Fuzzy Clustering with Evolutionary Algorithms

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Abstract

Objective function based fuzzy clustering aims at finding a fuzzy partition by optimizing a function evaluating a (fuzzy) assignment of a given data set to clusters, that are characterized by a set of parameters, the so-called prototypes. The iterative optimization technique usually requires the objective function not only to be differentiable, but prefers also an analytical solution for the equations of necessary conditions for local optima. Evolutionary algorithms are known to be an alternative robust optimization technique which are applicable to quite general forms of objective functions. We investigate the possibility of making use of evolutionary algorithms in fuzzy clustering. Our experiments and theoretical investigations show that the application of evolutionary algorithms to shell clustering, where the clusters are in the form of geometric contours, is not very promising due to the shape of the objective function, whereas they can be helpful in finding solid clusters that are not smooth, for example rectangles or cubes. These types of clusters play an important role for fuzzy rule extraction from data.

1 Introduction

Many fuzzy clustering algorithms are based on the idea to optimize an objective function. Usually, this objective function depends on the distances of the data to the clusters weighted by the membership degrees. By taking the first derivative of the objective function with respect to the cluster parameters, one obtains necessary conditions for the objective function to have an optimum. These conditions are then applied in an iteration procedure and define a clustering algorithm. This requires the objective function to be differentiable and

the iteration procedure can be computationally efficient only if the derived conditions lead to explicit equations for the cluster parameters. The set of cluster parameters, that determine the size and the shape of a cluster, depends on the specific application field. We mainly distinguish between fuzzy clustering as an explorative data analysis method, especially for unsupervised classification tasks, techniques for rule extraction (for instance for fuzzy controllers), and shell clustering algorithms, that are designed for boundary detection in image recognition.

Unfortunately, in many cases the restrictions that are enforced on the cluster parameters to be able to derive the iteration procedure are too narrow and exclude a lot of interesting possibilities. Thus it is desirable to have an alternative optimization technique that allows for more freedom in the choice of the cluster parameters.

In this paper we investigate whether evolutionary algorithms may be a solution to this problem. After a short introduction on fuzzy clustering in Section 2, we discuss the principal approach to use evolutionary algorithms for fuzzy clustering in Section 3. Section 4 is devoted to some theoretical and experimental results on applying evolutionary algorithms on the basis of the objective function introduced in Section 2. Evolutionary algorithms using other objective functions are discussed in Section 5.

2 Objective Function Based Fuzzy Clustering

Let us briefly review some fuzzy clustering methods based on objective functions. (For an overview see for example [8].) The cluster algorithm aims at minimizing the objective function

$$J(X, U, v) \;=\; \sum_{i=1}^{c} \sum_{k=1}^{n} (u_{ik})^m d^2(v_i, x_k) \tag{1}$$

under the constraints

$$\sum_{k=1}^{n} u_{ik} \;>\; 0 \qquad \text{for all } i \in \{1, \ldots c\} \tag{2}$$

and

$$\sum_{i=1}^{c} u_{ik} \;=\; 1 \qquad \text{for all } k \in \{1, \ldots n\}. \tag{3}$$

$X = \{x_1, \ldots, x_n\} \subseteq \mathbb{R}^p$ is the data set, c is the number of fuzzy clusters, $u_{ik} \in [0, 1]$ is the membership degree of datum x_k to cluster i, v_i is the prototype or the vector of parameters for cluster i, and $d(v_i, x_k)$ is the distance between prototype v_i and datum x_k. The parameter $m > 1$ is called fuzziness index. For $m \to 1$ the clusters tend to be crisp, i.e. either $u_{ik} \to 1$ or $u_{ik} \to 0$, for $m \to \infty$ we have $u_{ik} \to 1/c$. Usually $m = 2$ is chosen.

The objective function (1) to be minimized uses the sum over the quadratic distances of the data to the prototypes weighted with their membership degrees. (2) guarantees that no cluster is completely empty, (3) ensures that for each datum its classification can be distributed over different clusters, but the sum of the membership degrees to all clusters has to be 1 for each datum.

Differentiating (1), taking the constraints into account by Lagrange multipliers, one obtains

$$u_{ik} = \frac{1}{\sum_{j=1}^{c} \left(\frac{d^2(v_i, x_k)}{d^2(v_j, x_k)} \right)^{\frac{1}{m-1}}} \tag{4}$$

as a necessary condition for (1) to have a (local) minimum. The Equation (4) is therefore used for updating the membership degrees u_{ik} in an iteration procedure until the difference between the matrix (u_{ik}^{new}) and the matrix (u_{ik}^{old}) in the previous iteration step is less than a given tolerance bound ε.

The most simple fuzzy clustering algorithm is the fuzzy c-means (FCM) (see f.e. [2]) where the distance d is simply the Euclidean distance and the prototypes are vectors $v_i \in \mathbb{R}^p$. It searches for spherical clusters of approximately the same size and by differentiating (1) we obtain the necessary condition

$$v_i = \frac{\sum_{k=1}^{n} (u_{ik})^m x_k}{\sum_{k=1}^{n} (u_{ik})^m} \tag{5}$$

for the prototypes that are used alternatingly with (4) in the iteration procedure.

Gustafson and Kessel [7] designed a fuzzy clustering method that is looking for hyper-ellipsoidal clusters. The prototypes consist of the cluster centers v_i as in the FCM and (positive definite) covariance matrices C_i. The Gustafson and Kessel algorithm replaces the Euclidean distance by the transformed Euclidean distance

$$d^2(v_i, x_k) = (\rho_i \det C_i)^{1/p} \cdot (x_k - v_i)^\top C^{-1} (x_k - v_i).$$

These or similar algorithms can also be applied to learn fuzzy rules from data for classification problems [6, 12] or function approximation [11, 13, 21]. The fuzzy rules are derived from projections of the clusters leading to a certain loss of information which could be minimized if the clusters had the shapes of rectangles or cubes. Unfortunately, such shapes lead to non-differentiable objective functions.

