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[Baohua Liang](#) ^{*}, Erli Jin, Liangfen Wei, [Rongyao Hu](#)

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Article

Attribute Reduction Algorithm of Knowledge Granularity for Incomplete System under the Background of Clustering

Baohua Liang ^{1,2,3}, Erli Jin ³, Liangfen Wei ³ and Rongyao Hu ^{4,*}

¹ Guangxi Key Lab of Multi-Source Information Mining & Security, Guangxi Normal University, Guilin 541004, China;

² School of Computer science and Engineering, Guangxi Normal University, Guilin 541004, China;

³ School of Computer and Artificial Intelligence, Chaohu University, Hefei, 238000, China;

⁴ CBICA, University of Pennsylvania, Philadelphia, PA 19104, USA.

* Correspondence: rongyao.hu@pennmedicine.upenn.edu

Abstract: The phenomenon of missing data can be seen everywhere in reality. Most typical attribute reduction models are only suitable for complete systems. But for incomplete systems, we can't obtain the effective reduction rules. Even if there are a few reduction approaches, the classification accuracy of their reduction sets still needs to be improved. In order to overcome these shortcomings, this paper firstly defines the similarities of intra-cluster objects and inter-cluster objects based on the tolerance principle and the mechanism of knowledge granularity. Secondly, attributes are selected on the principle that the similarity of inter-cluster objects is small and the similarity of intra-cluster objects is large, and then the knowledge granularity attribute model is proposed under the background of clustering; Then, the IKAR algorithm program is designed. Finally, a series of comparative experiments about reduction size, running time and classification accuracy are conducted with twelve UCI data sets to evaluate the performance of IKAR algorithms, then the stability of Friedman test and Bonferroni-Dunn test are conducted. The experimental results indicate that the proposed algorithms are efficient and feasible.

Keywords: attribute reduction; knowledge granularity; clustering; similarity

1. Introduction

Rough set theory (RST) [1], initiated by Pawlak, is an effectively mathematical tool to deal with imprecise, fuzzy and incomplete data. RST has been successfully applied in machine learning [2–4], knowledge discovery [5,6], expert system [7], disease diagnostics [8–10], decision support [11,12] and other areas [13–15]. Attribute reduction is one of the research hotspots in RST. As an important technology in the process of data preprocessing, attribute reduction has captured researchers' attention in big data and knowledge discovery [16–18]. The main objective of attribute reduction is to remove some irrelevant or non-important attributes while keeping the original distinguishing ability unchanged. In this way, the effect of data dimension reduction can be achieved, and a lot of time and space resources can be saved for the process of knowledge discovery and rule extraction.

With the rapid development of network and information technology, we have gradually entered the era of big data. The data sets have the characteristics of large volume, rapid change, and diverse data forms [19]. At the same time, due to the influence of the collection method and environment during the data collection process, there are a large number of missing data or wrong data in the data sets. Because of the existence of these disturbed data, it will seriously affect the decision-making and judgement of big data, and even mislead decision makers. After a long period of unremitting endeavor, scholars have achieved outstanding results in attribute reduction [20–24,40]. For example, Dai [20] proposed a semi-supervised attribute reduction based on attribute indiscernibility. Cao [21] put forward a three-way approximate reduction approach by using information-theoretic measure.

Yang [22] presented a novel incremental attribute reduction method via quantitative dominance-based neighborhood self-information. Lin [16] et al. developed a feature selection way by using neighborhood multi-granulation fusion. In variable precision rough set, Yu [24] et al. raised a novel attribute reduction based on local attribute significance. In literature [40], Devi proposed a new dimension reduce technology by considering the picture of fuzzy soft matrices in the decision-making process. Above these classic reduction models are suitable only for complete systems.

The phenomenon of missing valuable data set in reality is ubiquitous, and the traditional reduction models can't obtain effective reduction rules. In order to reduce the incomplete system, Zhang and Chen proposed a lambda-reduction approach based on the similarity degree respect to a conditional attribute subset for incomplete set-valued information systems [25]. For incomplete interval-valued information systems (in short IIIS), Li [35] proposed the concept of similarity degree and tolerance relation between two information values of a given attribute. Then three reduction algorithms based on theta-discernibility matrix, theta-information entropy and theta-significance are designed. Liu introduced a new attribute reduction approach by using conditional entropy based on the fuzzy alpha-similarity relation [28]. Subsequent, Dai [36] proposed interval-valued fuzzy min-max similarity relations and designed two attribute reduction algorithms based on interval-valued fuzzy discernibility pairs model. Song[38] put forward the similarity degree between information values on each attribute and an attribute reduction method is designed by using information granulation and information entropy. Zhou presented a heuristic attribute reduction algorithm with a binary similarity matrix and attribute significance as heuristic knowledge under incomplete information systems [26]. Zhang[37] presented a novel approach for knowledge reduction by using the discernibility techniques in multi-granulation rough set model. He and Qu [27] put forward the fuzzy-rough iterative computation model based on symmetry relations for an incomplete categorical decision information system. Sirekha et al. proposed an attribute reduction on SE-ISI concept Lattice based on concept of object ranking for incomplete information systems [29]. Cornelis et al. put forward a generalized model of attribute reduction using fuzzy tolerance relation within the context of fuzzy rough set theory [30]. Liu applied the concept of accurate reduction and reduced invariant matrix for reducing attribute under information systems [31]. To reduce unnecessary tolerance classes for the original cover, Nguyen [39] introduced a new concept of stripped neighborhood covers and proposed an efficient heuristic algorithm in mixed and incomplete decision tables.

Although the above reduction algorithms can effectively reduce the incomplete information systems, the classification accuracy of the reduced set is not ideal. The main reason is that only the importance of attributes is considered when selecting attributes, and the impact of attributes on classification is not considered. Usually, people take the best result of clustering as a reference standard for classification work, and classify similar samples into the same cluster. In order to solve the problems mentioned above, this paper proposes an attribute reduction method from the perspective of clustering.

At present, the studies of attribute reduction on using the clustering idea to construct a feature selection model are relatively infrequent. In order to avoid or reduce the loss of some original information after discretization of continuous values, Zhang [44] proposed a feature selection method based on fuzzy clustering, but this method has no reliable theoretical support. Jia [41] proposed a spectral clustering method based on neighborhood information entropy feature selection, which uses feature selection method to remove redundant features before clustering. In order to take the classification effect of the data set into consideration when selecting features, Zhao proposed a Fuzzy C-Means clustering fuzzy rough feature selection method [43], which can improve the classification accuracy of the reduced set in certain degree, but the effect is not obvious. Jia proposed a similarity attribute reduction in the context of clustering in literature [42], which can greatly improve the classification accuracy of the reduced set, but it needs to continuously adjust the parameters to achieve the best classification effect. Such a feature set has certain random limitations, and increase the time consumption of the system. Therefore, it is necessary to design a stable model with high classification accuracy for data preprocessing.

Although these existing approaches can effectively reduce incomplete systems, they only consider the importance of the attributes themselves, and do not consider the correlation between attributes. The influence of conditional attributes on decision classification is not considered. In order to improve the classification accuracy of reduction set, we apply the idea of clustering.

Based on the principle that the similarity of samples within a cluster is as large as possible and the similarity of samples between clusters is as small as possible, an attribute reduction algorithm for an incomplete system is designed under the background of clustering. First, according to the principle of tolerance and the theory of knowledge granularity, we define the similarity of intra-cluster and inter-cluster for an incomplete system. Secondly, a formula for calculating the similarity of intra-cluster and inter-cluster objects is designed. After normalizing the two similarities, we define the similarity of objects. Then, according to the corresponding similarity mechanism, a new attribute reduction algorithm for incomplete system is proposed. Finally, a series of experiments have verified that the algorithm proposed in this paper is significantly better than other similar algorithms in terms of running time, accuracy and the stability of algorithm is analyzed by using Friedman test and Bonferroni-Dunn test in statistics.

The contribution of this paper is embodied in the following four aspects:

- 1) A tolerance class calculation in incomplete information systems is proposed and applied to knowledge granularity calculation.
- 2) Knowledge granules is used as a measure of sample similarity to measure the similarity of inter-cluster samples and intra-cluster samples.
- 3) A knowledge granularity reduction algorithm based on clustering context is designed in incomplete information systems.
- 4) Lots of experiments have been done to verify the validity of the algorithm proposed in this paper, and the stability of the algorithm is verified by mathematical statistics.

The rest parts of this paper are constructed as follows. The principle of tolerance and related concepts of knowledge granularity are recalled in section II. In Section III, we propose a similarity measure of intra-cluster and inter-cluster objects, and discuss the reduction mechanism according to the clustering background for missing data set. We normalize the similarity of inter cluster and intra-cluster, then design the corresponding reduction model in section IX. In section V, a series of experiments are conducted and the performance of the algorithm is evaluated from the reduction size, running time, classification accuracy and stability. Then the feasibility and effectiveness of the algorithm are verified. Finally, the advantages and disadvantages of the proposed algorithm in this paper are concluded and unfolded in the future work.

2. Preliminaries

In this section, we review some basic concepts in rough set theory, the definitions of tolerance class, knowledge granularity, clustering metrics, and the significance of attribute for incomplete decision systems.

2.1. Basic concept of RST

A decision information system is a quadruple $DS = (U, A, V, f)$, where U is a non-empty finite set of objects and A is a finite nonempty attribute sets; If $A = C \cup D$, where C is the conditional attribute sets and D is the decision attribute set; V is the union of attribute domains, $V = \bigcup_{a \in A} V_a$, V_a is the value set of attribute a , called the domain of a ; $f: U \times A \rightarrow V$ is an information function with $f(x, a) = V_a$ for each $a \in A$ and $x \in U$. For every attribute subset $B \subseteq C$, a indiscernibility relation is defined as follows:

$$IND(B) = \{(x, y) \in U \times U \mid \forall a \in B, f(x, a) = f(y, a)\} \quad (1)$$

By the relation $IND(B)$, we can get the partition of U denoted by $U / IND(B)$ or U / B . If $B \subseteq A \wedge X \subseteq U$, the upper approximation denotes as

$$\overline{B}(X) = \{x \in U \mid [x]_P \cap X \neq \emptyset\} \quad (2)$$

The lower approximation of X can be denoted as

$$\underline{B}(X) = \{x \in U \mid [x]_P \subset X\} \quad (3)$$

Where the objects in $\overline{B}(X)$ may belong to X , while the objects in $\underline{B}(X)$ must belong to X .

