

A NOVEL APPROACH TO DATA MINING USING SIMPLIFIED SWARM OPTIMIZATION

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Noorhaniza Wahid

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ABSTRACT

Data mining has become an increasingly important approach to deal with the rapid growth of data collected and stored in databases. In data mining, data classification and feature selection are considered the two main factors that drive people when making decisions. However, existing traditional data classification and feature selection techniques used in data management are no longer enough for such massive data. This deficiency has prompted the need for a new intelligent data mining technique based on stochastic population-based optimization that could discover useful information from data.

In this thesis, a novel Simplified Swarm Optimization (SSO) algorithm is proposed as a rule-based classifier and for feature selection. SSO is a simplified Particle Swarm Optimization (PSO) that has a self-organising ability to emerge in highly distributed control problem space, and is flexible, robust and cost effective to solve complex computing environments. The proposed SSO classifier has been implemented to classify audio data. To the author's knowledge, this is the first time that SSO and PSO have been applied for audio classification.

Furthermore, two local search strategies, named Exchange Local Search (ELS) and Weighted Local Search (WLS), have been proposed to improve SSO performance. SSO-ELS has been implemented to classify the 13 benchmark datasets obtained from the UCI repository database. Meanwhile, SSO-WLS has been implemented in Anomaly-based Network Intrusion Detection System (A-NIDS). In A-NIDS, a novel hybrid SSO-based Rough Set (SSORS) for feature selection has also been proposed. The empirical analysis showed promising results with high classification accuracy rate achieved by all proposed techniques over audio data, UCI data and KDDCup 99 datasets. Therefore, the proposed SSO rule-based classifier with local search strategies has offered a new paradigm shift in solving complex problems in data mining which may not be able to be solved by other benchmark classifiers.

PUBLICATIONS

Refereed Journal Articles:

1. Wei-Chang Yeh, Noorhaniza Wahid, Yuk Ying Chung. “Effective Audio Classification Algorithm using Swarm-based Optimization Technique Approach”, submitted to Expert System with Applications (SCI journal with Impact Factor 2.908).
2. Wei-Chang Yeh, Noorhaniza Wahid, Yuk Ying Chung. “A Hybrid Network Intrusion Detection System using Simplified Swarm Optimization (SSO)”, submitted to Applied Soft Computing (SCI journal with Impact Factor 2.415).
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Refereed Conference Paper:

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ABBREVIATIONS

Acronym	Expended term
ABC	Artificial Bee Colony
ACO	Ant Colony Optimization
ADEL	Ankyloglossia with Deviation of the Epiglottis and Larynx
AI	Artificial Intelligence
AIS	Artificial Immune Systems
A-NIDS	Anomaly-based Network Intrusion Detection Systems
ANN	Artificial Neural Network
ARS	Average Rule Size
BPSO	Binary Particle Swarm Optimization
BW	Baum-Welch
CHD	Coronary Heart Disease
CLPSO	Comprehensive Learning PSO
CPSO	Constricted PSO
CPSO	Cooperative Particle Swarm Optimizer
DE	Differential Evolution
DFT	Discrete Fourier Transform
DoS	Denial of Service
DPSO	Discrete PSO
EAs	Evolutionary Algorithms
ELS	Exchange Local Search
EP	Evolutionary Programming
ES	Evolution Strategy
FCM	Fuzzy C-Means
FCPS-classifier	Fuzzy Controlled Particle Swarm-classifier
FDTs	Fuzzy Decision Trees
FIPS	Fully Informed Particle Swarm
FN	False Negatives
FP	False Positives
FPPSO	Feature Partitioning PSO
GA	Genetic Algorithm
GEPSVM	Generalized Eigenvalue Proximal SVM
GMM	Gaussian Mixture Model
GNB	General Naive Bayes
GP	Genetic Programming
GRBF	Gaussian RBF
HEHRS	Hierarchical Rule Sets
H-IDS	Host-based IDS

Acronym	Expended term
HMM	Hidden Markov Model
IA	Immune Algorithm
IDS	Intelligent Dynamic Swarm
IDSRS	Intelligent Dynamic Swarm with Rough Set
IkNN	Informative k Nearest Neighbor
IPS-classifier	Intelligent Particle Swarm-classifier
IS	Immune System
KDD	Knowledge Data Discovery
kNN	kNearest Neighbour
KPCM	Kernel-based Possibilistic C-Means
LCM	Linear time Closed itemset Miner
LDPSO	Linearly Decreasing PSO
LDWPSO	Linear Decreasing Weight PSO
LGP	Linear Genetic Programming
LMNN	Large Margin Nearest Neighbor
LPC	Linear Predictive Coding
MARS	Multivariate Adaptive Regression Splines
MFCC	Mel Frequency Cepstral Coefficient
MI	Mutual Information
MLP	MultiLayer Perceptron
MOPSO	Multi-Objective PSO
MSSE	Mean Sum Squared Error
NB	Naive Bayes
NFL	Nearest Feature Line
NIDS	Network Intrusion Detection Systems
NIP	Numerical Interval Pruning
NN	Neural Network
NN	Nearest Neighbor
OBSI	Octave Band Signal Intensities
P2PkNN	Peer-to-Peer kNN
PCA	Principal Component Analysis
PM	Partition Matrix
PSO	Particle Swarm Optimization
PSO-NCC	PSO-Nearest Centroid Classifier
PUNN	Product Unit Neural Network
QDE	Quantum-Inspired Differential Evolution
QPPs	Quadratic Programming Problems
R2L	Remote to Local
RBF	Radial Basis Function
RHMM	Reverse HMM
RJMCMC	Reversible Jump Markov Chain Monte Carlo