In contrast to the methods that are designed for solid clusters, the so-called shell clustering algorithms are tailored for clusters in the form of boundaries of circles, ellipses, etc. (For an overview on shell clustering see [14].) Davé [5] developed one of the first shell clustering algorithms for the detection of circles. Each prototype consists of the cluster center v_i and the radius r_i. The distance function for the fuzzy c-shells algorithm (FCS) is

$$d^2((v_i, r_i), x_k) = (\| x_k - v_i \| - r_i)^2$$

so that exactly those data have zero distance to the cluster that are located on the circle with radius r_i and center v_i. Unfortunately, this distance function leads to a set of coupled non-linear equations for the v_i and r_i that cannot be solved in an analytical way. Thus an additional numerical iteration procedure to solve non-linear equations is necessary in each iteration step of the clustering algorithm which causes a lot of computational costs. Although this specific problem is solved for circles by the fuzzy c-spherical shells algorithm (FCCS) [16] using the distance function

$$d^2((v_i, r_i), x_k) = (\| x_k - v_i \|^2 - r_i^2)^2,$$

the general problem for other shell shapes remains.

3 Evolutionary Algorithms for Fuzzy Clustering

Evolutionary algorithms are a class of optimization methods that are inspired by the process of biological evolution. The principal idea is to have a collection or *population* of possible problem solutions encoded as parameter vectors – the *chromosomes* – that define a solution. From this population a new population – the next *generation* – is generated by first producing offspring from the old chromosomes by changing some components of the chromosomes, the *genes*, randomly and sometimes also by a mixing of genes of different chromosomes (*crossover*), and then by selecting the best chromosomes for the next generation. For details we refer to books like [1, 17].

As a quite general optimization strategy, evolutionary algorithms might be applicable to objective function based fuzzy clustering. Thus it is necessary to find a suitable coding of the parameters to be determined in fuzzy clustering. Obviously, the parameters to be optimized are the prototypes v_i and the membership degrees u_{ik}. In [3] it was proposed to perform hard clustering (i.e. $u_{ik} \in \{0, 1\}$), with genetic algorithms by taking the u_{ik} as the parameters for the evolutionary algorithm. For fuzzy clustering this does not seem to be suitable, since this means that besides the prototypes $c \cdot n$ real-valued parameters have to be optimized, where c is the number of clusters and n the number of data vectors. Since formula (4) is generally valid independent of the special type of prototype, it is not necessary to optimize the parameters u_{ik} by an evolutionary algorithm. In addition, the problems for the standard iterative optimization procedure are not caused by the membership degrees, but by the choice of the prototype parameters. Thus we restrict ourselves here to optimize only the prototypes by an evolutionary algorithm. The corresponding membership degrees are computed as in the usual algorithm on the basis of equation (4). In all our experiments in the following section, a chromosome is a vector with $c \cdot p$ real components. c is the number of clusters and p is the number of cluster (prototype) parameters, i.e. $p = 1$ for the FCM (an FCM-prototype is just the cluster centre v_i) and $p = 2$ for the FCCS (an FCSS-prototype consists of the cluster centre v_i and the radius r_i). Mutation is carried out by adding normally distributed values to the chromosome (vector) components. We applied various crossover operators (one-point, two-point, and uniform crossver). Although two-point crossover turned out to yield the best results, the general effect of crossover was almost neglectable.

The aim of fuzzy clustering depends on the application domain. In the case of data analysis and classification tasks as well as for rule extraction it is important to find an appropriate (gradual) assignment of the data to suitable prototypes, i.e., the emphasis is on the assignment and not on the exact prototype parameters. For shell clustering algorithms to detect geometrical objects in images the exact parameters, i.e. the prototype, play an important role. This means that it is not sufficient just to compile the data points on a circle in one cluster, but the exact values of the center and the radius of the circle are of great interest.

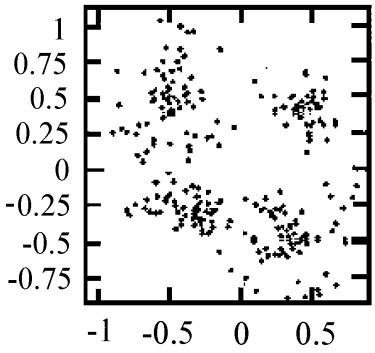

Figure 1: A test data set for the FCM

4 Experimental and Theoretical Results

Before we apply evolutionary algorithms to fuzzy clustering with non-standard prototypes, we first test their performance by two standard fuzzy clustering techniques, namely the FCM as an example for the search after solid clusters and the shell clustering algorithm FCS. Similar experiments were also carried out in [18] for the FCM. As in our approach, only the prototypes, i.e. for the FCM the cluster centers, are determined by the evolutionary algorithm in [18]. [18] also reports good results when the objective function (1) is replaced by the so-called partition coefficient, a validity measure that is sometimes used for the FCM to determine the number of clusters.

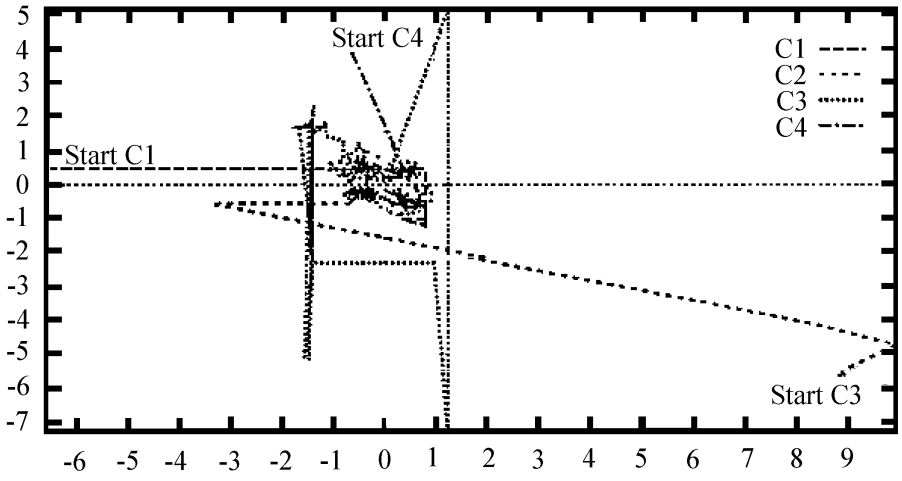

Figure 2: Tracking the prototypes

Figure 1 shows a test data set for the FCM with four clusters. Evolutionary algorithms with different parameters were all able to solve the problem of finding suitable prototypes. An example run is illustrated in Figure 2 where the tracks of the prototypes computed over the generations by the evolutionary algorithm are indicated. A comparison of different selection strategies (left to right: roulette wheel, remainder stochastic sampling, tournament) in Figure 3 shows that tournament selection has the best performance. In the figure the

dashed line indicates the average fitness and the other line the best fitness in each generation, averaged over 18 runs of the evolutionary algorithm.