Definition 1. Given an incomplete decision system $IDS = (U, C \cup D, V, f)$, $\forall P \subseteq C$, the binary tolerance relations between objects that are indiscernible in terms of value of attribute in P is defined as following:

$$T(P) = \{\forall a \in P, f(a, x) = f(a, y) \vee f(a, x) = * \vee f(a, y) = *\} \quad (4)$$

Where $*$ represents missing value. $T(B)$ is symmetric and reflexive, but not transitive.

Definition 2. Given an incomplete decision system $IDS = (U, C \cup D, V, f)$, $\forall P \subseteq C$, and $o \in U$, the tolerance class of object o with respect to attribute P is defined as follows:

$$T_P(o) = \{y \mid (o, y) \in T(P)\} \quad (5)$$

2.1. Basic concept of knowledge granularity

Definition 3. Suppose $IDS = (U, C \cup D, V, f)$ is an incomplete decision system, $\forall P \subseteq C$. $T_P(o_i)$ is the tolerance class of object o_i with respect to P . The knowledge granularity of P on U is defined as follows:

$$GK_U(P) = \frac{1}{|U|^2} \sum_{i=1}^{|U|} |T_P(o_i)| \quad (6)$$

Where $|U|$ represents the number of objects in data set U . Since the reflexivity and symmetry of $T_P(o_i)$, there are lot of repeated calculations when calculating $\sum_{i=1}^{|U|} |T_P(o_i)|$. In order to reduce the amount of calculation, we propose the definition 4 as following.

Definition 4. Suppose $IDS = (U, C \cup D, V, f)$ is an incomplete decision system, $\forall P \subseteq C$. $CT_P(o_i)$ is the simplified tolerance relation of object o_i with respect to P is defined as follows:

$$\begin{aligned} CT_P(o_i) = \{o_j \mid \forall a \in P, (f(a, o_i) = f(a, o_j) \vee (f(a, o_i) \\ = * \wedge f(a, o_j) = *) \wedge (i < j)) \vee (f(a, o_i) = * \wedge f(a, o_j) \neq *)\} \end{aligned} \quad (7)$$

Definition 5. Given an incomplete decision system $IDS = (U, C \cup D, V, f)$, $\forall P \subseteq C$ and $o \in U$, the simplified tolerance class of object o with respect to attribute P is defined as follows:

$$CT_P(o) = \{y \mid (o, y) \in CT(P)\} \quad (8)$$

We delete the symmetric element pair and reflexive element pair from definition 3 and obtain definition 5, so the definition 5 does not have the characteristics of symmetry and reflexivity.

Definition 6. Given an incomplete decision system $IDS = (U, C \cup D, V, f)$, $\forall P \subseteq C, o \in U$, $CT_P(o)$, is the simplified tolerance class of object o with respect to attribute P . The equal knowledge granularity of P on U is defined as following:

$$EGK_U(P) = \frac{1}{|U|^2} \sum_{o \in U} |CT_P(o)| \quad (9)$$

Theorem 1. Given an incomplete decision system $IDS = (U, C \cup D, V, f)$, $\forall P \subseteq C$. Let $U/P = \{X_1, X_2, \dots, X_l\}$, $X_i \subseteq U/P$, $|X_i| = n_i$ where $1 \leq i \leq l$. All objects in subdivision X_i are missing value on attribute P , $|X_l| = n_*$. For the convenience of the following description, we mark X_i as X_* . Objects with

all non-missing values on attribute P are marked with $\overline{X_*}$. $EGK_U(P)$ represents the equal knowledge granularity of P on U , we have:

$$EGK_U(P) = \frac{1}{|U|^2} \left(\sum_{i=1}^l C_{n_i}^2 + |X_*| |\overline{X_*}| \right) \quad (10)$$

Where $C_n^2 = \frac{n(n-1)}{2}$.

Proof. Suppose that $\forall o \in X_i$ and $\forall o_* \in X_*$. According to definition 4, we can get $\sum_{o \in X_i} |CT_P(o)| = \frac{n_i(n_i-1)}{2} + n_i n_*$ and $\sum_{o_* \in X_*} |CT_P(o_*)| = \frac{n_*(n_*-1)}{2}$. Suppose $\forall o \in U$, according to definition 5, then we

can get $EGK_U(P) = \frac{1}{|U|^2} \sum_{o \in U} |CT_P(o)| = \frac{1}{|U|^2} \left(\sum_{i=1}^{l-1} \sum_{o \in X_i} |CT_P(o)| + \sum_{o_* \in X_*} |CT_P(o_*)| \right) = \frac{1}{|U|^2} \cdot \left(\sum_{i=1}^{l-1} \frac{n_i(n_i-1)}{2} + n_* \sum_{i=1}^{l-1} n_i + \frac{n_*(n_*-1)}{2} \right) = \frac{1}{|U|^2} \left(\sum_{i=1}^l \frac{n_i(n_i-1)}{2} + n_* (|U| - n_*) \right)$. Since $C_{n_i}^2 = \frac{n_i(n_i-1)}{2}$, $C_{n_*}^2 = \frac{n_*(n_*-1)}{2}$ and $|\overline{X_*}| = |U| - n_*$, we can get $EGK_U(P) = \frac{1}{|U|^2} \left(\sum_{i=1}^l C_{n_i}^2 + |X_*| |\overline{X_*}| \right) \square$

Property 1. Given an incomplete decision system $IDS = (U, C \cup D, V, f)$, $\forall P \subseteq C$. If the knowledge granularity of P is $GK_U(P)$ on U and the equal knowledge granularity of P is $EGK_U(P)$, then we have:

$$EGK_U(P) = (GK_U(P) - \frac{1}{|U|}) / 2 \quad (11)$$

Proof. Let $U/P = \{X_1, X_2, \dots, X_{l-1}, X_*\}$, $|X_i| = n_i$, $|X_*| = n_*$ where X_i is the i -th subdivision of U/P and X_* stands for the subdivision of missing values on attribute P , $|X_i|$ stands for the number of object in subdivision X_i . We can get $|U| - \sum_{i=1}^{l-1} |X_i| = |X_*|$. According to definition 2, suppose that $\forall o \in X_i$ and $T_P(o) = \{x | x \in X_i \vee x \in X_*\}$, we can get $|T_P(o)| = n_i + n_*$. Since the $|T_P(o)|$ value of each object in X_i is $n_i + n_*$, we can get $\sum_{o \in X_i} |T_P(o)| = n_i(n_i + n_*)$. In the same way, we can get $\sum_{o_* \in X_*} |T_P(o_*)| = n_* |U|$.

According to definition 3, we can get $GK_U(P) = \frac{1}{|U|^2} \left(\sum_{i=1}^{l-1} \sum_{o \in X_i} |T_P(o)| + \sum_{o_* \in X_*} |T_P(o_*)| \right) = \frac{1}{|U|^2} \left[\sum_{i=1}^{l-1} n_i(n_i + n_*) + n_* |U| \right]$, then

$$GK_U(P) - \frac{1}{|U|} = \frac{1}{|U|^2} \left(\sum_{i=1}^{l-1} (n_i^2 - n_i) + \sum_{i=1}^{l-1} n_i + n_* \sum_{i=1}^{l-1} n_i + n_* |U| - |U| \right) = \frac{1}{|U|^2} \left(\sum_{i=1}^{l-1} (n_i^2 - n_i) - n_*^2 - n_* + 2n_* |U| \right)$$

$$|U| = \frac{2}{|U|^2} \left(\sum_{i=1}^{l-1} \frac{(n_i^2 - n_i)}{2} + \frac{n_*^2 - n_*}{2} + n_* (|U| - n_*) \right) = 2EGK_U(P) \cdot |U|$$

$$\frac{2}{|U|^2} \left(\sum_{i=1}^l \frac{(n_i^2 - n_i)}{2} + n_* (|U| - n_*) \right) = 2EGK_U(P) \square$$

Due to the time complexity of calculating $T_P(o)$ is $|U|$, we know that the time complexity of calculating $GK_U(P)$ is $|U|^2$. However, the time complexity of $CT_P(o)$ is $|U/P|$, the time complexity of calculating $EGK_U(P)$ is $|U/P|^2$ and $|U/P|^2 \ll |U|^2$. In addition, in the process of calculating $EGK_U(P)$, the sub-division with a cardinality of 1 is constantly pruned, which further speeds up the calculation. Therefore, the time of calculating $EGK_U(P)$ is less than $GK_U(P)$ for the same data set.

Example 1. Example of computing equivalent knowledgegranularity. Let $IDS = (U, C \cup D, V, f)$, $U = \{o_1, o_2, o_3, o_4, o_5, o_6, o_7, o_8, o_9\}$, $C = \{a, b, c, e\}$, $D = \{d\}$. The detail data shown in Table 1. Let $P = \{a, b\}$, we can get $f(o_6, P) = *$, $f(o_9, P) = *$. We use the following two methods to calculate $EGK_U(P)$.

1) According to definition 5, we obtain that $CT_P(o_1) = \{o_2, o_6, o_9\}$, $CT_P(o_2) = \{o_6, o_9\}$, $CT_P(o_3) = \{o_4, o_5, o_6, o_9\}$, $CT_P(o_4) = \{o_5, o_6, o_9\}$, $CT_P(o_5) = \{o_6, o_9\}$, $CT_P(o_6) = \{o_9\}$, $CT_P(o_7) = \{o_6, o_8, o_9\}$, $CT_P(o_8) = \{o_6, o_9\}$, $CT_P(o_9) = \{\emptyset\}$. According to definition 6, we can get $EGK_U(P) = \frac{1}{|U|^2}$.

$$\sum_{i=1}^9 |CT_P(o_i)| \cdot \frac{1}{9^2} (3+2+4+3+2+3+2+1+0) = \frac{20}{81}.$$

2) Since $U/P = \{\{o_1, o_2\}, \{o_3, o_4, o_5\}, \{o_7, o_8\}, \{o_6, o_9\}\}$, let $X_1 = \{o_1, o_2\}$, $X_2 = \{o_3, o_4, o_5\}$, $X_3 = \{o_7, o_8\}$, $X_* = \{o_6, o_9\}$, then $|X_*| = 2$, $|\overline{X_*}| = |U| - 2 = 7$. According to theorem 1, we can get $EGK_U(P) = \frac{1}{9^2} \cdot [C_2^2 + C_3^2 + C_2^2 + C_2^2 + 2 \cdot (9-2)] = \frac{20}{81}$.