Acronym	Expended term
RSC	Rough Set Classification
RSC-PGA	Rough Set Classification Parallel Genetic Algorithm
RST	Rough Sets Theory
RWS	Roulette Wheel Selection
SA	Simulated Annealing
SI	Swarm Intelligence
SMO	Sequential Minimal Optimization
SNNB	Selective Neighborhood Naive Bayes
SOM	Self Organizing Maps
SR	Silence Ratio
SSO	Simplified Swarm Optimization
SSO-ELS	SSO with Exchange Local Search
SSO-WLS	SSO with Weighted Local Search
SVDF	Support Vector Decision Function
SVM	Support Vector Machine
TN	True Negatives
TP	True Positives
TSP	Traveling Salesman Problem
TSVM	Twin SVM
U2R	User to Root
WEKA	Waikato Environment for Knowledge Analysis
WFPPSO	Weighted Feature Partitioning PSO
WLS	Weighted Local Search
WSVM	Weighted Support Vector Machine

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CHAPTER 1. Background Information and Problem Statements

In the emerging age of digital information we are overwhelmed with data, while our capability to analyse and interpret such huge datasets lags behind. Furthermore, it has been estimated that every 20 months the amounts of data stored in the world's databases are doubled, which caused difficulties when trying to justify this figure in a quantitative sense [1]. Often, traditional data analysis, and interpretation of changing data, has become insufficient for data processing as the data volumes grow exponentially. In addition, due to the advancement of software capabilities and hardware tools that enable the automated data collection, as well as the decreasing trend in their cost, there has been a dramatic increase in the data being collected and stored in databases. Although in recent years information collection and storage has become easier and more inexpensive, great effort is required to extract relevant knowledge information from such large-scale databases. Therefore, a new generation of computational techniques and tools is required to support the extraction of useful knowledge from the rapidly growing volume of data. Hence, data mining becomes the reliable solution for elucidating the patterns that underlie it.

Data mining is the application of specific algorithms that has been widely used for extracting patterns or models from data. Two main aspects in data mining are data classification and feature selection. Data classification classifies a data item into one of several predefined categorical classes. Feature selection can be defined as a process of choosing a small subset of features from the original set of features which is necessary and sufficient to describe the target concept. Other than the well-known

classical data mining techniques, heuristic approaches based on swarm intelligence algorithms have gained more attention and have been adopted in data classification problems in order to find a good solution. This thesis proposes and presents some new data mining approaches based on a population-based optimization algorithm for various data classification problems. The second topic being discussed in this thesis is about feature selection, which presents a new hybrid rough set reduction approach to feature selection. At this stage, this thesis is concerned with finding new approaches in both topics that contribute to the best classification accuracy; computation time is not taken into consideration. In this chapter, section 1.1 presents a brief discussion on data mining from the perspective of Knowledge Data Discovery (KDD). In section 1.2, the motivation is described and the problem statement in this thesis is identified. Next, the objective and contribution of the thesis is stated in section 1.3, followed by the organisation of the thesis in section 1.4.

1.1. *Data Mining*

Data mining is the process of analysing data from different perspectives and summarising it into useful information. It blends traditional data analysis methods with sophisticated algorithms for processing large volumes of data. It has been widely used and unifies research in fields such as statistics, databases, machine learning and Artificial Intelligence (AI). Regarding that, data mining has been seen as an explosion of interest from both academia and industry to improve the process of visualising and understanding the pattern of the data. Data mining is the core part of the Knowledge Discovery in Database (KDD) process, which is essential to solve a problem in a specific domain [2]. Generally, KDD is the overall process of identifying valid, novel,

potentially useful and ultimately understandable patterns in data and converting it into useful information [3]. An overview of the steps constituting the KDD process is depicted in Figure 1.1.

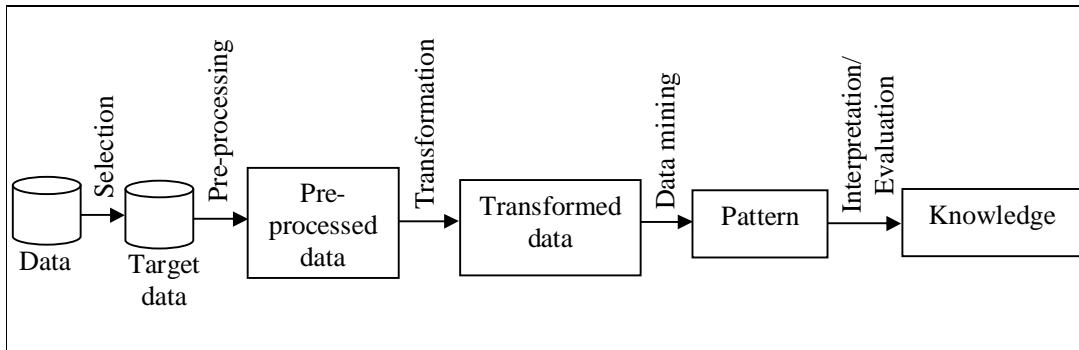


Figure 1.1. The processes of KDD [3]

Data mining in KDD applies a specific algorithm to extract meaningful knowledge so that the discovered knowledge can be applied in the related area to increase working efficiency and also to improve the quality of decision-making. Data mining involves several steps such as data integration from various databases, data pre-processing, and induction of a model using a learning algorithm. Based on the requirements of the problem domain, various techniques that expose diverse kinds of patterns from a given dataset have been implemented in data mining. The most common techniques learned in data mining include data classification, data clustering, association rule discovery, and outlier detection.