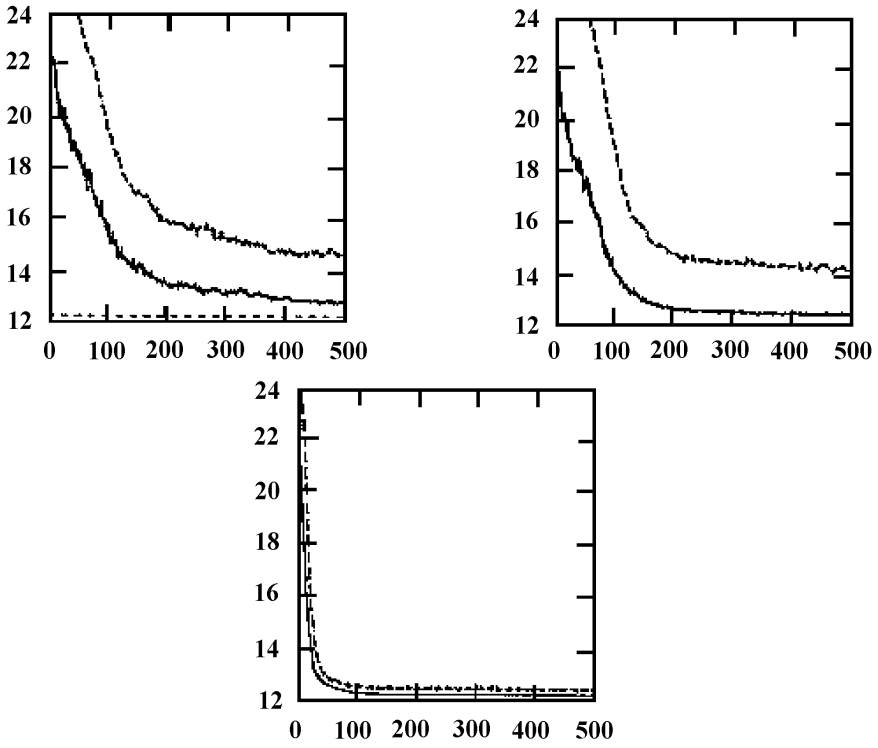

Figure 3: Comparison of selection strategies

Looking at these good results for the FCM we were quite optimistic also for shell clustering. However, the results were more or less disappointing. Figure 4 shows a test data set with five circles for the FCS and two results of the evolutionary algorithm. In one case no circle was detected correctly, in the other only one of the five. These results could not be improved, neither by experimenting with the mutation or crossover rate nor by introducing techniques like controlling the mutation rate on the basis of a span measure [17].

The typical results of the optimization of the objective function of the FCS by an evolutionary algorithm tend to yield larger circles – an observation which was also made in [4] where rectangular shells were considered. Thus we tested the evolutionary algorithm with a data set representing only one circle with center (0,0) and radius 2. In one case we limited the search space for the radius and the coordinates to the interval $[-2.2, 2.2]$, in the other case to the interval $[-22, 22]$. In the first case the evolutionary algorithm computed the correct radius and center in all test runs after about 30 generations. For the second case the circle was detected correctly only in about 60% of the cases after approximately 125 generations.

This motivated us to take a closer look at the fitness function. We consider again a circle with center (0,0) and radius 2 as the data set. In Figure 5 the evaluation of a chromosome is shown whose y-coordinate for the circle is the correct value 0, whereas the x-value is shifted to the right between 0 and 22. The different curves were drawn for radius values between 1 and 10. The

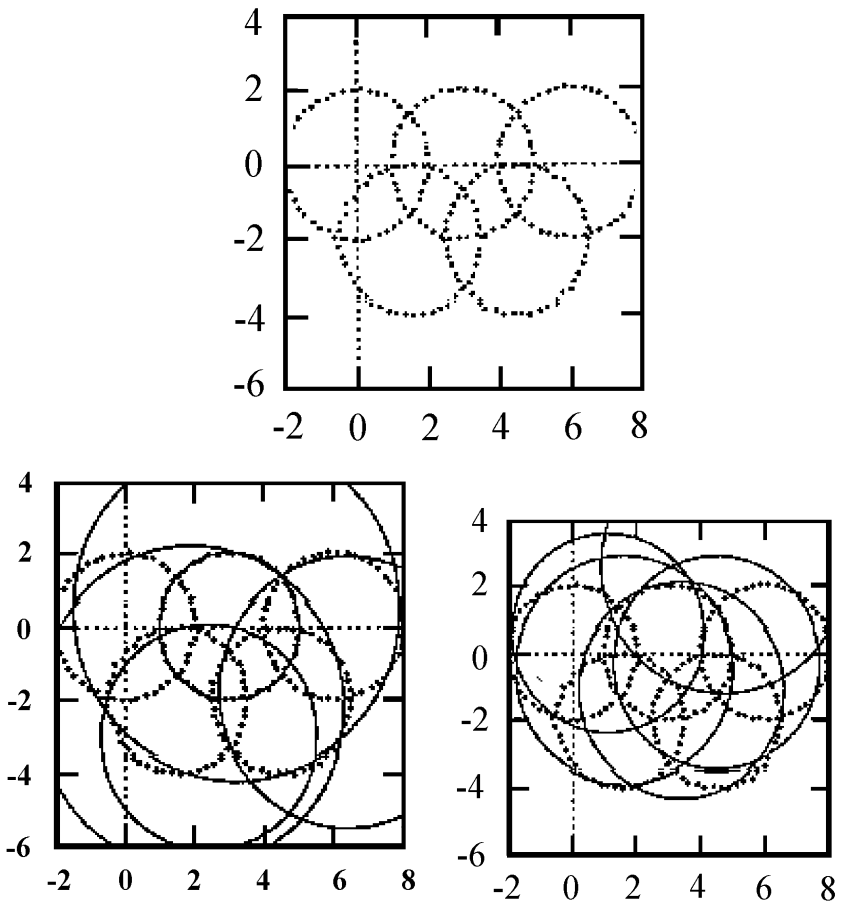

Figure 4: FCS: Test data and results

middle and lower diagram are just magnified scalings of the upper diagram.