Although the above two methods get the same results, the calculation time is different. Since method 1 needs to scan the data set multiple times, it consumes more time. However, method 2 only needs to scan the data set one time and obtain each subdivision of U/P . According to the number of objects in each subdivision, we can get the combination value quickly, then the value of equivalent knowledge granularity is calculated.

3. The mechanism of knowledge granularity attribute reduction in the background of clustering

Most of traditional attribute reduction models use equivalence class relation to compute the importance of conditional attribute. Although these methods can effectively deal with complete decision-making information systems, they can not obtain correct reduction rules in incomplete ones. In order to deal with the loss of information effectively, this paper focuses on the reduction of incomplete decision systems.

Table 1. Incomplete information system.

U	a	b	c	e	d
O1	0	0	0	1	0
O2	0	0	1	*	1
O3	0	1	0	1	0
O4	0	1	*	0	1
O5	0	1	0	*	1
O6	*	*	1	1	1
O7	1	0	*	0	2
O8	1	0	1	0	2
O9	*	*	0	1	2

The traditional reduction model does not consider the impact on the classification of the data set when deleting redundant attributes. If there are inconsistent objects in the data set, the classification accuracy of the reduced set will be affected. In order to improve the data quality, this paper uses the idea of clustering. Clustering is to divide all objects in the data set into different clusters according to a certain standard when the target category is unknown. Objects within a cluster are as similar as possible, and objects between clusters are as dissimilar as possible. Classification is to classify all objects in the data set according to a certain nature and level when the object category is known. Good clustering results can be used as a reference standard for accurate classification. The desired results of clustering are the objects of the same class are gathered in intra-clustering, otherwise they will be gathered in different inter-clustering. This paper studies the labeled data objects decision information system. Therefore, we use the results of the classification to guide the process of clustering the data objects. When the data objects are clustered, they follow the principle that the objects of intra-

clustering are as close as possible and the objects of inter-clustering are as far away as possible. Next, we discuss how to measure the distance of intra-clustering and inter-clustering objects.

3.1. The intra-cluster similarity for incomplete systems

Generally, there are two approaches for clustering calculations: distance and similarity. The closer is the distance between two different objects, the weaker is their ability to distinguish. On the contrary, the farther is the distance, the stronger is the ability to distinguish. In this paper, similarity method is used to measure the distinguish ability of objects. Since the knowledge granularity can measure the similarity between objects, the coarser is the knowledge granularity, the stronger is the distinguishing ability. The better is the knowledge granularity, the weaker is the distinguishing ability. Next, we discuss how to use knowledge granularity information to measure the similarity of objects in an incomplete system.

Definition 7 (The similarity of intra-cluster objects). Given an incomplete decision system $IDS = (U, C \cup D, V, f)$, $U / D = \{D_1, D_2, \dots, D_n\}$. For sake of convenience, let $U / D = \pi_D$, $D_i \subseteq \pi_D$, $P \subseteq C$. Suppose the equivalence division relationship of D_i under attribute set P is $\mathbb{R}_P = \{X_1, X_2, \dots, X_m\}$, the similarity of objects in the cluster of D_i about attribute P is defined as following (Where $o \in D_i$),

$$SIntra_{D_i}(P) = EGK_{D_i}(P) = \frac{1}{|D_i|^2} \sum_{o \in D_i} |CT_P(o)| \quad (12)$$

Definition 8 (The average similarity of intra-cluster objects). Given an incomplete decision systems $IDS = (U, C \cup D, V, f)$, $U / D = \{D_1, D_2, \dots, D_n\}$. $P \subseteq C$, $o \in D_i$. The Knowledge granularity similarity of intra-clustering objects for subdivision D_i with respect to attribute P is $SIntra_{D_i}(P)$, then the average intra-clustering similarity is defined as following,

$$ASIntra_{\pi_D}(P) = \frac{1}{n} \sum_{i=1}^n SIntra_{D_i}(P) \quad (13)$$

The desired effect of clustering is that the similarity of intra-clustering is high, and the similarity of inter-clustering is low.

Property 2. Given an incomplete system $IDS = (U, C \cup D, V, f)$, $U / D = \pi_D$, $D_i \subseteq \pi_D$, $P, Q \subseteq C$. If $P \subseteq Q$, we have

$$ASIntra_{\pi_D}(P) \geq ASIntra_{\pi_D}(Q) \quad (14)$$

Proof. Let $D_i / P = \{X_1, X_2, \dots, X_k \cup X_{k+1}, X_{k+2}, \dots, X_n, X_* \cup Y\}$, where X_*, Y represents the object sets with missing values on attribute set P . Since $P \subseteq Q$, we can get $D_i / Q \subseteq D_i / P$. This to say, each subdivision of D_i / Q is a subset of some subdivision of D_i / P . Let $D_i / Q = \{X_1, X_2, \dots, X_k, X_{k+1}, X_{k+2}, \dots, X_n, Y, X_*\}$. According to theorem and definition 7, we can get

$$\begin{aligned} SIntra_{D_i}(P) &= \frac{2}{|D_i|^2} \left(\sum_{j=1}^{k-1} C_{|X_j|}^2 + C_{|X_k|+|X_{k+1}|}^2 + \sum_{j=k+2}^n C_{|X_j|}^2 + C_{|X_*|+|Y|}^2 + (|X_*|+|Y|)(|D_i|-|X_*|-|Y|) \right) \\ &= \frac{2}{|D_i|^2} \left(\sum_{j=1}^{k-1} C_{|X_j|}^2 + C_{|X_k|}^2 + C_{|X_{k+1}|}^2 + |X_k||X_{k+1}| + \sum_{j=k+2}^n C_{|X_j|}^2 + C_{|X_*|}^2 + C_{|Y|}^2 + |X_*||Y| + (|X_*|+|Y|)(|D_i|-|X_*|-|Y|) \right) \\ &\quad - |Y|) = \frac{2}{|D_i|^2} \left(\sum_{j=1}^n C_{|X_j|}^2 + |X_k||X_{k+1}| + C_{|X_*|}^2 + C_{|Y|}^2 + |X_*||Y| + (|X_*|+|Y|)(|D_i|-|X_*|-|Y|) \right). \quad \text{Since} \\ SIntra_{D_i}(Q) &= \frac{2}{|D_i|^2} \left(\sum_{j=1}^n C_{|X_j|}^2 + C_{|X_*|}^2 + C_{|Y|}^2 + |X_*|(|D_i|-|X_*|) \right), \quad \text{then} \quad \frac{2}{|D_i|^2} (|X_k||X_{k+1}| + |Y||D_i|-|X_*||Y|-|Y|^2) = \frac{2}{|D_i|^2} (|X_k||X_{k+1}| + |Y|(|D_i|-|X_*|-|Y|)). \quad \text{Since } |D_i| \geq |X_*|+|Y|, |X_k||X_{k+1}| \geq 0 \text{ So, the results} \end{aligned}$$

of $SIntra_{D_i}(P) \geq SIntra_{D_i}(Q)$ and $\sum_{i=1}^n SIntra_{D_i}(P) \geq \sum_{i=1}^n SIntra_{D_i}(Q)$ are obtained. Above all, we can get $ASIntra_{\pi_D}(P) \geq ASIntra_{\pi_D}(Q)$.

According to property 2, we conclude that the intra-cluster similarity is monotonic when the conditional attribute set changes.

Example 2. In Table 1, let $P = \{a, b\}$, we have $D_1 = \{o_1, o_3\}$, $D_2 = \{o_2, o_4, o_5, o_6\}$, $D_3 = \{o_7, o_8, o_9\}$, $SIntra_{D_1}(P) = 0$, $SIntra_{D_2}(P) = \frac{0+C_2^2+1 \times 3}{4^2} = \frac{1}{4}$, $SIntra_{D_3}(P) = \frac{C_2^2+1 \times 2}{3^2} = \frac{1}{3}$, $ASIntra_{\pi_D}(P) = \frac{1}{3} \sum_{i=1}^3 SIntra_{D_i}(P) = \frac{7}{36}$.

3.2. The inter-cluster similarity for incomplete systems

Definition 9. (The inter-cluster similarity for incomplete systems) Given $IDS = (U, C \cup D, V, f)$ be an incomplete decision system, Let $\pi_D = \{D_1, D_2, \dots, D_n\}$. Suppose $P \subseteq C$, $D_i, D_j \subset U / D$, then the inter-cluster similarity of D_i and D_j respect of attribute set P for incomplete systems is defined as following,

$$SInter_{D_i, D_j}(P) = \frac{1}{(|D_i| + |D_j|)^2} \left(\sum_{o \in D_i \cup D_j} |CT_P(o)| - \sum_{o \in D_i} |CT_P(o)| - \sum_{o \in D_j} |CT_P(o)| \right) \quad (15)$$

Assuming that D_i and D_j are objects of two different clusters, the inter-cluster similarity between D_i and D_j is calculated in two steps. The first step is to calculate the similarity after the two clusters are merged, and the second step is to remove the similarity information of the same cluster. The rest is the similarity information between objects in different clusters.