1.2. Motivation and Problem Statements

As mentioned earlier, data mining has been widely used to solve various kinds of data classification problems. However, data classification has turned out to be one of the most pervasive problems that encompasses many diverse applications in the data mining field. These problems have attracted more active research in order to find

efficient approaches to address them, and the outcome of the research is still unsatisfactory.

The ultimate goal of classification is to discriminate new data into the most likely of the specific categorical variable (the class) based on the induction model generated by the classifier. However, the classification problem has become very complicated and computationally infeasible when the number of possible different combinations of variables is so high. Hence, Swarm Intelligence (SI) algorithms are generally more suitable to solve these difficult problems because they are based on stochastic population-based approaches. In addition, they are also capable of avoiding becoming stuck in a local optimal and can find a global optimal solution [4].

Many data mining approaches have been proposed using stochastic population-based algorithms such as Particle Swarm Optimization (PSO), Immune Algorithm (IA), Artificial Bee Colony (ABC) and Ant Colony Optimization (ACO). Nevertheless, there are problems in real-world that are NP-hard and combinatorial. Thus, evolutionary algorithm like PSO is generally more suitable to solve these difficult problems because of its stochastic nature. PSO is a well-known, biologically inspired computational search and optimization algorithm which is based on the social behaviours of bird flocks or schools of fish [5]. Because of its easy implementation, PSO has been successfully applied in many fields, particularly in optimization applications [6] and data mining [6-8]. This is due to its simplicity and efficiency when navigating a search space for optimal solutions. In terms of data mining, PSO has emerged as a promising technique to discover useful and interesting knowledge from databases [8]. Because of these advantages, the motivation of this thesis is to develop a new data mining technique based on the original PSO algorithm.

In many applications, people are dealing with massive data that contains multi-dimensional attributes such as network intrusion data, stock market data, medical data, weather forecast data and much more. Thus, data classification is faced with a problem when it has to generate rules with many attributes or features. Obviously, the time required to generate rules is proportional to the number of features. In addition, irrelevant and redundant features can reduce both the predictive accuracy and comprehensibility of the induced rule and degrade the classifier speed (due to its high dimensionality). Thus, selecting the most relevant features is necessary, and this strategy is implemented to simplify the rules and reduce its computational time while retaining the quality of classification, as it represents the original features set.

This thesis proposes and investigates the application of a new efficient population-based optimization algorithm for data mining based on the PSO algorithm. The new technique is referred as a Simplified Swarm Optimization (SSO) algorithm. Like PSO, SSO solves data mining problems by simulating the social interaction among agents or particles in their population, such as birds flocking or fish schooling. To deal with the problem of feature selection, a new hybrid swarm intelligence-based rough set theory for feature selection using SSO is proposed as a way to improve some performance criterion, such as accuracy of data classification. In this thesis, the author is concerned with introducing a new population-based optimization technique for data mining and feature selection that contributes to maximising the classification accuracy. Therefore, computation time is not taken into consideration.

Throughout this thesis, each algorithm was implemented using Java NetBeans IDE 6.1 on the following system: 1.8GHz Pentium (R) processor and 2GB RAM running in Windows XP Professional. Five traditional classifiers were involved in the

experiments for comparison with the proposed technique. Those benchmark classifiers were implemented from Waikato Environment for Knowledge Analysis (WEKA) [1]. The employed classifiers were set with their default parameters as set in WEKA.

1.3. Objectives and Contributions of the Thesis

This section outlines the main objectives and contributions to the area of data mining, particularly in feature selection and classification problems.

- i. To develop and implement an efficient data classification technique based on an SSO algorithm to be implemented in audio datasets.
- ii. To develop and implement a novel Exchange Local Search (ELS) strategy to improve the performance of the SSO rule-based classifier on various datasets.
- iii. To develop and implement a new hybrid SSO-based Rough Set for feature selection and Weighted Local Search (WLS) strategy with SSO classifier to improve the Anomaly-based Network Intrusion Detection System (A-NIDS).

1.4. Outline of the Thesis

The remainder of this thesis is structured as follows:

“CHAPTER 2 Literature Review of Feature Selection and Data Mining” comprehensively presents two main topics that cover the foundations of feature selection and data mining. These topics provide a review of recent work that has been conducted in feature selection (with more emphasis on Rough set theory) and in data mining (with more emphasis on traditional classification techniques).

“CHAPTER 3 Data Mining using the Particle Swarm Optimization Algorithm” provides some introduction to four population-based optimization algorithms (with more emphasis on the PSO). This chapter also reviews the implementation of PSO algorithms for data mining purposes in various applications and problem domains. Four approaches have been highlighted in PSO-based classification including: PSO for Rule-based Classification Model, Nearest Neighbor Classification, PSO as Optimizer within Other Learning Algorithms, and Clustering with PSO Algorithms. Also, some PSO variants for data classification are discussed in this chapter.

“CHAPTER 4 Data Mining using Simplified Swarm Optimization Algorithm” presents the Simplified Swarm Optimization (SSO) algorithm that is based on traditional PSO for data mining. This is followed by the principle of the SSO algorithm; the SSO rule mining scheme; the SSO rule evaluation; and SSO rule pruning. The proposed algorithm is then applied to audio data and it is compared with Support Vector Machine (SVM) to investigate its competitiveness.