From this figure it is obvious that the correct radius 2 gets the best evaluation only as long as the shifted circle center is not shifted to far away from the original circle center. The smaller radius 1 does never yield better values than a larger radius. And with increasing distance of the shifted center to the original center, a larger radius gets a better evaluation. This implies that in early generations where the random circle centers are still far away from the correct circle centers, chromosomes with smaller radius values are not being selected. But when these values are missing in the population, a random mutation to the correct radius without mutating also the center to the correct value leads to a very bad evaluation. Thus these genes are so strongly dependent on each other that only a simultaneous random jump of all genes to the correct values can lead to good results.

When we applied the evolutionary algorithm again to the data set of Figure 1 and limited the search range for radius genes to the interval $[0, 2.2]$, the results were satisfactory. In all test runs the circles were detected correctly after about 80 generations in average.

Our theoretical considerations and experiments show that the applicability of evolutionary algorithms to shell clustering is not very promising except when the parameter range can be restricted to quite limited bounds.

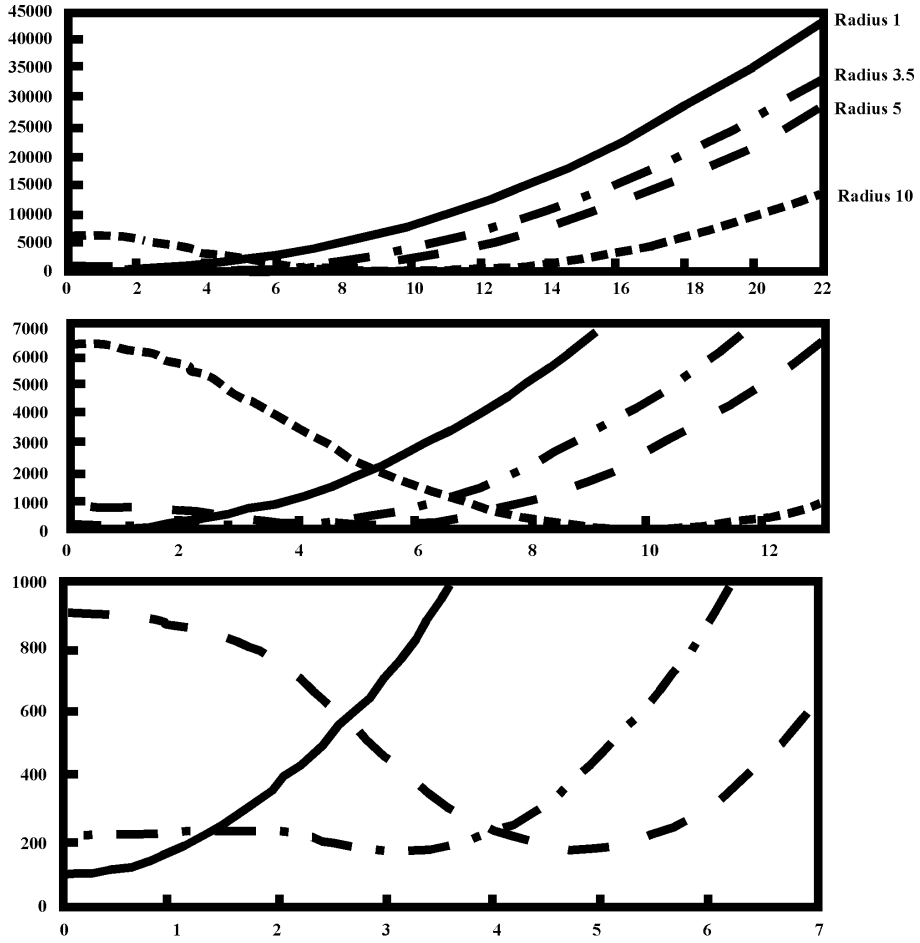

Figure 5: Evaluation of chromosomes with shifted circle centers

As mentioned in Section 2, the standard iteration procedure is difficult or impossible not only for certain types of shell clusters, but also for instance for solid rectangular clusters. Such clusters would be ideal for fuzzy rule extraction, especially when they are restricted to axes parallel rectangles or cubes.

In order to obtain clusters of this type we define as prototypes for p-dimensional data cluster centers $z_i \in \mathbb{R}^p$ and diagonal matrices B_i. As the distance function we choose

$$\begin{aligned} d^2(x_k, (z_i, B_i)) &= \| B_i(z_i - x_k) \|_\infty^2 \\ &= \left(\max_{1 \le j \le p} \{ b_{ij} \cdot |z_{ij} - x_{kj}| \} \right)^2 . \end{aligned}$$

In opposition to [4] we use the supremum norm $\| \ . \ \|_\infty$ instead of the 1-norm. In order to avoid the undesired solution $b_{ij} = 0$ for all i, j, we have to enforce a restriction on the matrices B_i. Analogously to the Gustafson and Kessel algorithms we require the matrix B_i to have a fixed value ϱ_i for the determinant, which determines the size or the volume of the cluster i. If nothing is known about the data, we simply choose $\varrho_i = 1$ for all $i = 1, \ldots, c$.