Property 3. Given an incomplete decision system $IDS = (U, C \cup D, V, f)$, $\pi_D = \{D_1, D_2, \dots, D_n\}$, $P \subseteq C$. Let $D_i / P = \{X_1, X_2, \dots, X_k, X_{k+1}, \dots, X_m, X_*\}$, $\overline{X_*} = D_i - X_*$, $D_j / P = \{Y_1, Y_2, \dots, Y_k, Y_{k+1}, \dots, Y_n, Y_*\}$, $\overline{Y_*} = D_j - Y_*$, $D_i \cup D_j / P = \{X_1 \cup Y_1, X_2 \cup Y_2, \dots, X_k \cup Y_k, X_* \cup Y_*, X_{k+1}, \dots, X_m, Y_{k+1}, \dots, Y_n\}$, where $\forall o \in X_* \cup Y_*, f(o, P) = *$, $*$ is the flag of missing value. We have :

$$SInter_{D_i, D_j}(P) = \frac{1}{(|D_i| + |D_j|)^2} \left(\sum_{l=1}^k |X_l| |Y_l| + |X_*| |Y_*| + |X_*| |\overline{Y_*}| + |\overline{X_*}| |Y_*| \right) \quad (16)$$

Proof. According definition 5 and theorem 1, we can get $\sum_{o \in D_i \cup D_j} |CT_P(o)| = \sum_{l=1}^k C_{|X_l|+|Y_l|}^2 + C_{|X_*|+|Y_*|}^2 + \sum_{l=k+1}^m C_{|X_l|}^2 + \sum_{l=k+1}^n C_{|Y_l|}^2 + (|X_*| + |Y_*|)(|\overline{Y_*}| + |\overline{X_*}|)$, $\sum_{o \in D_i} |CT_P(o)| = \sum_{l=1}^m C_{|X_l|}^2 + C_{|X_*|}^2 + |X_*| |\overline{X_*}|$, $\sum_{o \in D_j} |CT_P(o)| = \sum_{l=1}^n C_{|Y_l|}^2 + C_{|Y_*|}^2 + |Y_*| |\overline{Y_*}|$. Since $\sum_{l=1}^k C_{|X_l|+|Y_l|}^2 = \sum_{l=1}^k \left(\frac{|X_l|^2 - |X_l| + |Y_l|^2 - |Y_l|}{2} + |X_l| |Y_l| \right) = \sum_{l=1}^k C_{|X_l|}^2 + \sum_{l=1}^k C_{|Y_l|}^2 + \sum_{l=1}^k |X_l| |Y_l|$, according definition 9, we can conclude that $(|D_i| + |D_j|)^2 SInter_{D_i, D_j}(P) = \sum_{o \in D_i \cup D_j} |CT_P(o)| - \sum_{o \in D_i} |CT_P(o)| - \sum_{o \in D_j} |CT_P(o)| =$

$$\sum_{o \in D_i \cup D_j} |CT_P(o)| = \sum_{l=1}^k |X_l| |Y_l| + |X_*| |Y_*| + |X_*| |\overline{Y_*}| + |\overline{X_*}| |Y_*| \quad . \quad \text{Then we have}$$

$$SInter_{D_i, D_j}(P) = \frac{1}{(|D_i| + |D_j|)^2} \left(\sum_{l=1}^k |X_l| |Y_l| + |X_*| |Y_*| + |X_*| |\overline{Y_*}| + |\overline{X_*}| |Y_*| \right) \square$$

Property 4. Given an incomplete decision system $IDS = (U, C \cup D, V, f)$, $\pi_D = \{D_1, D_2, \dots, D_n\}$, $P \subseteq Q \subseteq C$, $D_i, D_j \subset U / D$, then we have $SInter_{D_i, D_j}(P) \geq SInter_{D_i, D_j}(Q)$.

Proof. Let $D_i / Q = \{X_1, X_2, \dots, X_k, X_{k+1}, \dots, X_m, X_\Delta, X_* - X_\Delta\}$, $\overline{X_* - X_\Delta} = D_i - X_* + X_\Delta$, $D_j / Q = \{Y_1, Y_2, \dots, Y_k, Y_{k+1}, \dots, Y_n, Y_\Delta, Y_* - Y_\Delta\}$, $\overline{Y_* - Y_\Delta} = D_j - Y_* + Y_\Delta$. Since $P \subseteq Q \subseteq C$, then D_i / Q is a refinement of D_i / P , D_j / Q is a refinement of D_j / P . Let $D_i / P = \{X_1, X_2, \dots, X_k \cup X_{k+1}, \dots, X_m, X_*\}$, $\overline{X_*} = D_i - X_*$, $D_j / P = \{Y_1, Y_2, \dots, Y_k \cup Y_{k+1}, \dots, Y_n, Y_*\}$, $\overline{Y_*} = D_j - Y_*$. Suppose $D_i \cup D_j / Q = \{X_1 \cup Y_1, X_2 \cup Y_2, \dots, X_k \cup Y_k, X_{k+1} \cup Y_{k+1}, X_\Delta \cup Y_\Delta, (X_* - X_\Delta) \cup (Y_* - Y_\Delta), X_{k+2}, \dots, X_m, Y_{k+2}, \dots, Y_n\}$, $D_i \cup D_j / P = \{X_1 \cup Y_1, X_2 \cup Y_2, \dots, X_k \cup Y_k, X_{k+1} \cup Y_{k+1}, X_* \cup Y_*, X_{k+2}, \dots, X_m, Y_{k+2}, \dots, Y_n\}$, then

$$\begin{aligned} SInter_{D_i, D_j}(P) &= \sum_{l=1}^{k-1} |X_l| |Y_l| + (|X_k| + |X_{k+1}|)(|Y_k| + |Y_{k+1}|) + |X_*| |Y_*| + |X_*| |\overline{Y_*}| + |\overline{X_*}| |Y_*|, SInter_{D_i, D_j}(Q) \\ &= \sum_{l=1}^{k-1} |X_l| |Y_l| + |X_k| |Y_k| + |X_{k+1}| |Y_{k+1}| + |X_\Delta| |Y_\Delta| + |X - X_\Delta| |Y - Y_\Delta| + |X_* - X_\Delta| |\overline{Y_* - Y_\Delta}| + |\overline{X_* - X_\Delta}| |Y_* - Y_\Delta|. \\ SInter_{D_i, D_j}(P) - SInter_{D_i, D_j}(Q) &= |X_k| |Y_{k+1}| + |X_{k+1}| |Y_k| + |X_*| |Y_*| + |X_*| |\overline{Y_*}| + |\overline{X_*}| |Y_*| - |X_\Delta| |Y_\Delta| \\ &- |X - X_\Delta| |Y - Y_\Delta| - |X_* - X_\Delta| |\overline{Y_* - Y_\Delta}| + |\overline{X_* - X_\Delta}| |Y_* - Y_\Delta|. \text{ Since } |\overline{X_*}| = |D_i| - |X_*|, |\overline{Y_*}| = |D_j| - |Y_*|, \\ &|\overline{X_* - X_\Delta}| = |D_i| - |X_*| + |X_\Delta|, |\overline{Y_* - Y_\Delta}| = |D_j| - |Y_*| + |Y_\Delta|, |X_k| |Y_{k+1}| + |X_{k+1}| |Y_k| \geq 0, \text{ then } SInter_{D_i, D_j}(P) \\ &- SInter_{D_i, D_j}(Q) \geq |X_*| |Y_*| + |X_*| |\overline{Y_*}| + |\overline{X_*}| |Y_*| - |X_\Delta| |Y_\Delta| - |X - X_\Delta| |Y - Y_\Delta| - |X_* - X_\Delta| |\overline{Y_* - Y_\Delta}| + |\overline{X_* - X_\Delta}| |Y_* - Y_\Delta| \\ &= |X_\Delta| (|D_j| + |Y_\Delta| |D_i| - |X_*| |Y_\Delta| - |X_\Delta| |Y_*|) = |X_\Delta| (|D_j| - |Y_*|) + |Y_\Delta| (|D_i| - |X_*|). \end{aligned}$$

Since $|D_j| \geq |Y_*|$, $|D_i| \geq |X_*|$, then we have $SInter_{D_i, D_j}(P) - SInter_{D_i, D_j}(Q) \geq 0$.

From property 4, it can be concluded that the similarity of inter-clusters has mono-tonicity with the change of conditional attributes.

Definition 10 (The average similarity of inter-cluster objects). Given an incomplete decision system $IDS = (U, C \cup D, V, f)$, $\pi_D = \{D_1, D_2, \dots, D_n\}$, $P \subseteq C$. In n different clusters, the inter-cluster similarity is calculated between every two clusters, and the time of comparisons is $\frac{1}{2}n(n-1)$, then the average knowledge granularity similarity of inter-cluster in data set U respect of attribute set P is defined as following,

$$ASInter_{\pi_D}(P) = \frac{2}{n(n-1)} \sum_{i=1}^{n-1} \sum_{j=i+1}^n SInter_{D_i, D_j}(P) \quad (17)$$

Example 3. In Table 1, the same conditions as example 2, let $P = \{a, b\}$, we have $D_1 = \{o_1, o_3\}$, $D_2 = \{o_2, o_4, o_5, o_6\}$, $D_3 = \{o_7, o_8, o_9\}$, $SInter_{D_1, D_2}(P) = \frac{1 \times 1 + 1 \times 2 + 2 \times 1}{(2+4)^2} = \frac{5}{36}$, $SInter_{D_1, D_3}(P) = \frac{0+0+1 \times 2}{(2+3)^2} = \frac{2}{25}$, $SInter_{D_2, D_3}(P) = \frac{0+1 \times 2 + 3 \times 1}{(4+3)^2} = \frac{5}{49}$, $ASInter_{\pi_D}(P) = \frac{1}{3} \sum_{i=1}^2 \sum_{j=i+1}^3 SInter_{D_i, D_j}(P) = \frac{14153}{132300}$.

4. Attribute reduction of knowledge granularity for incomplete systems

Traditional attribute reduction methods are mostly aimed at data sets with no missing data. Various data sets in reality are often incomplete due to various subjective or objective factors. Therefore, we research information systems with missing data and propose corresponding algorithms to improve the data quality of incomplete system reduction sets.

This paper discusses how to design a reduction method with the idea of clustering. For an ideal clustering effect, the objects of inter-cluster should be far away, and the objects of intra-cluster should be close together. Here, similarity is used to measure the distance between two different objects. The higher is the similarity, the closer are the objects. Conversely, the lower is the similarity, the farther are the objects. Based on the above analysis, we design a formula to measure the importance of attributes as following,

$$SIM_R = SAIntra + \lambda \cdot (1 - ASInter) \quad (18)$$

Where λ is the weight, $1 - ASIntra$ is the dissimilarity of intra-cluster objects. We can set the importance of the intra-cluster similarity and inter-cluster similarity by using the size of the

parameter λ . This method needs to adjust the parameters continuously, which consumes a lot of time. To this end, we first normalize $ASIntra$ and $ASInter$. Then, the two similarity calculations can be measured within a unified range, avoiding the parameter adjustment process.

