators and the local search promoted by K-Means (i.e., the memetic framework) may represent a robust approach to explore and exploit the problem search space, leading to better optimization performances. By the other hand, the use of just one K-Means step for each

Data set	GA-k-means	DE-k-means	PSO-k-means	GSO-k-means	BSA-k-means	FMKGSO	MKGSO	TMKGSO
Disp01	$0.9902 \pm$ 0.0536	$0.9805 \pm$ 0.0741	1.0000 ± 0	1.0000 ± 0	$0.9906 \pm$ 0.0512	$1.0000 \pm$ $\mathbf{0}$	1.0000 ± 0	1.0000 ± 0
Disp02	$0.9598 \pm$ 0.0682	$0.9321 \pm$ 0.1036	$0.9601 \pm$ 0.0669	$0.9685 \pm$ 0.0500	$0.9435 \pm$ 0.0889	$0.9777 \pm$ $\mathbf{0}$	0.9777 ± 0	0.9777 ± 0
Disp03	$0.9462 \pm$ 0.0993	$0.9615 \pm$ 0.0671	$0.9695 \pm$ 0.0499	$0.9677 \pm$ 0.0582	$0.9696 \pm$ 0.0512	$0.9775 \pm$ $\mathbf{0}$	$0.9775 + 0$	0.9775 ± 0
Disp04	$0.9370 \pm$ 0.0800	$0.8743 \pm$ 0.1412	0.9580 ± 0	$0.9157 \pm$ 0.1097	$0.9368 \pm$ 0.0805	$0.9580 \pm$ $\mathbf{0}$	0.9580 ± 0	0.9580 ± 0
Disp05	$0.9469 \pm$ 0.0613	$0.9081 \pm$ 0.1154	0.9581 ± 0	$0.9454 \pm$ 0.0691	$0.9469 +$ 0.0610	$0.9581 \pm$ $\mathbf{0}$	0.9581 ± 0	0.9581 ± 0
Prox01	$0.9765 \pm$ 0.0893	$0.9616 \pm$ 0.1000	1.0000 ± 0	1.0000 ± 0	$0.9925 \pm$ 0.0409	$1.0000 \pm$ $\mathbf{0}$	1.0000 ± 0	1.0000 ± 0
Prox02	$0.8923 +$ 0.1178	$0.8839 \pm$ 0.1211	$0.9007 \pm$ 0.1138	$0.8923 +$ 0.1179	$0.8605 \pm$ 0.1255	$0.9681 \pm$ $\mathbf{0}$	0.9681 ± 0	$0.9597 \pm$ 0.0463
Prox03	$0.8533 \pm$ 0.1294	$0.8021 \pm$ 0.1255	$0.8588 \pm$ 0.1191	$0.8596 \pm$ 0.1179	$0.8386 \pm$ 0.1213	$0.9416 \pm$ $\mathbf{0}$	$0.9253 \pm$ 0.0622	$0.9326 \pm$ 0.0495
Prox04	$0.8298 \pm$ 0.1068	$0.8068 \pm$ 0.1164	$0.8276 \pm$ 0.1103	$0.8287 \pm$ 0.1050	$0.8594 \pm$ 0.0839	$0.8874 \pm$ $\bf{0}$	$0.8781 +$ 0.0509	$0.8433 \pm$ 0.1005
Prox05	$0.8452 \pm$ 0.0055	$0.8288 \pm$ 0.0567	$0.8435 \pm$ 0.0051	$0.8420 \pm$ 0.0064	$0.8453 \pm$ 0.0052	$0.8459 +$ $\mathbf{0}$	0.8459 ± 0	0.8459 ± 0

Table 13 Experimental results for synthetic data sets: correct Rand index (CR^{\dagger})

Table 14 Experimental results for synthetic data sets: weighted quantization error $(J_{e_2}^{\downarrow})$