“CHAPTER 5 The Proposed SSO with Exchange Local Search for Data Classification” presents a proposed Exchange Local Search (ELS) strategy to be incorporated with SSO for data classification. To show the applicability of the proposed approach, SSO with ELS (SSO-ELS) is then applied to 13 datasets obtained from public sources such as the UCI repository database. The performance is compared with and without ELS for SSO and PSO, and four other traditional classifiers including SVM, J48, PART and kNearest Neighbor.

“CHAPTER 6 A Hybrid SSO-based Rough Set Reduction Method for Network Intrusion Detection Systems” introduces a proposed hybrid SSO-based rough set

reduction for features dimensionality reduction. This approach is specifically applied to solve the classification problem in Anomaly-based Network Intrusion Detection Systems (A-NIDS) due to its large amount of attributes. This is followed by the principle of the proposed SSO with Weighted Local Search (WLS) strategy for mining anomaly intrusion patterns. The experimental results, when compared with original SSO and PSO, and also with SVM and Naive Bayes, show the effectiveness of hybridizing SSO-WLS approaches for A-NIDS detection.

“CHAPTER 7 Conclusions and Future Work” contains a summary, conclusion, limitations and future direction of the research conducted in this thesis.

CHAPTER 2. Literature Review of Feature Selection and Data Mining

In recent years, the field of automated data mining has emerged as an important area of applied research when dealing with the voluminous data collected in various industries. This is due to the low cost and availability of larger storage devices. Thus, two major data mining tasks that must be solved are feature selection and classification. In this chapter, the investigation on several techniques of feature selection and data classification is continued, and comprehensive reviews on both topics are presented in section 2.1 and section 2.2. In this thesis, the new data mining algorithms based on the population-based optimization algorithm are proposed in Chapter 3.

2.1. Feature Selection Overview

Feature selection plays an important role in data pre-processing technique for data mining [2]. It is a process of finding a subset of features from the original set of features, and forming patterns in a given dataset to obtain the optimal one according to the given goal of processing and criterion. It reduces the number of features, removes irrelevant, redundant, or noisy data and brings immediate effects for applications: speeding up a data mining algorithm, improving mining performance such as classification accuracy, and improving results comprehensively.

In the context of classification, feature selection can be structured into three fractions: *filter* method, *wrapper* method and *embedded* method [9]. Filter methods rely on the intrinsic properties of the training data to select some features without involving any

learning algorithm. Each feature is ranked according to some univariate metric, and only the highest ranking features are used while the remaining low ranking features are eliminated. Afterwards, this subset of features is presented as input to the classification algorithm. Therefore, feature selection is allowed to be performed only once, and then different classifiers can be evaluated. A number of multivariate filter techniques were introduced to overcome univariate problems in filter methods.

Wrapper methods embed the model hypothesis search within the feature subset space. These methods begin by looking for the dependency from a suboptimal subset. Then this value is fed into the fitness function of the selected learning algorithm and evaluated in order to find the suitable features. These methods suffer from a high risk of overfitting and require huge computational cost.

Meanwhile, in the third category of feature selection, namely, embedded methods, the search for an optimal subset of features is built into the classifier construction, which can be seen as a search in the combined space of feature subsets and hypotheses. Thus, their function is seen as more specific to a given learning algorithm. Embedded methods are less computationally intensive than wrapper methods due to the internal interaction with the classification model during the feature selection process.

2.1.1. Feature Selection Problems

The feature selection problem is more or less a special case of a much broader problem of subset selection. Suppose a large set of M items $\{x_k, y_k\}$ where $k = 1, 2, \dots, M$ consisting of n input variables $x_{k,i}$ where $i = 1, 2, \dots, n$ and one output variable y_k is given from which we need to find a small m subset being optimal in a certain sense. Fitness function (F_i) is computed from the values $x_{k,i}$ and y_k , $k = 1, 2, \dots, m$ to rank the

variables with a high F_i value. Then, the features with a high F_i value will be selected as the most relevant features. In statistical subset selection, being optimal usually means being most suitable for classification or data approximation. Although it may not be apparent, the subset selection problem may become prohibitive because of its computational complexity [10]. It is not possible to rank the items in our set simply according to their individual properties and select only the best. The item properties may depend strongly on each other and a subset of individually “bad” features may prove to be rather “good” because of positive interaction effects. Because of this uncertainty, the only apparent way of searching for optimal subsets is to evaluate all the 2^n possible item combinations.

Generally, feature selection is composed of two parts. First, there is a searching strategy that selects a feature subset. By searching all the feature subset, we shall ensure the number of $\binom{N}{M}$ where N is the original features, M is the feature subset, and $M < N$. If we optimized an M , we should be able to estimate the combination number of 2^N . Therefore, we must select a good searching strategy to evaluate quite a number of combinations because it is definitely inefficient to evaluate the whole feature sets. In addition, testing all subsets is a combinatorial NP-hard [11] problem that requires an exponential amount of computational time. Second, there is a fitness function that evaluates the feature subsets, returns their scores to the corresponding subsets and finally selects the best subset.