4.1. Normalization of inter-cluster similarity and intra-cluster similarity

Given an incomplete decision system $IDS = (U, C \cup D, V, f), \pi_D = \{D_1, D_2, \dots, D_n\}, P \subseteq C, D_i, D_j \subset U/D$. Since the number of elements in each sub-division may be different, then the value range of its equivalent knowledge granularity may also be different. To calculate the average similarity of all subdivisions, it must be calculated in the same domain. For the sake of generality, we normalize the inter-cluster similarity and intra-cluster similarity. According to definition 7 and theorem 1, we can get $SIntra_{D_i}(P) = EGK_{D_i}(P)$. When all data objects in the subdivision D_i are indistinguishable with respect to the attribute set P , $EGK_{D_i}(P) = \frac{1}{2}(1 - \frac{1}{|D_i|})$ takes the maximum value.

When all data objects in D_i can be distinguished from each other, $EGK_{D_i}(P) = 0$ takes the minimum value. So, the result of $EGK_{D_i}(P) \in \left[0, \frac{|D_i|-1}{2|D_i|}\right]$ is obtained. If the value of $EGK_{D_i}(P)$ is mapped to the range $[0,1]$, the correction formula of $SIntra_{D_i}(P)$ is defined as

$$SIntra_{D_i}(P)' = EGK_{D_i}(P) / \left(\frac{|D_i|-1}{2|D_i|} \right) = EGK_{D_i}(P) \frac{2|D_i|}{|D_i|-1} \quad (19)$$

The average similarity of intra-cluster objects is corrected as following,

$$ASIntra_{\pi_D}(P) = \frac{1}{n} \sum_{i=1}^n SIntra_{D_i}(P)' \quad (20)$$

D_i and D_j are two object sets of different clusters. Suppose $D_i/P = \{X_1, X_2, \dots, X_n, X_*\}$, $D_j/P = \{Y_1, Y_2, \dots, Y_m, Y_*\}$. If $f(X_l, P) = f(Y_l, P) (1 \leq l \leq k)$, then $D_i \cup D_j/P = \{X_1 \cup Y_1, X_2 \cup Y_2, \dots, X_k \cup Y_k, \dots, X_n, Y_{k+1}, \dots, Y_m, X_*, Y_*\}$. According to property 3, the similarity of inter-cluster respect to D_i and D_j denotes as $SInter_{D_i, D_j}(P) = \frac{1}{(|D_i|+|D_j|)^2} \left(\sum_{l=1}^k |X_l||Y_l| + |X_*||Y_*| + |\overline{X_*}||\overline{Y_*}| \right)$.

When $\bigcup_{i=1}^k X_i \cup Y_i = \emptyset$, $X_* = \emptyset, Y_* = \emptyset$, $SInter_{D_i, D_j}(P) = 0$, takes minimum value. When all data objects in D_i and D_j are indistinguishable, $SInter_{D_i, D_j}(P) = \frac{|D_i||D_j|}{(|D_i|+|D_j|)^2}$ takes maximum value. Then, we can get $SInter_{D_i, D_j}(P) \in \left[0, \frac{|D_i||D_j|}{(|D_i|+|D_j|)^2}\right]$. The normalized formula of $SInter_{D_i, D_j}(P)$ is as following,

$$SInter_{D_i, D_j}(P)' = SInter_{D_i, D_j}(P) \frac{(|D_i|+|D_j|)^2}{|D_i||D_j|} \quad (21)$$

The definition of the average similarity of inter-cluster objects is revised as following,

$$ASInter_{\pi_D}(P) = \frac{n(n-1)}{2} \sum_{i=1}^{n-1} \sum_{j=i+1}^n SInter_{D_i, D_j}(P)' \quad (22)$$

After the similarities of inter-cluster and intra-cluster objects are normalized $SIM_R = ASInter + \lambda \cdot (1 - ASIntra)$ is revised as following,

$$SIM_R = AIntra_{\pi_D}(P) + 1 - AInter_{\pi_D}(P) \quad (23)$$

Since $ASInter_{\pi_D}(P)$ represents the similarity of intra-cluster objects, then the dissimilarity is $1 - ASInter_{\pi_D}(P)$. If you use formula of SIM_R to measure the effect of clustering, the larger is the value of SIM_R , the better is the effect.

4.2. The knowledge granularity attribute reduction algorithm for incomplete systems(IKAR)

In section IV.A, we discussed the similarities of inter-cluster and intra-cluster objects from the perspective of clustering, which provided a clear goal for the next step of attribute selection.

Definition 11 (Equal knowledge granularity attribute reduction). Given an incomplete decision system $IDS = (U, C \cup D, V, f)$, an attribute subset $R \subseteq C$ is an equal knowledge granularity attribute reduction if and only if:

$$\begin{aligned} 1) \quad & R = \min_{P \subseteq C} \{SIM_P\} \\ 2) \quad & \forall R' \subset R, SIM_{R'} > SIM_R \end{aligned} \quad (24)$$

In this definition 11, condition (1) is the jointly sufficient condition that guarantees the equal knowledge granularity value induced from the reduction is minimal, and condition (2) is the individual necessary condition that guarantees the reduction is minimal.

According to definition 11, we can find a smaller reduction set with an ideal classification effect. Property 2 proves that when the attribute decreases, the similarity of intra-cluster increases monotonically. Property 4 proves that the similarity of inter-cluster also increases when decreasing the attribute, so that the dissimilarity will decrease. Obviously, the formula SIM_R cannot determine its monotonicity. To find the R when the SIM_R is the largest in the conditional attribute C is a combinatorial optimization problem, and trying the methods one by one is not the best way to solve the problem, so we use a heuristic method to find the smallest reduced set. For any attribute $a \in C$, its inner significance is defined as follows:

$$Sig_C^a = SIM_C - SIM_{C-a} \quad (25)$$

The bigger is the value of Sig_C^a , the more important is the attribute.

In order to get the optimal reduction set quickly, we adopt the deletion strategy. Firstly, the importance Sig_C^a of attribute a is defined by the formula of SIM , then sort the different Sig_C^a . Secondly, let $R=C$, the value of SIM_C is used as the initial condition, which ensures that the clustering effect after reduction is better than the raw data set. Then, remove the unimportant attribute a from the remaining attributes $C - R$ and calculate the value of SIM_{R-a} . If $SIM_{R-a} \geq SIM_R$, delete the attribute a and continue, otherwise the algorithm terminates. The detail of IKAR algorithm is shown in Figure 1.

Example 4. In Table 1, Since $D_1 = \{o_1, o_3\}$, $D_2 = \{o_2, o_4, o_5, o_6\}$, $D_3 = \{o_7, o_8, o_9\}$. Let $R = C$, according to the definition of SIM_R , we can get $SIM_R = \frac{72}{75}$, $SIM_{C-a} = \frac{71}{72}$, $SIM_{C-b} = \frac{100}{72}$, $SIM_{C-c} = \frac{60}{72}$ and $SIM_{C-e} = \frac{80}{72}$, then $Sig_C^a = \frac{4}{72}$, $Sig_C^b = -\frac{25}{72}$, $Sig_C^c = \frac{15}{72}$, $Sig_C^e = -\frac{5}{72}$. Since $SIM_{C-b} = \frac{100}{72} > SIM_R$, we delete the attribute b from R and let $R = R - b$, $SIM_R = \frac{100}{72}$. In the same way, we obtain that $SIM_{R-a} = \frac{94}{72}$, $SIM_{R-c} = \frac{92}{72}$, $SIM_{R-e} = \frac{108}{72}$, $Sig_R^a = \frac{6}{72}$, $Sig_R^c = \frac{8}{72}$ and $Sig_R^e = -\frac{8}{72}$. Since $SIM_{R-e} = \frac{108}{72} > SIM_R$, we delete the attribute e from R and let $R = R - e$. We calculate the value of SIM_{R-a} , SIM_{R-c} , get the results of $SIM_{R-a} = \frac{82}{72}$ and $SIM_{R-c} = \frac{100}{72}$. Now, we have $Sig_R^a = \frac{18}{72}$ and $Sig_R^c = \frac{8}{72}$. If the attribute c is deleted $SIM_{R-c} = \frac{100}{72} < SIM_R$, algorithm is terminated. We have $R = \{a, c\}$.

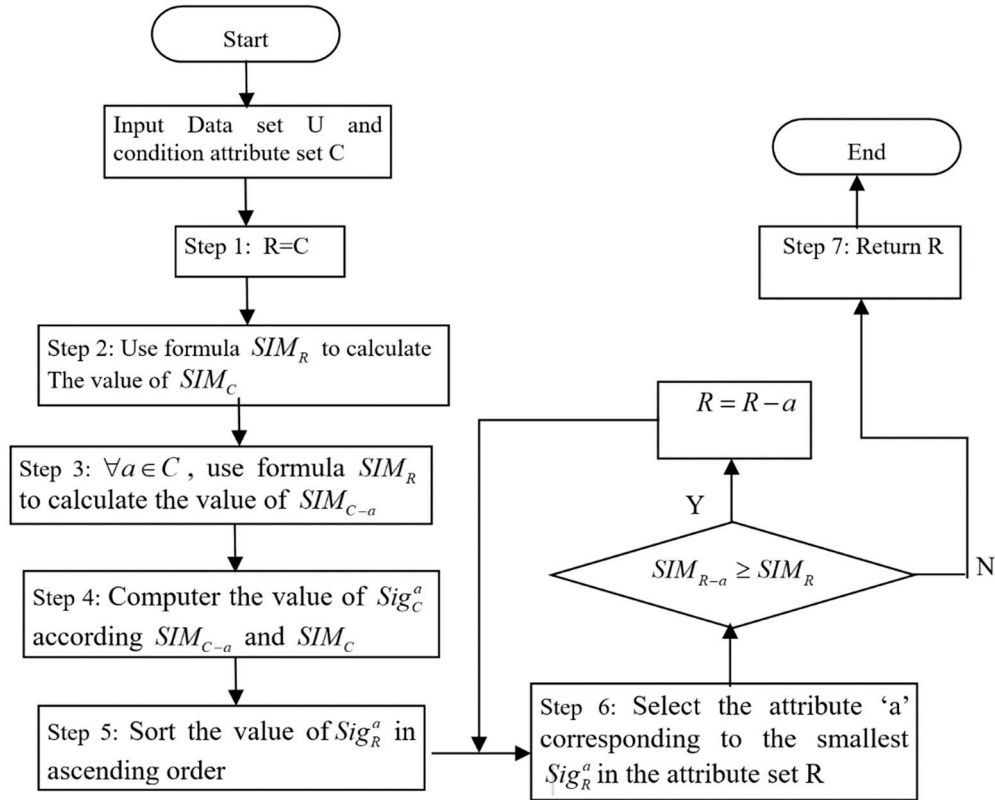


Figure 1. The execution process of algorithm (IKAR).