Data set	GA-k-means	DE-k-means	PSO-k-means	GSO-k-means	BSA-k-means	FMKGSO	MKGSO	TMKGSO
Disp01	$3.2328 \pm$ 1.3206	3.4689 \pm 1.8158	2.9917 ± 0	2.9917 ± 0	$3.2413 \pm$ 1.3673	$2.9917 \pm$ $\mathbf{0}$	2.9917 ± 0	2.9917 ± 0
Disp02	7.1389 \pm 1.8318	$7.8209 \pm$ 2.6467	7.1258 \pm 1.7812	6.9053 \pm 1.3555	$7.6312 \pm$ 2.5251	$6.6578 +$ $\mathbf{0}$	6.6578 ± 0	6.6578 ± 0
Disp03	8.7565 \pm 1.7919	$8.6020 \pm$ 1.6466	8.4021 \pm 1.2750	8.3638 \pm 1.0658	$8.3862 \pm$ 1.1877	8.1691 \pm $\mathbf{0}$	8.1691 ± 0	8.1691 ± 0
Disp04	9.8094 \pm 1.3951	$10.905 \pm$ 2.4666	9.4428 ± 0	$10.173 \pm$ 1.8938	9.8259 \pm 1.4582	$9.4428 \pm$ $\mathbf{0}$	9.4428 ± 0	9.4428 ± 0
Disp05	$11.942 \pm$ 1.0423	$12.706 \pm$ 2.1751	11.752 ± 0	$11.937 \pm$ 1.0157	$11.923 +$ 0.9407	$11.752 +$ $\mathbf{0}$	11.752 ± 0	11.752 ± 0
Prox01	3.4074 \pm 1.1860	$3.7009 \pm$ 1.5691	3.0958 ± 0	3.0958 ± 0	$3.2518 \pm$ 0.8547	$3.0958 \pm$ $\mathbf{0}$	3.0958 ± 0	3.0958 ± 0
Prox02	$3.2264 +$ 0.5268	$3.2662 +$ 0.5451	3.1889 \pm 0.5088	$3.2271 +$ 0.5280	3.3829 \pm 0.5771	$2.8873 +$ $\mathbf{0}$	2.8873 ± 0	$2.9247 \pm$ 0.2052
Prox03	$3.3046 \pm$ 0.5948	$3.5301 \pm$ 0.5442	3.2661 \pm 0.4986	$3.2670 \pm$ 0.5000	$3.3933 \pm$ 0.5537	$2.9198 \pm$ $\mathbf{0}$	$2.9875 \pm$ 0.2578	$2.9561 \pm$ 0.1990
Prox04	$3.0787 +$ 0.4468	3.1564 \pm 0.4602	$3.0606 \pm$ 0.4115	$3.0668 \pm$ 0.4223	$2.9208 \pm$ 0.2409	$2.8419 \pm$ $\mathbf{0}$	$2.8680 +$ 0.1431	$2.9836 \pm$ 0.3271
Prox05	$2.7254 +$ 0.0003	$2.7735 \pm$ 0.1829	$2.7254 \pm$ 0.0003	$2.7253 \pm$ 0.0002	$2.7254 \pm$ 0.0002	$2.7252 \pm$ $\mathbf{0}$	2.7252 ± 0	2.7252 ± 0

individual from the population in each generation (MKGSO approach), represented the worst combination for the memetic models, in terms of computational costs (measured by the average execution time), but such combination still kept the good performances of all other evaluated memetic approaches (FMKGSO and TMKGSO),