Feature selection has been used as a measure to decide the importance and necessity of features. Hence, a broad range of sub-optimal feature selection methods have been developed over the last decade to overcome its limitation. Previous research on

feature selection which employed a heuristic approach, such as hill-climbing methods, has proven more efficient when dealing with little noise and a small number of features [12, 13]. On the other hand, due to the reason that heuristic methods fail to find an optimal reduct [14], many researchers have shifted to metaheuristic approaches such as Genetic Algorithm (GA), Simulated Annealing (SA), Tabu Search, Ant Colony Optimization (ACO) and Particle Swarm Optimization (PSO) [15].

Previous research has introduced various techniques of feature selection with the use of SI algorithms, either in single or hybrid approaches. Bello et al. [16] proposed a two-stage heuristic search performed by PSO and ACO in order to find the optimal feature subsets. In [17], the authors used chaotic BPSO with logistic map to determine the inertia weight to solve the feature selection problem. As a result, their approach has successfully reduced the computation time as well as improved the quality of reducts and classification accuracies. Moreover, several hybrid methods that used SI algorithms and rough sets for feature selection have been reported. Methods which combine ACO and rough set to find reducts with promising results were proposed by Ke et al. and Chen et al. in [15] and [18], respectively. Later, inspired from the work of Wang et al. [14], a new algorithm to find minimal rough set reducts by using modified PSO, namely, Intelligent Dynamic Swarm (IDS), was introduced by Bae et al. [19]. The experimental results from that work showed that IDS is efficient for rough set-based feature selection. In the next section, this study only limit to the concept of rough set theory for the purpose of the feature selection approach.

2.1.2. Rough Sets Theory

Rough sets theory (RST), which was proposed by Pawlak [20], has been successfully applied as a feature selection method to discover data dependencies and reduce the number of attributes contained in a dataset. It involves mathematical theory that is able to deal with imprecise, uncertain and vague information [21]. At the first stage of the feature selection procedure, rough set is used to reduce and clean data with minimal model assumption. Then the result is used as a basis for further analysis performed with our proposed method. Rough set has been proven as an efficient tool for feature selection as it does not require any preliminary or additional knowledge except for the supplied data.

Table 2.1. A decision table

Object	a_1	a_2	a_3	a_4
x_1	2	3	1	4
x_2	1	4	3	5
x_3	2	3	1	4
x_4	3	2	1	1
x_5	1	2	1	2
x_6	1	2	3	2
x_7	1	4	3	5
x_8	2	1	1	2
x_9	1	2	3	4
x_{10}	1	4	3	5

In rough set theory, the data is organised into a table called a decision table. Table 2.1 shows the structure of the table which consists of rows and columns. The rows of the decision table correspond to objects, and columns correspond to attributes. In the dataset, a class label indicates the class to which each row belongs. The class is called

the *decision attribute*, while the rest of the attributes are the *condition attributes*. Some basic concepts of rough set are explained below.

The information system of a decision table is defined as a tuple $T = (U, A)$, where U is a non-empty set of finite objects (the universe) and A is a non-empty finite set of attributes. As in Table 2.1, the information table assigns an object x to the universe U with $U = \{x_1, x_2, \dots, x_{10}\}$. Whereas each attribute or feature $a \in A$ is associated with a set $\forall a$ of its value denoted as $A = \{a_1, a_2, \dots, a_4\}$. It is called the domain of a . The attribute set A can be partitioned into two subsets, C and D , which are called condition and decision attributes, respectively.

For a subset of attributes $R \subseteq A$, the R -indiscernibility relation is denoted by $IND(R)$, is an equivalence relation defined as:

$$IND(R) = \{(x, y) \in U \times U : \forall a \in R, a(x) = a(y)\} \quad \text{Eq. 1}$$

where $a(x)$ denotes the value of feature a of object x . If $(x, y) \in IND(R)$, x and y are said to be indiscernible with respect to R . The family of all equivalence classes of $IND(R)$ is denoted by $U/IND(R)$. Each element in $U/IND(R)$ is a set of indiscernible objects with respect to R . Equivalence classes $U/IND(C)$ and $U/IND(D)$ are called condition and decision classes. According to the objects from the example in Table 2.1, we consider the full set of attributes $A = \{a_1, a_2, a_3, a_4\}$ and the sets of $U/IND(R)$ can be listed as $U/IND(R) = \{\{x_1, x_3\}, \{x_2, x_7, x_{10}\}, \{x_4\}, \{x_5\}, \{x_6\}, \{x_8\}, \{x_9\}\}$.

For any concept $X \subseteq U$ and attribute subset $R \subseteq A$, X could be approximated by the R -lower approximation and R -upper approximation using the knowledge of R . The lower approximation of X is the set of elements of U that are surely in X is defined in

Eq. 2. The upper approximation of X is the set of elements of U that are possibly in X is defined as in Eq. 3.

$$\underline{R}(X) = \bigcup \{E \in U / IND(R) : E \subseteq X\} \quad \text{Eq. 2}$$

$$\overline{R}(X) = \bigcup \{E \in U / IND(R) : E \cap X \neq \emptyset\} \quad \text{Eq. 3}$$

Consider the target set $X = \{x_1, x_2, x_3, x_4\}$, therefore the R -lower approximation and R -upper approximation can be represented as $\underline{R}X = \{x_1, x_3, x_4\}$ and $\overline{R}X = \{x_1, x_3, x_4, x_2, x_7, x_{10}\}$, respectively. Let R and S be equivalence relations over U , then the positive, negative and boundary regions can be defined as:

$$\begin{aligned} POS_R(S) &= \bigcup_{X \in U/S} \underline{R}X \\ NEG_R(S) &= U - \bigcup_{X \in U/S} \overline{R}X \\ BND_R(S) &= \bigcup_{X \in U/S} \overline{R}(X) - \bigcup_{X \in U/S} \underline{R}(X) \end{aligned}$$