4.3. Time complexity analysis

The time complexity of Step 1 is $O(1)$. In Step 2, we calculate the value of SIM_C which include $ASIntra_{\pi_D}(C)$ and $ASInter_{\pi_D}(C)$. The computational time complexity of intra-cluster $SIntra_{D_i}(C)$ similarity is $O(|C||D_i|)$, then the time complexity of calculating $ASIntra_{\pi_D}(C)$ is $O(|U||C|)$. Since the time complexity of calculating the similarity $SInter_{D_i, D_j}(C)$ about inter-cluster D_i and D_j is $O((|D_i| + |D_j|)|C|)$, then time complexity of $ASInter_{\pi_D}(C)$ is $O(\sum_{i=1}^{|U/D|-1} \sum_{j=i+1}^{|U/D|} (|D_i| + |D_j|)|C|) = O((|U/D|-1)|U||C|)$. We can get the time complexity of Step 2 is $O(|U/D||U||C|)$; In Step 3, we consume time is $O(|U/D||U||C|^2)$. Since Step 4 utilizes the results of Step 3, the time complexity is $O(1)$. Step 5 is to sort the importance Sig_R^a of each attribute, and the time complexity is $O(\frac{1}{2}|C|^2)$. Since the time complexity of calculating Sig_R^a is $O(|U/D||U||R|)$, then the time complexity of Step 5 is $O(|U/D||U||R|) + O(\frac{1}{2}|C|^2)$. In Step 6, the time complexity of deleting a redundant attribute is $O(|U/D||U||R|^2)$, and Step 6 needs to be executed $|C| - |R|$ times, $R \subseteq C$, then the time complexity of Step 6 is $O(|U/D||U|(|C|^3 - |R|^3))$; In summary, the time complexity of IKAR is $\max\{O(|U/D||U||C|^2), O(|U/D||U|(|C|^3 - |R|^3))\}$.

5. Experiments results analysis

In order to evaluate the feasibility and effectiveness of the proposed algorithm in this paper, the complete data set is preprocessed to obtain incomplete information systems, and many different attribute reduction algorithms are used for reduction. The reduction set obtained in the previous stage is classified and analyzed by multiple classifiers in the Weka Tool. It is compared with other

three existing algorithms in terms of reduction set size, running time, accuracy and algorithm stability. The specific experimental framework is shown in Figure 2.

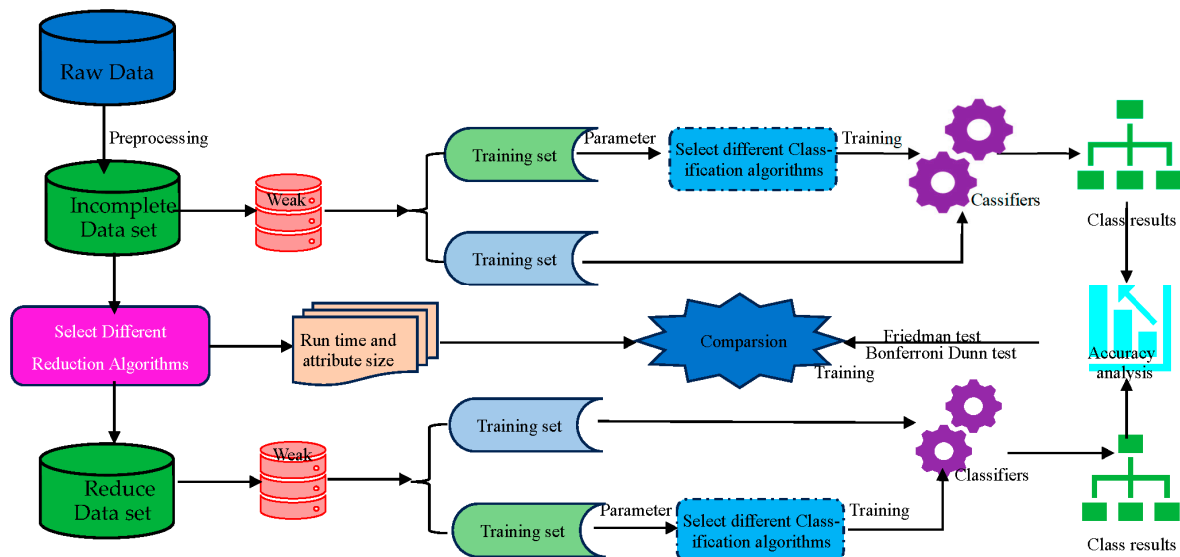


Figure 2. The frame work chart of experiments.

All of the datasets are displayed in Table 2. The twelve datasets selected were downloaded from UCI. In Table 2, $|U|$, $|C|$ and $|D|$ represents the number of objects, conditional attributes and the categories of decision attribute respectively. In order to generate an incomplete information system, we delete 12% attribute values from the raw datasets, and the missing values are randomly and uniformly distributed. The missing value of each attribute is removed with equal probability, which eliminates the impact on the classification accuracy of the later reduction set due to different attribute selection. The three classifiers J48, SMO of the weka 3.8 software platform are used to demonstrate the classification effect of the reduction set. In subsequent experiments, the reduced sets of each dataset were analyzed for accuracy using the cross method. All the objects in the reduced set are randomly divided into 10 equal parts, one of which is used as the test set, and the remaining 9 are used as the training set. In this way, each reduction set is repeated 10 times for the cross experiment. Finally, the size of reduction set, running time and classification accuracy obtained by the 10 experiments are averaged. We execute all experiments on a PC with Windows 10, Intel(R) Core(TM) i7-10710U CPU @ 1.10GHz, 1.61 GHz and 8GB memory. Algorithms are code in python and the software being used is PyCharm Community Edition 2020.2.3 x64.

Table 2. Description of twelve data sets.

ID	Data sets	Abbreviation	$ U $	$ C $	$ D $
1	Promoters	Prom	106	57	2
2	Heart-statlog	Hear	270	13	2
3	Hepatitis	Hepa	155	19	2
4	HandWritten	Hand	5620	64	10
5	Chess kr-kp	Ches	3196	36	2
6	Splice	Spli	3190	61	3
7	Letters	Lett	20000	17	26
8	Vote	Vote	435	16	2
9	Mushroom	Mush	8124	22	2
10	Qsar	Qsar	1055	42	2
11	Shuttle	Shut	43500	9	7
12	Satimage	Sati	6435	36	6

5.1. Reduction set size and running time analysis

This section mainly verifies the feasibility of IKAR for data set reduction from perspective of reduction size and calculation time. Here, we have selected other three representative incomplete system attribute reduction approaches. For the convenience of the following description, these three methods are referred to as NGLE[32], IEAR[33] and PARA[34] respectively. NGLE is the neighborhood multi-granulation rough sets-based attribute reduction using Lebesgue and entropy, and the method considers both algebraic view and information view. IEAR is the information entropy attribute reduction for incomplete set-valued data using the similarity degree function, and proposed λ -information entropy. PRAR is the positive region attribute reduction using indiscernibility and discernibility relation.

The attribute reduction size of four algorithms is shown in Table 3. Table 4 shows the time consumed by the reduction process of the four algorithms in seconds. Ave represents the average value, Best in Table 3 represents the number of minimum reduction sets obtained, and Best in Table 4 stands for the number of times which the running time is the shortest. From Table 3, it can be seen that the average reduction set size of the IKAR algorithm is 11.833, and the average reduction set size obtained by the NGLE, LEAR and PARA algorithms are 11.917, 11.00 and 12.083, respectively. IKAR obtained the minimum reduction set on the Ches, Spli, Mush and Shut data sets, the reduction effect is slightly better than the NGLE, PARA, but not so ideal as the LEAR algorithm. In the 12 data sets in Table 2, the LEAR algorithm obtains the minimum reduction set in 11 data sets.

From the consumption time of Table 4, we can find that the reduction advantage obtained by the IKAR algorithm is not obvious, but the IKAR algorithm is obviously better than the other three algorithms. When calculating the reduction set of 12 data sets, the average time required by IKAR is 15.20 seconds, the average time consumed by the other three algorithms are (155.40, 2003.56, 260.38) seconds respectively. From the experimental results in Table 4, it can also be found that the IKAR algorithm only needs 25.30 seconds to reduce the Shut data set, and the running time of the other three algorithms of NGLE, LEAR, and PRAR is (886.08, 3974.02, 133.25) seconds. When reducing the Mush dataset, IKAR takes 4.45 seconds, and the other three algorithms take (411.09, 4544.04, 827.16) seconds. Of course, the IKAR algorithm also has shortcomings, when the number of sample categories is more, the algorithm is more time-consuming. For example, if there are 10 different categories of samples in the Hand dataset, the calculation time of the IKAR algorithm takes 92.93 seconds. At this time, the time-consuming of the NGLE and PARA algorithms are (81.61, 91.28) seconds, respectively. It takes less time than that of the IKAR algorithm.

Table 3. The attribute reduction size with the four methods on the twelve UCI data sets.

Data sets	IKAR	NGLE	LEAR	PRAR
Prom	5	4	5	5
Hear	10	10	9	10
Hepa	9	10	7	10
Hand	12	11	10	11
Ches	29	30	29	30
Spli	9	10	9	9
Lett	9	9	8	9
Vote	10	9	8	10
Mush	4	5	5	5
Qsar	31	30	29	31
Shut	4	5	4	5
Sati	10	10	9	11
Ave	11.833	11.917	11.00	12.083
Best	4	1	11	1

Table 4. The run time with the four methods on the twelve UCI datasets.

Data sets	IKAR	NGLE	LEAR	PRAR
Prom	0.98	2.042	3.231	1.553
Hear	0.03	0.16	2.09	0.33
Hepa	0.06	0.11	2.13	0.25
Hand	92.93	81.61	3222.89	91.28
Ches	6.35	140.98	3075.25	590.79
Spli	35.43	205.86	6557.61	309.79
Lett	3.98	6.93	117.87	8.47
Vote	0.06	0.38	6.09	1.01
Mush	4.45	411.09	4544.04	827.16
Qsar	3.84	4.97	98.76	6.61
Shut	8.98	886.08	3974.70	1133.25
Sati	25.30	125.61	2322.02	154.04
Ave	15.20	155.49	1993.89	260.38
Best	11	1	0	0

From the above analysis, we can get that the algorithm IKAR can effectively reduce the data set, and it is obviously better than similar algorithms in terms of reduction speed.