Data set	GA-k-means	DE-k-means	PSO-k-means	GSO-k-means	BSA-k-means	FMKGSO	MKGSO	TMKGSO
Disp01	$3.8415 \pm$ 3.2624	4.4236 \pm 4.4823	3.2458 ± 0	3.2458 ± 0	$3.8622 \pm$ 3.3763	$3.2458 \pm$ $\bf{0}$	3.2458 ± 0	3.2458 ± 0
Disp02	13.375 \pm 3.4581	14.280 \pm 4.2780	13.378 \pm 3.4693	$12.921 \pm$ 2.4885	14.133 \pm 4.4080	$12.466 \pm$ $\mathbf{0}$	12.466 ± 0	12.466 ± 0
Disp03	$12.778 \pm$ 1.9092	12.834 \pm 2.6421	$12.630 \pm$ 2.5661	$12.411 \pm$ 1.3675	$12.581 \pm$ 2.3002	$12.162 \pm$ $\mathbf{0}$	12.162 ± 0	12.162 ± 0
Disp04	$12.033 \pm$ 1.9258	13.876 \pm 4.0761	11.527 ± 0	$12.576 \pm$ 2.7213	$12.225 \pm$ 2.6688	$11.527 \pm$ Ω	11.527 ± 0	11.527 ± 0
Disp05	13.070 \pm 2.0900	14.538 \pm 4.2234	12.689 ± 0	12.997 \pm 1.6910	$13.082 \pm$ 2.1547	$12.689 \pm$ $\mathbf{0}$	12.689 ± 0	12.689 ± 0
Prox01	$3.8775 +$ 2.2080	4.7609 \pm 3.7982	3.2974 ± 0	3.2974 ± 0	$3.6763 \pm$ 2.0760	$3.2974 +$ $\bf{0}$	3.2974 ± 0	3.2974 ± 0
Prox02	4.0095 \pm 1.2998	4.1025 \pm 1.3371	3.9165 \pm 1.2543	$4.0095 \pm$ 1.2998	4.3813 \pm 1.4055	$3.1729 +$ $\mathbf{0}$	3.1729 ± 0	$3.2658 \pm$ 0.5092
Prox03	3.9949 \pm 1.3633	4.6991 \pm 1.4728	3.9949 \pm 1.3633	3.9949 \pm 1.3633	4.3013 \pm 1.4605	3.0491 \pm $\bf{0}$	$3.2318 \pm$ 0.6953	$3.1512 \pm$ 0.5595
Prox04	3.5836 \pm 1.0632	3.7080 \pm 1.0301	$3.5448 \pm$ 0.9919	$3.5292 \pm$ 0.9737	$3.1859 \pm$ 0.4594	$3.0354 \pm$ $\bf{0}$	$3.0927 \pm$ 0.3139	$3.3622 \pm$ 0.7700
Prox05	$2.8617 \pm$ 0.0056	$2.9290 \pm$ 0.2535	$2.8609 \pm$ 0.0053	$2.8612 \pm$ 0.0055	$2.8612 \pm$ 0.0054	$2.8574 +$ $\bf{0}$	2.8574 ± 0	2.8574 ± 0

Table 15 Experimental results for synthetic data sets: intra-cluster distance (D_{max}^{\downarrow})

Table 16 Experimental results for synthetic data sets: inter-cluster separation (D_{min}^{\uparrow})

Data set	GA-k-means	DE-k-means	PSO-k-means	GSO-k-means	BSA-k-means	FMKGSO	MKGSO	TMKGSO
Disp01	$70.431 \pm$ 12.756	68.121 \pm 17.653	72.760 ± 0	72.760 ± 0	$70.417 \pm$ 12.825	$72.760 \pm$ $\mathbf{0}$	72.760 ± 0	72.760 ± 0
Disp02	$70.993 \pm$ 17.191	64.584 \pm 24.870	$71.257 \pm$ 16.153	73.106 \pm 13.117	$68.622 \pm$ 21.143	75.501 \pm $\mathbf{0}$	75.501 ± 0	75.501 ± 0
Disp03	$65.536 \pm$ 18.336	70.374 \pm 12.010	69.540 \pm 11.051	69.607 \pm 10.889	69.506 \pm 10.952	$71.723 +$ $\bf{0}$	71.723 ± 0	71.723 ± 0
Disp04	67.101 \pm 13.943	55.316 \pm 26.032	70.749 ± 0	62.980 \pm 20.149	$66.824 \pm$ 14.937	$70.749 \pm$ $\mathbf{0}$	70.749 ± 0	70.749 ± 0
Disp05	71.039 \pm 11.266	$70.418 \pm$ 15.102	73.096 ± 0	$71.290 \pm$ 9.8895	$71.105 \pm$ 10.899	73.096 \pm $\mathbf{0}$	73.096 ± 0	73.096 ± 0
Prox01	44.267 \pm 11.207	42.920 \pm 13.593	47.211 ± 0	47.211 ± 0	47.235 \pm 0.1302	47.211 \pm $\mathbf{0}$	47.211 ± 0	47.211 ± 0
Prox02	$9.9199 \pm$ 4.5234	9.5492 \pm 4.7216	$10.244 \pm$ 4.3635	9.8985 \pm 4.5569	$10.010 \pm$ 8.1720	$12.831 \pm$ $\mathbf{0}$	12.831 ± 0	$12.514 \pm$ 1.7377
Prox03	$8.9750 \pm$ 4.0264	7.6806 \pm 4.2878	8.9680 \pm 4.0372	8.9385 \pm 4.0787	$9.0281 \pm$ 4.0169	11.774 \pm $\mathbf{0}$	$11.205 \pm$ 2.1630	$11.497 \pm$ 1.5176
Prox04	$10.065 \pm$ 3.2354	9.8634 \pm 3.6289	$9.6850 \pm$ 3.6083	$10.000 \pm$ 3.4498	$10.838 \pm$ 2.4424	11.638 \pm $\mathbf{0}$	$11.365 \pm$ 1.4917	$10.253 \pm$ 3.1510
Prox05	$11.617 \pm$ 0.0477	$12.002 \pm$ 1.5962	$11.582 \pm$ 0.0650	$11.536 \pm$ 0.0531	$11.608 \pm$ 0.0566	$11.537 \pm$ Ω	11.537 ± 0	11.537 ± 0