The positive region, $POS_R(S)$ is a set of objects of U that can be classified with certainty to classes $U/IND(S)$ employing attributes of R . The negative region $NEG_R(S)$, is the set of all elements U that cannot be classified to classes of U/S . On the other hand, the boundary region $BND_R(S)$ is the set of all elements that can be possibly, but not certainly, classified in this way. If the boundary region is empty, that is, $\overline{R}(X) = \underline{R}(X)$, X is said to be R -definable or crisp. In the opposite case, if $BND_R(S) \neq \emptyset$, the S is referred to as *rough (indefinable)* set with respect to R .

Another important issue in data analysis is discovering dependencies between attributes. Let R and S be subsets of A . We will say that S depends on R in a degree k ($0 \leq k \leq 1$), denoted $R \Rightarrow_k S$ if

$$k = \gamma_R(S) = \frac{|POS_R(S)|}{|U|} \quad \text{Eq. 4}$$

If $k = 1$, S depends totally on R , and if $k < 1$, S depends partially on R , and if $k = 0$, then S does not depend on R . In other words, S depends totally (partially) on R , if all (some) elements of the universe U can be surely classified to blocks of the partition U/S employing R [14]. Dependency degree k can be used as heuristic in greedy algorithms to compute feature reduction.

The reduction of attributes is achieved by comparing equivalence relations generated by a set of attributes. Let $A = (U, C \cup D)$ be a decision table where C is the set of condition features, and D is the set of decision features. The degree of dependency between condition and decision features, $\gamma_R(D)$, is called the equality of approximation of classification induced by the set of decision features [22]. The aim of feature reduction is to remove redundant features so that the reduced set provides the same quality of classification as the original set. A *reduct* is defined as a subset of minimal cardinality R_{min} of the conditional attribute set C in such a way that $R \subseteq C$ is said to be a D -reduct of C if $\gamma_R(D) = \gamma_C(D)$. A given decision table may have many feature reducts, thus the set of all reducts is defined in Eq. 5.

$$Red = \{R \subseteq C \mid \gamma_R(D) = \gamma_C(D) \forall B \subset R, \gamma_B(D) \neq \gamma_C(D)\} \quad \text{Eq. 5}$$

For instance, suppose that $C = \{a_1, a_2, a_3, a_4\}$, $R_1 = \{a_2, a_3, a_4\}$ and $R_2 = \{a_1, a_2, a_4\}$.

As a reduct is the minimal set of attributes preserving the positive region such that

$$\gamma_{R_1}(D) = \gamma_C(D) \text{ and } \gamma_{R_2}(D) = \gamma_C(D), \text{ then } Red = \{\{a_2, a_3, a_4\}, \{a_1, a_2, a_4\}\}.$$

The intersection of all the sets in R_{min} is called the *core*. The core is the set of attributes that are contained by all reducts, defined as

$$CORE_D(C) = \bigcap RED_D(C) \text{ where } RED_D(C) \text{ is the } D\text{-reduct of } C. \text{ In other words,}$$

the core is the set of attributes that cannot be removed without changing the positive region. This means that all attributes present in the core are indispensable. Therefore, from the above example of *Red*, we conclude that $core(C) = \{a_2, a_4\}$.

The applications of the rough set model in feature selection problems have demonstrated its usefulness in a wide range of techniques from rough set [12], fuzzy rough set [10], and neural network [23]. Recently, the use of the metaheuristic approach has been widely implemented and has significantly increased the ability of finding very high quality solutions to select more dominating features in order to perform the classification tasks optimally in the data mining process. For example, hybrid rough with GA [24] and Genetic Programming (GP) [25] was proposed by Wa'el et al. for choosing the most relevant features in a network intrusion dataset. Chen et al. [18] and Ke et al. [15] proposed a feature selection mechanism based on rough-Ant Colony Optimization (ACO). Inspired from the work of Wang et al. [14, 26], which used PSO to find optimal features subset based rough set reducts, Bae et al. [19] proposed a new discrete PSO called Intelligent Dynamic Swarm (IDS) with rough set (IDSRS) for feature reduction. Instead of successfully reducing the massive

number of features from the original set, IDSRS could improve the computation speed when compared with the original PSO.

2.2. Data Mining Overview

In practice, the two primary goals of data mining tend to be *prediction* and *description*. Prediction involves using some variables or fields in the dataset to predict unknown or future values of other variables of interest. On the other hand, description emphasises finding patterns describing the data that can be interpreted by humans. The goal of predictive data mining is to produce a model expressed as an executable code which can be used to perform data mining tasks such as classification, prediction or other similar tasks. On the other hand, the goal of descriptive data mining is to gain an understanding of the analysed system by uncovering patterns and relationships in large datasets. Thus, the goals of prediction and description are achieved by using data mining techniques to solve data mining problems by realising several data mining tasks such as data classification, estimation, prediction, and association rules. The following section particularly focuses on data classification, which is our main consideration in this thesis.

2.2.1. Classification Task in Data Mining

Data classification is the most studied task in data mining. It is a process of discovering a predictive learning function that classifies a set of data items into one of several predetermined classes. All these data are typically represented by matrices whose rows are associated with the *objects* (or records), while the columns are associated with many measures taken on each object. Such measures are often referred to as *variables* (or features). The relationship in each record is tied up by a

specific attribute called a class *attribute*. The goal of the classification algorithm is to induce a model so that it can classify any new instances into classes according to class values. Data classification can be divided into two categories: *unsupervised classification* and *supervised classification*.