5.2. Changes of the classification accuracy when missing value

It is not enough to evaluate the overall performance of a reduction algorithm only from the size and running time of the reduction set. Here we further analyze the performance of the above four algorithms from the perspective of the classification accuracy of the reduction set. In order to find out the influence of missing values on IKAR algorithm, we selected four data sets in Table 2, including Heas, Prom, Hepa, Stai. During the experiment, we divided the missing values into 5 different levels, 2%, 4%, 6%, 8%, and 10%. First, 2% of the data objects in the original data set were randomly selected, random deletion is conducted in these data objects of attribute values for different attributes to generate an incomplete data set. In order to reduce the bias of attribute selection that may be caused by random data deletion, we use the 2% data just obtained as a basis, and then select other attributes with the same probability, and then delete 2% data, generate an incomplete data set with a missing value of 4%, and so on, generate a data set with a missing value of 6%, 8%, and 10% respectively. After the incomplete data set is ready, we use IKAR, NGL, LEAR, PRAR four algorithms for reduction, and use SMO, J48 classifiers to analyze the classification accuracy of the reduced set. The specific results are shown in Figures 3 and 4. The horizontal axis in Figures 3 and 4 represents the proportion of deletion data objects in the data set, and the vertical axis represents the classification accuracy.

Figures 3 and 4 show the change trend diagrams obtained by using the SMO and J48 classifiers to analyze the accuracy of the reduced data set. From the results of Figures 3 and 4, it can be seen that when increasing of missing data, the classification accuracy of the above four algorithms has a downward trend. For example, under the SMO classifier, when the proportion of missing values in the data set Heas changes from 2% to 4%, the accuracy changes of IKAR and other NGL, LEAR and PRAR algorithms are (84.01→83.69), (80.32→79.69), (77.01→76.86), (79.24→78.92). Under the J48 classifier, when the proportion of missing values in the data set Heas changes from 2% to 4%, the IKAR and other NGL, LEAR, PRAR, the accuracy changes are (81.43→80.94), (77.85→76.28), (74.79→73.33), (76.24→74.86) respectively. The main reason for this phenomenon is that as the proportion of missing data increasing, more and more data objects cannot be distinguished, resulting in a decrease in classification accuracy. From the trend diagrams in Figures 3 and 4, it can be seen that the IKAR algorithm changes smoothly. The other three algorithms have a greater impact on the classification accuracy of the reduced set with the proportion of missing data increasing. For example, under the J48 classifier, when the missing proportion of the Hepa dataset changes from 2% to 10%, the accuracy of IKAR changes to 2.09, while the accuracy changes of the NML, LEAR and PRAR

algorithms are 3.2, 3.19 and 4.49 respectively. Under the SMO classifier, the classification accuracy of the Heas dataset, the IKAR algorithm changes value is 1.12, and the other three algorithms are 3.37, 2.28 and 2.93 respectively.

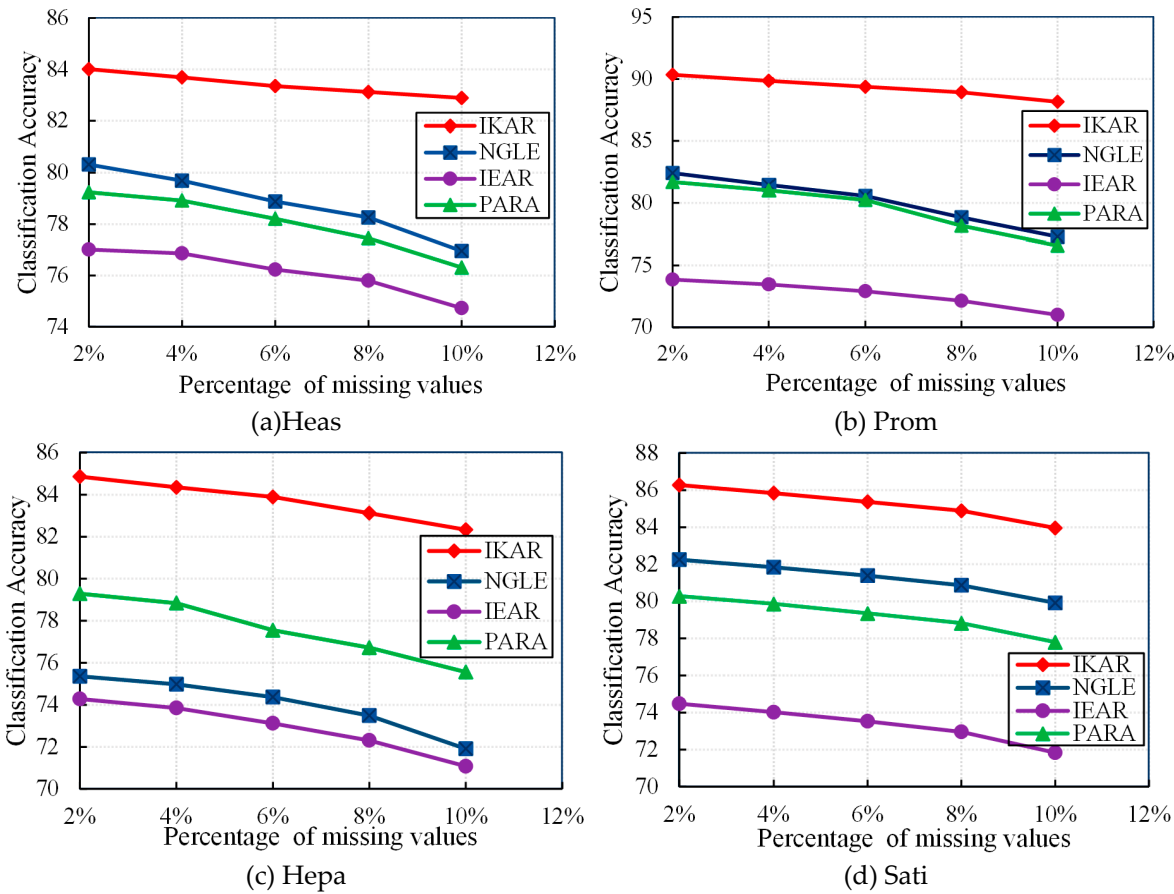
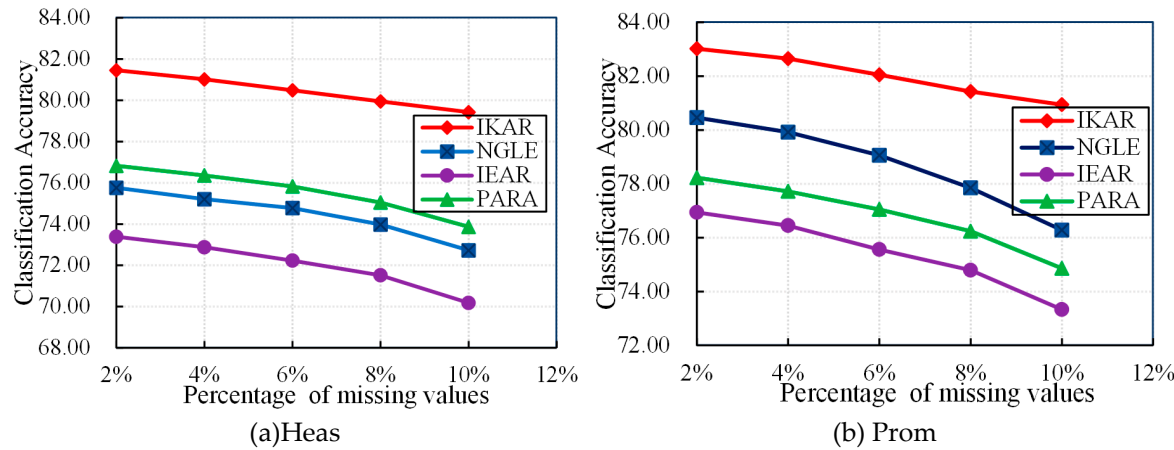


Figure 3. Variation of classification accuracy of different percentage of missing values with classifiers SMO.



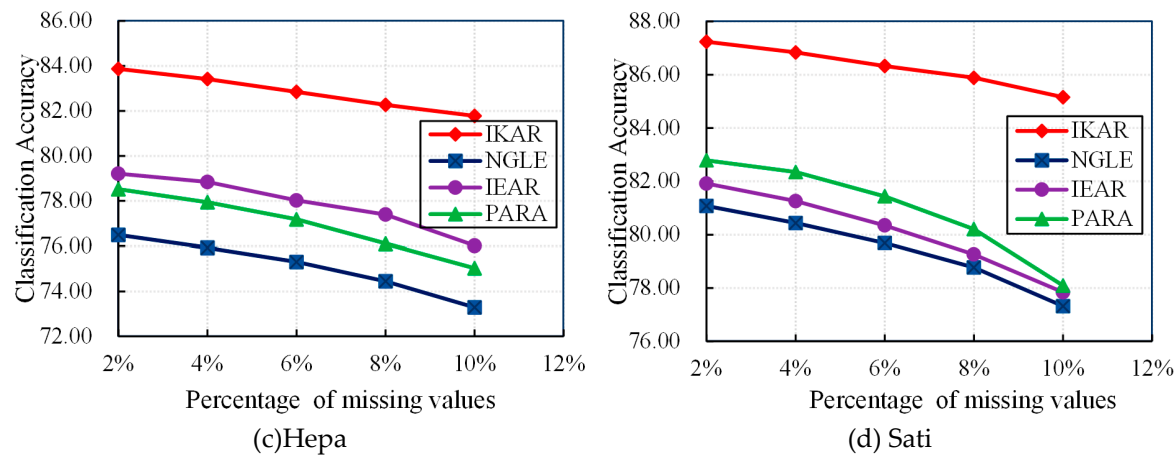


Figure 4. Variation of classification accuracy of different percentage of missing values with classifiers J48.