All results were quite satisfactory. In 90-100 percent of the test runs the evolutionary algorithm was able to assign the data to the correct clusters and the cluster centers were detected approximately correct. Figures 6 and 7 show

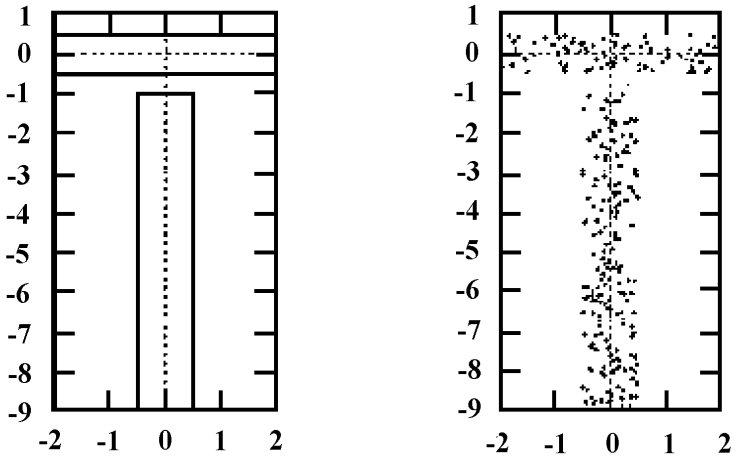

Figure 6: Two separated rectangular clusters (not filled and filled)

two 2-dimensional examples. In both cases we used two data sets – one in which data points were only placed on the edges of the rectangles and one where the rectangles were filled with data points.

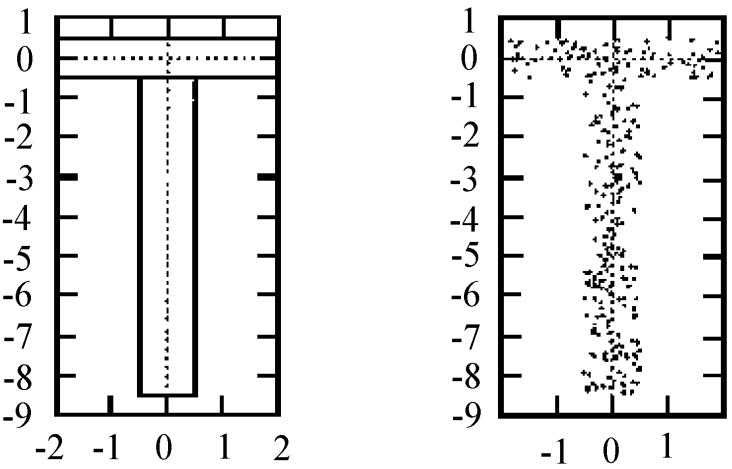

Figure 7: Two contiguous rectangular clusters (not filled and filled)

It is worth noticing that the best chromosome in Figure 7 for the data set with points only one the edges had the vectors $(0, -0.1)$ and $(0, -5.1)$ for the cluster centers and and the values 0.6 and 1.6 for the entree in the upper left corner of the diagonal matrices. The other non-zero value of the 2-dimensional diagonal matrix can then be calculated, since the determinants are fixed. This chromosome was assigned the error value 626.5 whereas the 'desired' solution with centers $(0, 0)$ and $(0, -4.5)$ and matrix values 0.5 and 2.0 gets a larger error value of 740.8. Thus this method is not suited for computing the correct parameters of rectangular shells. Nevertheless, the assignment of the data to the clusters was satisfactory even if data points were only present on the edges. Also the 3-dimensional case in Figure 8 and 9 caused no problems.

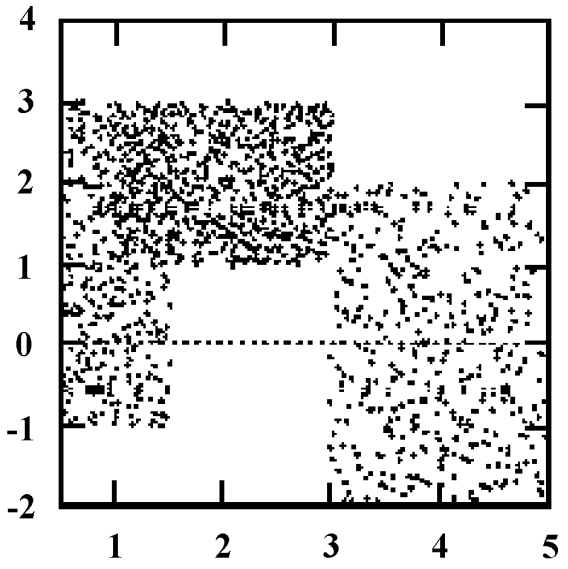

Figure 8: x/y-projections of three 3-dimensional clusters

5 Using Other Objective Functions

The good results on rectangular clusters described in the previous section can be applied to deriving fuzzy rules from data. In this section we take a closer look at evolutionary algorithm based fuzzy clustering and learning fuzzy rules. The principal idea to apply fuzzy clustering in order to derive if-then rules from data is that each cluster induces a rule by projecting the cluster to the corresponding coordinate spaces [11, 21]. The projection of a cluster to the i-th domain is obtained by taking the i-th coordinate of each data point and assigning to it the membership degree of the original data point to the cluster. In this way a discrete fuzzy set is defined on the i-th coordinate space. To extend this fuzzy set to the whole i-th domain, a piecewise linear fuzzy set can be defined on the basis of these discrete points, an enveloping fuzzy set or a suitable approximation by a parameterized fuzzy set like a triangular or trapezoidal membership function can be chosen [21].

In this way a cluster induces the rule

$$\text{If } \xi_1 \text{ is } \mu_1 \text{ and } \ldots \text{ and } \xi_{p-1} \text{ is } \mu_{p-1} \text{ then } \xi_p \text{ is } \mu_p.$$

where μ_i denotes the (extension of the) i-th projection of the considered cluster and $\xi_1, \ldots, \xi_{p-1}$ are input variables and ξ_p is the output variable. In this way a Mamdani-type fuzzy controller is defined [11, 21].