- i. ***Unsupervised classification***: A method of unsupervised classification uses only the matrix of dissimilarity. No information of the class of an object is provided to the method. The objective is then to build a set of groups where each group contains a set of sufficiently similar objects. The number of classes and their contents are not known at the beginning. In data mining, unsupervised classification is also known as *clustering* [1]. Clustering divides the data into groups of similar objects and each group consists of objects that are similar between themselves and dissimilar to objects of other groups. There are many clustering algorithms that exist such as hierarchical clustering, K-Means clustering, Fuzzy C-Means clustering and many more. The two most popular classes of clustering algorithms are K-Means and Fuzzy C-Means (FCM) [27].
- ii. ***Supervised classification***: On the other hand, in a method of supervised classification, the objects are labelled and each value of the label represents a class. The objective of this method is to build hyperplanes separating the objects according to their class. Supervised classification is performed on a set of training examples. Each training example $\langle x_i, y_i \rangle$ is composed of a feature vector, x_i and a corresponding class label, y_i . The feature vector contains measurable characteristics of the object under consideration.

In this section we are interested in the problem of supervised classification, therefore, the application of unsupervised classification in data mining will not be discussed in further detail. Support Vector Machines, kNearest Neighbor, Decision Tree (i.e.: PART and J48) and Naive Bayes classifiers are in the group of supervised classification and they have been used as state-of-the-art benchmarking classifiers throughout this thesis. The next sections, section 2.2.2.1 to section 2.2.2.4, will briefly describe the basic principle for each of the supervised classifiers and their applications in data classification problems.

2.2.2. Traditional Data Classification Methods

2.2.2.1. Support Vector Machines Classifier

Support Vector Machines (SVMs) are a group of supervised learning methods that can be applied to classification or regression. Theoretically, SVM is a well-motivated algorithm that is based on the statistical learning proposed by Vapnik [28]. SVM has shown promising empirical good performance and successful application in many fields such as bioinformatics, text categorisation, speaker verification, handwritten digit recognition, face detection, engineering and science, financial market evaluation [29], pattern recognition [30], image recognition [31] and many more. SVM has received overwhelming attention as a classification technique from diverse research communities due to its suitability, and works very well with high-dimensional data as well as avoiding the curse of the dimensionality problem [32].

The emergence of SVM in data mining technology has improved the efficiency of data classification that helps people to make decisions. Nevertheless, in real data mining applications, real-life databases sometimes contain outliers or noise data that

lead to the degradation of the classification performance. In order to improve the outlier sensitivity problem of standard SVM for two-class data classification, Yang et al. [33] introduced Weighted Support Vector Machine (WSVM). Different weights generated by Kernel-based Possibilistic C-Means (KPCM) are assigned to different data points such that WSVM learns the decision surface according to the relative importance of the data point in the training set. Furthermore, the training example is usually non-fuzzy where the output is either 1 or 0. However, due to the noise and error of measurement, the training examples are usually uncertain (fuzzy) [34]. Ji et al. [35] introduced the concept of fuzzy linear separable and approximately fuzzy linear separable for training two-class classification. Other works that related to fuzzy training data sets are also presented by Kikuchi and Abe [36] and Lin and Wang [37].

Recently, Mangasarian and Wild [38] proposed Generalized Eigenvalue Proximal SVM (GEPSVM) to perform a binary classification by obtaining two non-parallel hyperplanes, one for each class. In this approach, datapoints of each class are clustered around the corresponding hyperplane. The new datapoints are assigned to a class based on its proximity to one of the two hyperplanes. This formulation leads to two generalised eigenvalue problems, whose solutions are obtained as eigenvectors corresponding to the smallest eigenvalues. In 2007, Twin SVM (TSVM), which is similar in spirit to GEPSVM, was proposed by Jayadeva et al. [39]. The proposed TSVM obtains two non-parallel hyperplanes by solving two novel formulations of Quadratic Programming Problems (QPPs). The idea is to solve two dual QPPs of smaller size rather than solving a single dual QPP with a large number of parameters in conventional SVM. Experimental results in [39] showed the effectiveness of TSVM over GEPSVM and standard SVM on UCI datasets.

SVM was originally designed for binary classification in order to construct an optimal or maximal hyperplane so that the margin of separation between the negative and positive dataset will be maximised. There are many hyperplanes that might classify the data. The best hyperplane is chosen based on the one that represents the largest separation, or margin, between the two classes. So we chose the hyperplane so that the distance from it to the nearest data point on each side is maximised. Figure 2.1 illustrates how SVM finds the unique hyperplane having the maximum margin of linearly separable classes of objects denoted with δ .