5.3. Classification accuracy analysis

The previous experiments have compared the changes of the classification accuracy when missing value. With the change of the proportion of missing value, the accuracy of the IKAR not only can change smoothly, but also get better classification effect on multiple data sets. We use the SMO and J48 classifier to analyze the accuracy of the previous reduction set, the details content is shown in Table 5 and Table 6 respectively. Under the classifier SMO, the IKAR algorithm obtained an average accuracy of 90.14 and the average accuracy obtained by the other three algorithms NGLE, IEAR and PRAR was only (83.05,82.03,82.98) respectively. Among the 12 data sets of Table 2, 11 times of highest accuracy are achieved with the IKAR algorithm. For example, the classification accuracy of IKAR is (94.76,99.68,93.86) in the data sets of Hand, Shut and Spli. The classification accuracy of IEAR is (81.59,85.93,81.53) and the average accuracy is 10 percentage points lower than IKAR. Similarly, on the J48 classifier, the average classification accuracy of the IKAR is 90.16, and the other three algorithms were obtained (84.79,83.53,84.98) respectively. Through the analysis in section 5.2, we can conclude that when the proportion of missing value changes, the trend of accuracy of IKAR algorithm is relatively stable, and it has a higher classification accuracy than other similar reduction models, which can effectively improve the classification quality of the reduction set.

Table 5. Comparison of the classification accuracies on reduced data sets with SMO(%).

Data sets	IKAR	NMLE	IEAR	PRAR
Prom	90.34(1)	83.33(2)	81.37(4)	82.15(3)
Heas	84.25(1)	80.02(3)	79.88(4)	81.39(2)
Hepa	85.81(1)	76.34(4)	84.62(4)	80.42(3)
Hand	94.76(1)	88.73(2)	81.59(4)	83.17(3)
Krkp	87.34(2)	81.29(4)	87.86(1)	82.83(3)
Spli	93.86(1)	84.86(2)	81.53(4)	82.49(3)
Lett	80.78(1)	73.49(3)	70.29(4)	74.26(2)
Vote	95.64(1)	90.21(2)	86.82(4)	89.33(3)
Mush	100.00(1)	98.34(3)	97.49(4)	98.96(2)
Qsar	82.95(1)	73.58(3)	70.61(4)	72.45(2)
Shut	99.68(1)	87.92(3)	85.93(4)	88.89(2)
Sati	86.27(1)	78.47(3)	76.39(4)	79.39(3)
Ave	90.14(1)	83.05(2)	82.03(4)	82.98(3)
AveRank	1.08	2.75	3.58	2.58

Table 6. Comparison of the classification accuracies on reduced data sets with J48(%).

Data sets	IKAR	NMLE	IEAR	PRAR
Prom	83.02(1)	80.46(2)	76.94(4)	78.23(3)
Heas	81.46(1)	75.76(3)	73.39(4)	76.83(2)
Hepa	83.87(1)	76.49(4)	79.21(2)	78.52(3)
Hand	94.69(1)	90.43(2)	88.36(4)	89.55(3)
Krkp	91.50(1)	85.31(3)	83.96(4)	86.45(2)
Spli	93.54(1)	88.51(2)	85.57(4)	87.37(3)
Lett	89.23(1)	83.86(3)	81.39(4)	84.07(2)
Vote	95.63(1)	87.91(2)	85.86(4)	86.99(3)
Mush	99.89(1)	98.75(3)	97.83(4)	98.92(2)
Qsar	83.96(1)	80.39(2)	78.27(4)	79.41(3)
Shut	97.92(1)	88.53(4)	89.61(3)	90.58(2)
Sati	87.24(1)	81.08(4)	81.92(3)	82.79(2)
Ave	90.16(1)	84.79(3)	83.53(4)	84.98(2)
AveRank	1.00	2.83	3.67	2.50

5.4. Lgorithm stability analysis

To indicate the statistical significance of classification results, non-parametric test methods Friedman test and Bonferroni-Dunn test [68] are used to analyze classification accuracy of each classifier with different methods in 5.2 section, where Friedman test is a statistical test that uses the ranking of each method on each dataset. The Friedman statistic is described as following:

$$\chi_F^2 = \frac{12N}{t(t+1)} \sum_{i=1}^s R_i^2 - 3t(t+1) \quad (26)$$

$$F_F = \frac{(N-1)\chi_F^2}{N(t-1) - \chi_F^2} \quad (27)$$

Where N is the number of data sets, t is the categories of algorithms. R_i represents the average rank of the classification effect ranking of the i-th algorithm on all data sets, and the statistics F_F obey the Fisher distribution with t-1 and (t-1)(N-1) degrees of freedom. If the value of F_F is bigger than $F_{\alpha}(t-1, N-1)$, then the original hypothesis does not hold.

The F_F test value can judge where these algorithms are different, but it cannot indicate the superiority of the algorithm. In order to explore which algorithm is better, we use the Bonferroni-Dunn test to calculate the critical value range of the average sequence value difference, which is defined as following,

$$CD_{\alpha} = q_{\alpha} \sqrt{\frac{t(t+1)}{6N}} \quad (28)$$

If the difference between the average ranking values of the two algorithms exceeds the critical region CD_{α} , the hypoth-esis that 'the performance of the two algorithms is the same' will be rejected with the corresponding confidence. Otherwise, the two algorithms perform differently, and the algorithm with the higher average rank is statistically better than the algorithm with the lower average rank. Generally, we set $\alpha=0.05$.

In order to compare the stability of the IKAR algorithm in this paper, we choose three other similar algorithms NGLE, LEAR and PRAR to reduce the data sets in Table 2. Then the previous reduction results are analyzed by the classifiers SMO and J48 in the Weka tool. The classification accuracy is detected by Friedman test and Bonferroni-Dunn test. When t=4 and N=12, then $\chi_F^2=23.4$ and $F_F=20.429$. Under the classifier SMO, the classification accuracy of the four algorithm reduction sets is sorted, and the number 1 represents the most ideal. The detail of sorting results is shown in

Table 5. The average ranking values of IKAR, NMLE, IEAR and PRAR algorithms are 1.08, 2.75, 3.58 and 2.58 in turn.

Since the critical value of $F_{0.05}(3,33)$ is 2.892 and $F_F > 2.892$, we can reject the original hypothesis at $\alpha = 0.05$ under the Friedman test. So, there are statistical differences in classification accuracy among the above four algorithms. Next, $q_{0.05} = 2.569$ and $CD_{0.05} = 1.354$, the results of Bonferroni-Dunn test for these four algorithms at $\alpha = 0.05$ is shown in Figure 5. The accuracy of the algorithm on the left side of the coordinate axis is relatively high in Figure 5. From Figure 5, we know that the average accuracy ranking of IKAR is $s_1 = 1.08$. Among the other three algorithms, the better-ranking algorithm PRAR has an average ranking value of $s_2 = 2.58$. Since $|s_1 - s_2| = 1.5$ and $|s_1 - s_2| > CD_{0.05}$, we can find that the IKAR algorithm is significantly superior to the PRAR. In the same way, the IKAR is better on accuracy than NGLE and LEAR. However, there is no obvious difference in the ranking of the NMLE, LEAR and PRAR algorithms.

For the same reason, under the classifier J48, the average ranking values of the classification accuracy of the above four algorithms are 1.00, 2.83, 3.67 and 2.50 as shown in Table 6. Since $\chi_F^2 = 26.8$ and $F_F = 32.043$, then $F_F > F_{0.05}$. We can get that the classification accuracy of these four algorithms is significantly different in the statistical sense under the J48 classifier. From Figure 6, we can get that the IKAR algorithm is significantly different from the NGLE classification, while the ranking of the NGLE, LEAR and PRAR algorithms have no obvious differences among each other.

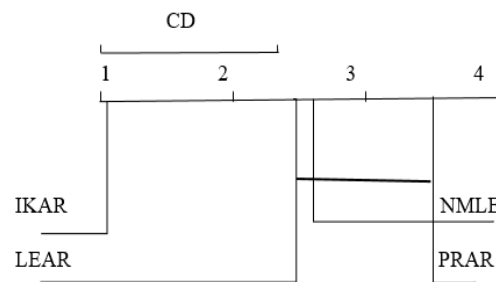


Figure 5. Bonferroni-Dunn test with SMO.

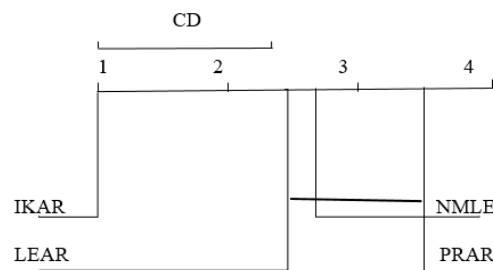


Figure 6. Bonferroni-Dunn test with J48.

Therefore, all the test results demonstrate that there is no consistent evidence to denote statistical differences between any two of four approaches under the SMO and J48 classifier. In general, the IKAR model is better than the other models in stability.

6. Conclusion

In the face of incomplete systems, most of the traditional attribute reduction models cannot obtain effective reduction rules, and affect the classification accuracy of the reduced set. Therefore, we propose a new attribute reduction method IKAR based on the clustering background for incomplete system. IKAR uses the tolerance principle to calculate the information of knowledge granularity, and to measure the similarity of data objects. When selecting attributes, the similarity of intra-cluster objects should be as large as possible, and the similarity of inter-cluster objects should

be as small as possible. The innovations of this paper are manifested in the following four aspects: (1) Use the tolerance principle to quickly calculate knowledge granularity; (2) Use knowledge granularity to calculate the similarity of inter-cluster and intra-cluster objects; (3) Use the idea of clustering to calculate the importance of attributes; (4) In addition to conventional time, space and precision analysis, it also analyzes the stability of data sets with different percent missing value. All experiments show that the IKAR algorithm is not only superior in reducing time compared with the other three algorithms, it also has an excellent performance in terms of accuracy and stability. Of course, the IKAR algorithm also has some shortcomings. For example, it is unsuitable for data sets with multiple decision values, complex data type and the dynamical changing of the data sets are not considered.

For further researches, we focus on attribute reduction of incomplete systems with mixed data types, especially how to deal with missing data? In addition, we will integrate the incremental learning method into the knowledge granularity reduction model in the background of clustering. Even to address big data sets more efficiently, applying GPU and MapReduce technologies to design parallel attribute reduction model or acceleration strategy is a very popular research topic.

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