A number of variants of this principle were proposed by different authors to solve control, function approximation and classification problems with fuzzy rules (for a brief overview see [10]).

Fuzzy clusters are usually not bounded in the sense, that even data very far away from the cluster center have non-zero membership degree, although this degrees tend to be small as long as the number of clusters is not too small. In [9] an evolutionary algorithm based fuzzy clustering algorithm was proposed that constructs membership functions in the form of hyper-cones for the clusters so that the membership degree to a cluster becomes zero out

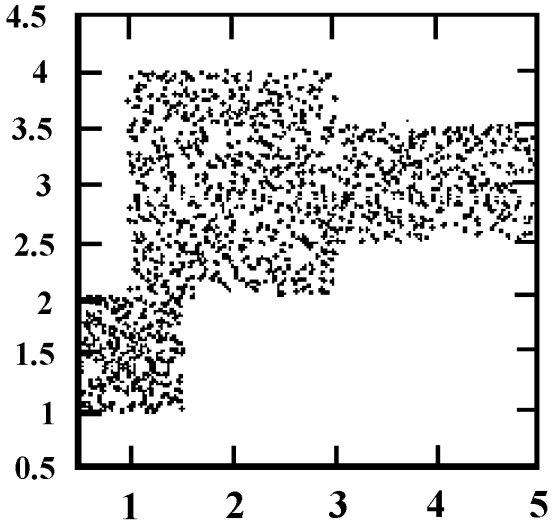

Figure 9: x/z-projections of three 3-dimensional clusters

of a region in the form of a hyper-ellipsoid. Nevertheless, the more serious problem appearing when extracting rules from fuzzy clusters is not solved by this approach, namely the problem of a certain loss of information enforced by the projection of a multidimensional cluster.

The approach described in [20] constructs for each cluster a collection of triangular fuzzy sets – one for each dimension – with an evolutionary algorithm and avoids the projection of the clusters in this way. However, there are no restrictions for the fuzzy sets with triangular membership functions so that it is possible that one might be contained in another one or that they strongly overlap.

A method for constructing rules by fuzzy clustering that restricts to well-behaved triangular membership functions (in the sense that the membership degrees at each point add up to 1) is the so-called grid clustering [10], a fuzzy clustering algorithm that aims at finding fuzzy partitions for the single domains on the basis of multidimensional data. For the grid clustering we assume that we are given a data set $\{x_1, \ldots, x_n\} \subset \mathbb{R}^p$. We are looking for fuzzy partitions on the single domains consisting of triangular membership functions with the restrictions that for each domain at most two supports of different fuzzy sets have a non-empty intersection and the sum of the membership degrees is one at any point of the domain. This means that we have to determine a suitable grid in the multidimensional space. We assume that for each domain the number of triangular membership functions is predefined. We define the membership degree of a data point to a cluster represented by a grid point as the minimum of the membership degrees of the triangular membership functions whose tips are the projections of the grid point. The grid clustering algorithm introduced in [10] is not based on an objective function, but relies on a heuristic strategy for constructing the clusters.

In order to improve this grid clustering algorithm, we have designed a suitable objective function that can be optimized by an evolutionary algorithm.

The aim of our strategy is to place the prototypes on a suitable grid in the

multidimensional space in order to get fuzzy partitions on the single domains consisting of triangular membership functions. For this reason our objective function should not depend on the order in which we analyze the single dimensions. The coordinates of the prototypes should only be influenced by those coordinates of the data that are relatively near (in this case near means a small Euclidean distance) to the prototype's coordinate. A simple way in single dimensions is to let the data between two prototypes only influence these two. On a grid there are a lot of prototypes with the same coordinates in a single dimension. In the objective function each coordinate has only once to be taken into account. These considerations led us to the following objective function:

$$J = \sum_{r=1}^{p} \sum_{j=1}^{k_r} \left(\sum_{\substack{s \in \{1,\ldots,n\}: \\ c_{j-1}^{(r)} < x_{sr} < c_j^{(r)}}} \left(\frac{c_j^{(r)} - x_{sr}}{c_j^{(r)} - c_{j-1}^{(r)}} \right)^m + \sum_{\substack{\ell \in \{1,\ldots,n\}: \\ c_j^{(r)} < x_{\ell r} < c_{j+1}^{(r)}}} \left(\frac{x_{\ell r} - c_j^{(r)}}{c_{j+1}^{(r)} - c_j^{(r)}} \right)^m \right). \quad (6)$$

x_{sr} and $x_{\ell r}$ are the r-th coordinate of the data x_s and x_ℓ, respectively ($s, \ell \in \{1, \ldots, n\}$). We assume that we have in the r-th dimension ($r \in \{1, \ldots, p\}$) k_r triangular fuzzy sets. Each triple $c_{j-1}^{(r)}, c_j^{(r)}, c_{j+1}^{(r)}$ induces a triangular membership function with the interval $(c_{j-1}^{(r)}, c_{j+1}^{(r)})$ as the support and $c_j^{(r)}$ as the point with the membership degree one. Thus the fractions in the sums (without the power m) provide the value one minus the membership degree to the triangular membership function with the tip at $c_j^{(r)}$ of the data (or better: their r-th projection) that lie on the support of the membership function. Since we add up the values for all triangular fuzzy sets (and the sum of the membership degrees of a datum to neighbouring fuzzy sets yields one), we obtain the smallest considerable value of (6), when all the memberships degree are 0.5 (as long as $m > 1$ is chosen) and the largest considerable value, when the membership degrees are either zero or one. We want the $c_j^{(r)}$ to be in the center of data clusters, i.e. the membership degree is high (near one) for data in the cluster and low (near zero) for data in other clusters. Thus we aim at maximizing (6). Note that (6) is a measure very similar to the partition coefficient [2].