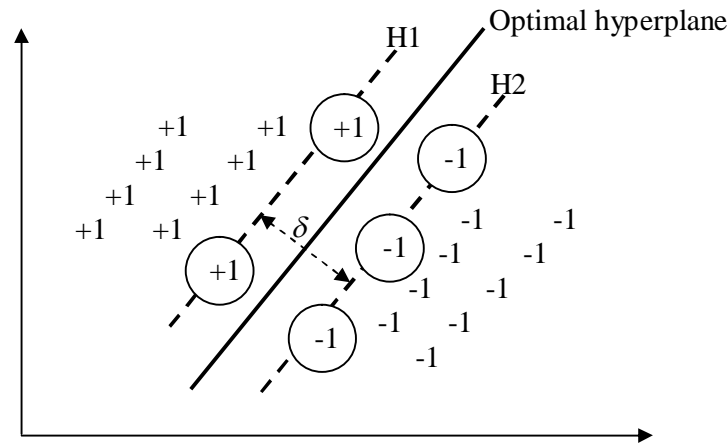


Figure 2.1. Maximum margin hyperplane denoted with δ for linearly separable case [40]

SVM works by mapping the input vectors into a higher dimensional feature space and then obtains the optimal separating hyperplane in the higher dimensional feature space. As depicted in Figure 2.1, two objects from class +1 – defined by the hyperplane H1 – and three objects from class -1 – defined by hyperplane H2 which were represented inside the circle – are called *support vectors*. For classification problems, SVM plots the training vectors in the high-dimensional feature space, and labels each vector by its class. SVM classifies data by determining a set of support

vectors, which are members of the set of training inputs that outline the maximum margin hyperplane in the feature space.

A binary classification problem to find a maximal margin classifier in SVM is presented as the following explanation. Suppose we are given some training data N where each training data is denoted by a pair of (x_i, c_i) . Each x_i is a real number with $x_i = (x_{i1}, x_{i2}, \dots, x_{id})^T$ corresponds to the attribute set for the i^{th} training data, and $i = (1, 2, \dots, N)$. Each object x_i belongs to a class $c_i \in \{-1, +1\}$. Any hyperplane can be written in the following form:

$$w \cdot x_i + b = 0, \quad \text{Eq. 6}$$

where w and b are parameters of the model. A training set of objects is linearly separable if at least one linear classifier exists defined by the pair (w, b) , which correctly classifies all training data as shown in Figure 2.2.

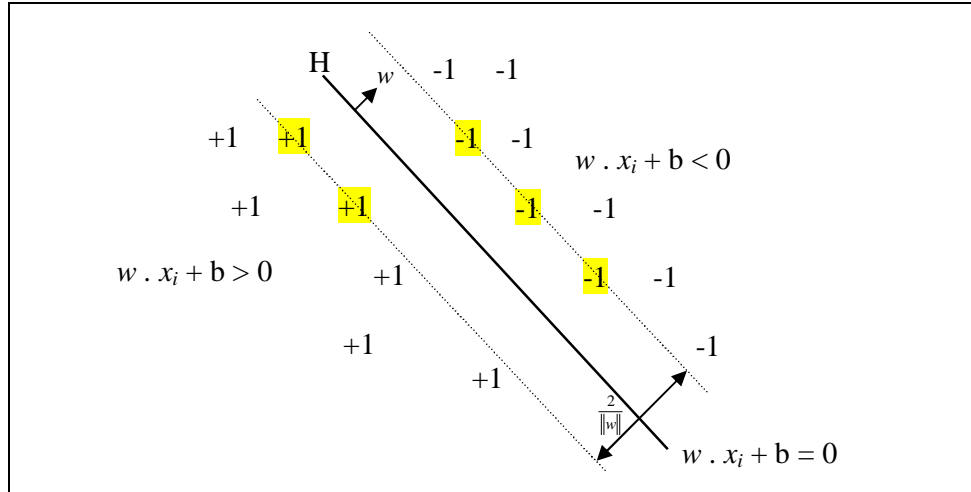


Figure 2.2. The classification hyperplane which defines two linear separable classes in +1 region and -1 region [40]

As the objective of training SVM is to find a pair of (w, b) with the largest margin, the parameters are subject to the following two constraints:

$$\begin{aligned}
w \cdot x_i + b &> 0, \text{ if } c_i = 1 \\
w \cdot x_i + b &< 0 \text{ if } c_i = -1
\end{aligned}
\tag{Eq. 7}$$

In this case, w and b in Eq. 7 can be rewritten in a more compact form as below:

$$c_i(w \cdot x_i + b) \geq 1, i = 1, 2, \dots, N \tag{Eq. 8}$$

The margin separation between these two hyperplanes is defined by (w, b) is $\frac{2}{\|w\|}$ and

therefore we want to minimise $\|w\|$, where $\|w\|$ is the norm of the vector w . However,

it is difficult to solve the optimization problem as it depends on $\|w\|$ that involves

square roots, which is known as a convex optimization problem. Hence, the Lagrange

multiplier is used by substituting $\|w\|$ with $\frac{1}{2}\|w\|^2$ without changing the solution.

Combining the Lagrange multiplier with Eq. 8 produces dual formulation of the optimization as in the following equation:

$$L_D = \sum_i \lambda_i - \frac{1}{2} \sum_{i,j=1}^N \lambda_i \lambda_j c_i c_j x_i \cdot x_j, \tag{Eq. 9}$$

where $\lambda_i \geq 0$ and under constraint $\sum_{i=1}^N c_i \lambda_i = 0$.

In a case of binary classification, SVM can easily find a hyperplane that separates two-dimensional objects which belong to two classes. However, in a real world problem, the dataset of interest usually cannot be separated with a linear classifier as shown in Figure 2.3(a). Therefore, SVM can solve this difficulty by implementing special non-linear functions to non-linearly map the input data to a high dimensional feature space as illustrated in Figure 2.3(b). These non-linear functions are called *kernels*, which are computed by feature functions denoted with ϕ . Here, the linear

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