showing its potential in terms of performance (considering all five selected clustering metrics) when dealing with data clustering problems, in comparison to all other state-of-theart non-memetic hybrid algorithms. In a future investigation, some mechanisms, like a selection process on the population individuals that would be improved by K-Means in each GSO generation, would be performed in an attempt to improve MKGSO computational costs.

Fig. 4 Results for Prox05 data set: a Original data set, b GA-k-means, c DE-k-means, d PSO-k-means, e GSO-k-means, f BSA-k-means, g FMKGSO, h MKGSO, i TMKGSO

As future works, we intend to evaluate the influence of the adopted distance function on the behavior of the proposed memetic algorithms, so the proposed approaches would be more robust to deal with clusters with different formats and shapes, overcoming some limitations of the standard Euclidean distance. Also, new fitness functions will be introduced and tested, using a larger testing bed, that will be obtained by the development of alternative synthetic data set configurations and scenarios. Finally, mechanisms for the automatic determination of the best number of clusters will be implemented, in such a way that such parameter would be estimated by the memetic evolutionary model itself, instead of being provided as an a priori input parameter to the algorithm.

Data set	GA-k-means	DE-k-means	PSO-k-means	GSO-k-means	BSA-k-means	FMKGSO	MKGSO	TMKGSO
Disp01	0.8063	0.7808	0.7798	0.9301	0.7955	0.7533	3.2231	0.5825
Disp02	0.7987	0.7803	0.7792	0.9301	0.7988	0.8456	3.2337	0.6636
Disp03	0.7968	0.7810	0.7832	0.9303	0.7990	0.9501	3.2367	0.7013
Disp04	0.8013	0.7821	0.7828	0.9320	0.8014	0.9412	3.2257	0.7482
Disp05	0.8005	0.7832	0.7817	0.9344	0.8016	1.0081	3.2252	0.8167
Prox01	0.7945	0.7813	0.7819	0.9320	0.7985	0.7967	3.2203	0.6293
Prox02	0.7999	0.7817	0.7782	0.9353	0.7990	0.9621	3.2320	0.7646
Prox03	0.8027	0.7832	0.7791	0.9371	0.8004	1.1342	3.2330	0.8668
Prox04	0.8018	0.7824	0.7819	0.9328	0.7988	1.2774	3.2399	0.9039
Prox05	0.8059	0.7888	0.7821	0.9352	0.8026	1.3448	3.2589	0.9195

Table 18 Average execution times for the synthetic data sets (in seconds)

Table 19 Overall evaluation on all data sets: average Friedman-Nemenyi ranks (and position) for each metric. The critical distance is $CD = 1.9169$. The best results are bold faced

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