A special treatment is needed for the data on the left/right of the leftmost/rightmost prototype (the values $c_1^{(r)}$ and $c_{k_r}^{(r)}$). In the beginning, we assume that the $c_j^{(r)}$ are uniformly distributed, i.e. equidistant. We add in each dimension two additional prototypes $c_0^{(r)}$ and $c_{k_r+1}^{(r)}$, again equidistant to the left and right of $c_1^{(r)}$ and $c_{k_r}^{(r)}$. The values $c_0^{(r)}$ and $c_{k_r+1}^{(r)}$ are assumed to be fixed and must not be changed by the evolutionary algorithm. Nevertheless, we have to take these additional prototypes into account in the objective function so that the data at the edge of the universe have the same influence as the data in the middle. This means that the second sum in (6) actually goes from $j = 0$ to $j = k_r + 1$. For the construction of the prototypes we only need the grid coordinates in each dimension.

We apply a standard $(1, b)$, respectively $(1 + b)$ evolution strategy, i.e. our population consists of just one chromosome from which generate b children.

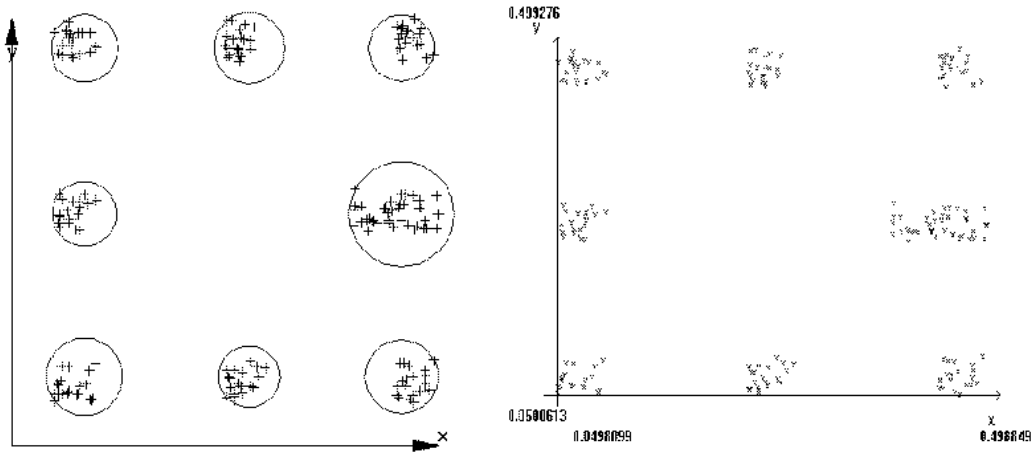

Figure 10: Left: Original grid clustering. Right: ES-based grid clustering

The chromosome in the succeeding generation is the best chromosome of the b children in case of the ,-strategy and the best chromosome of the b children and the parent chromosome in case of the +-strategy.

If we have a p-dimensional domain of interest, a chromosome of the evolutionary algorithm consists of p vectors. Projecting the grid points (prototypes) into dimension r, leads to the attributes of vector r. The p vectors do not need to have the same number of components. If we consider for instance two-dimensional data, i.e. $p = 2$, with three grid points in the x-dimension ($k_1 = 3$) and two grid points in the y-dimension ($k_2 = 2$) (not including the two fixed boundary points in each dimension), then a typical chromosome would look like $((1.2, 2.1, 3.4), (0.8, 2.7))$. This would mean that we have the six (inner) grid points $(1.2, 0.8), (2.1, 0.8), (3.4, 0.8), (1.2, 2.7), (2.1, 2.7), (3.4, 2.7)$ (plus 14 grid points at the boundary).

The descendants are generated from the parent chromosome by adding normally distributed values to the components, ensuring that all grid coordinates remain within the boundaries determined by the data at the very boundaries in each dimension.

Since the objective (6) consists of a sum over the dimensions, each dimensions contributes a value to the sum independent of the other dimensions. Therefore, the grid coordinates can be optimized independently in each dimension. Thus we do not just choose the best chromosome for the next generation, but consider each sub-vector, specifying the grid coordinates in one dimension, separately and choose the best one in each dimension.

In most cases the initialization with equidistant $c_j^{(r)}$ is good enough that the +-strategy does not fall into local optima. So the +-strategy leads to good results. In case of the ,-strategy the population of descendants has to be much bigger than the population so that the evaluation time increases drastically. All combinations of coordinates have to be computed in each dimension. If k_r is the size of the population and b is the number of descendants in this dimension

$$\sum_{r=1}^{p} \binom{b}{k_r} \text{ (,-strategy), respectively} \qquad \sum_{r=1}^{p} \binom{b+k_r}{k_r} \text{ (+-strategy)}$$

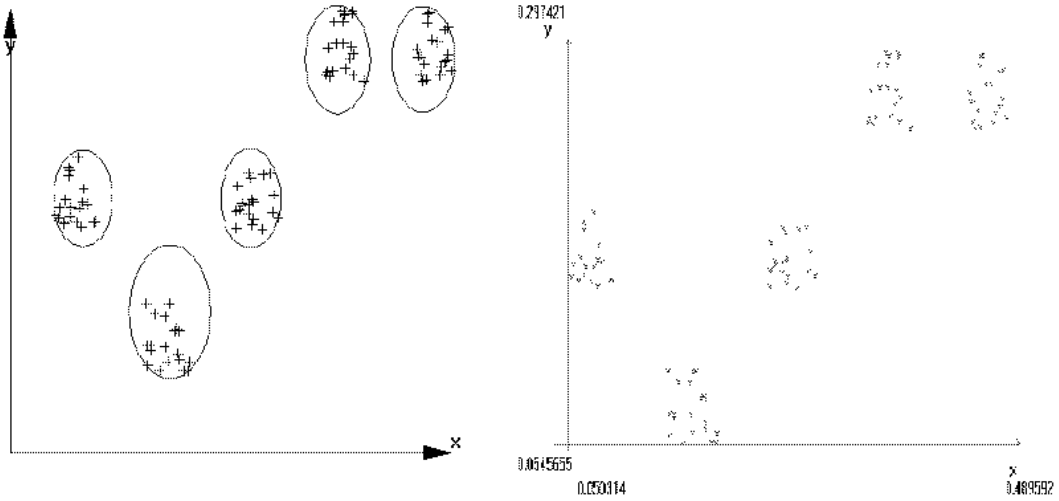

Figure 11: Left: Original grid clustering. Right: ES-based grid clustering

possible combinations have to be evaluated. Maybe heuristics like tabu search can help to reduce the evaluation time. The chosen strategy for the examples is the +-strategy with variance 0.9 and 100 Iterations. Figures 10 and 11 illustrate the results for two examples (showing only the grid points).

To compare the evolutionary algorithm with the original heuristic grid clustering algorithm described in [10], the results of the latter one were evaluated with the same objective function. The advantages of the evolutionary computation with the proposed objective function are the same as the advantages of the grid clustering algorithm. The result of the evolution strategy (Figure 10: 310.465, Figure 11: 159.515) is in both examples better than the results of the original grid clustering algorithms (for the data set in Figure 10: 303.906, in Figure 11: 159.438). Using the +-strategy, the evolutionary algorithm has the advantage of monotone increasing objective function values.

The progress of (6) is shown in Figure 12. The upper diagram shows the value of (6) for the dataset from Figure 10.

The first example has its best value after the first iteration, when the heuristic grid clustering algorithm is evaluated. This good value is not reached again during the computation. Faster evaluation is an advantage of the grid clustering algorithm. The values of the objective function for the heuristic grid clustering are only slightly smaller than the results obtained from the evolution strategy. It has to be taken into account that the grid clustering algorithm is not objective function based and therefore not tailored for the particular objective function of the evolutionary algorithm. Nevertheless, the results show that it is a very good heuristic method being much faster than the evolution strategy.

6 Conclusions

Our investigations show that shell clustering with evolutionary algorithms seems to be quite problematic, since there exist a lot of local optima and the correct solution often looks like a very narrow optimum. The situation is

better for solid clusters. It should however be noted that evolutionary algorithms require a much longer computation time to solve the problem compared to the standard iteration procedure so that the application of evolutionary algorithms makes only sense when the standard iteration procedure cannot be applied according to a non-differentiable distance function or when an analytical solution for the single iteration steps cannot be found. Even in that a case, as our grid clustering example shows, a good heuristic algorithm can lead to results almost as good as the ones obtained by evolutionary computation in a much shorter time.

We have restricted our investigations to probabilistic clustering, requiring that the membership degrees of a datum to the clusters add up to one. In principal, we can as well use the possibilistic version of the objective function [15] dropping the probabilistic constraint.

Another important question is a strategy for determining the number of clusters. This can be done in the usual way on the basis of suitable validity measures. However, this requires to carry out the clustering for varying numbers of cluster and increases the computation time.

We obtained the most promising results for clusters suitable for constructing fuzzy rules. Of course, there are other techniques of learning rules, like neuro-fuzzy approaches (for an overview see [19]). However, it turns out that most of these approaches are well suited for tuning the fuzzy sets, but not so for detecting rules.

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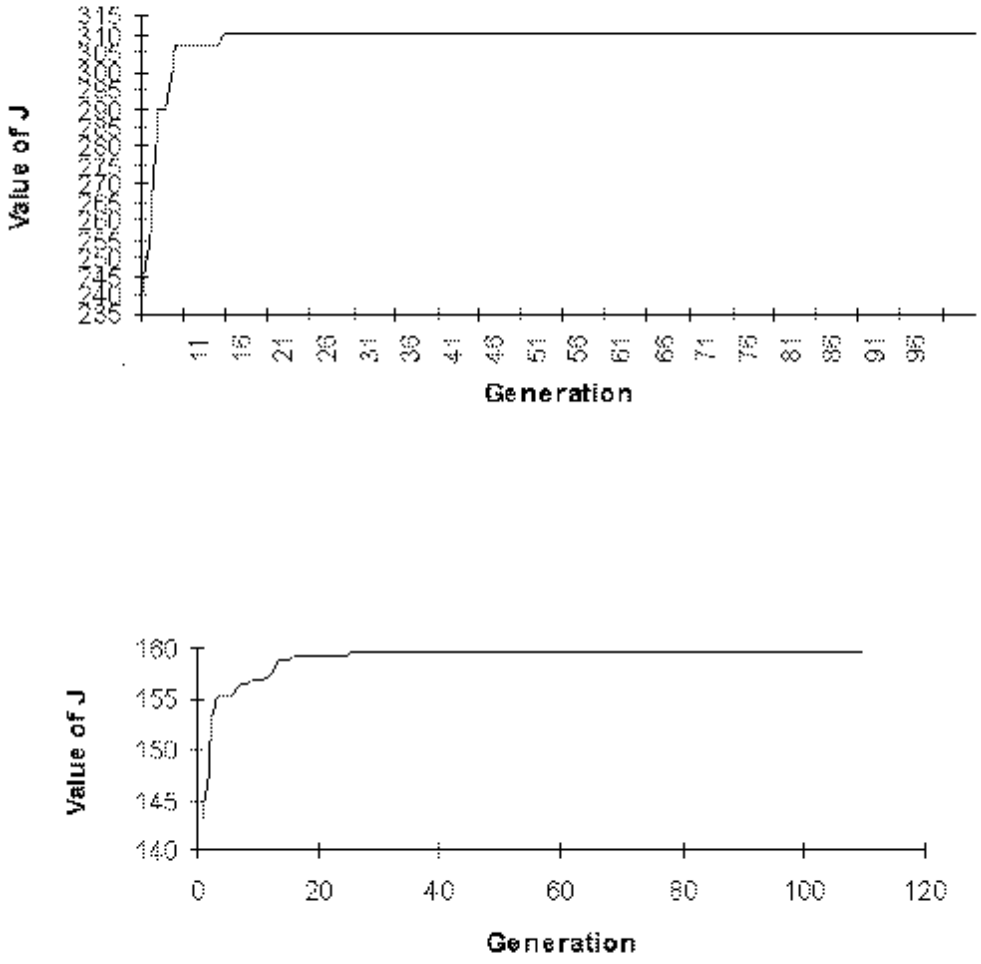

Figure 12: Improvement of (6) by the evolutionary algorithm for the data sets shown in Figures 10 and 11