Random Quasi Intersections with Applications to Instant Machine Learning

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Abstract: Random quasi intersections method was introduced. The number of such intersections grows exponentially with the increasing amount of pattern features, so that a non-polynomial problem in some machine learning applications emerges. However, the paper experimentally shows that randomness allows finding solutions to some visual machine learning tasks using a random quasi intersection-based fast procedure delivering 100% accuracy. Also, the paper discusses implementation of instant learning, which is, unlike deep learning, a non-iterative procedure. The inspiration comes from search methods and neuroscience. After decades of computing only one method was found able to deal efficiently with big data, - this is indexing, which is at the heart of both Google-search and large-scale DNA processing. On the other hand, it is known from neuroscience that the brain memorizes combinations of sensory inputs and interprets them as patterns. The paper discusses how to best index the combinations of pattern features, so that both encoding and decoding of patterns is robust and efficient.

1 INTRODUCTION

Most machine learning methods make use of iterative training, gradient descend and many adaptive coefficients resulting in slow learning. At the same time, human brain can often memorize new visual patterns at first glance. This paper investigates how to build non-iterative, instant learning by partly borrowing ideas from text search engines (Brin and Page, 1998). For some previous work on numerical data indexing please refer to (Sivic and Zisserman, 2009, Mikhailov et al., 2017, Mikhailov and Karavay, 2021). However, the novelty of the approach discussed in this paper comes from the suggested method of random quasi intersections, which makes indexing of noisy, numerical patterns a possibility. Random quasi intersections (hereinafter referred to as quasi intersections) emerge as a result of random pairing of **close** elements of two sets that are defined in a metric space. Depending on applications, quasi intersections can aid machine learning leading up to 100% accurate classification. Besides, quasi intersections allow for pattern inversion that drastically cuts down the computational complexity leading to instant non-iterative learning. Similar to deep learning, quasi intersections can work on raw features for some computer vision application. Whereas the classical intersection size of sets X, Y is defined as the number of pairs of identical elements

$$|X \bigcap Y| = |\{(x \in X, y \in Y): x = y\}|$$

the quasi intersection size is defined as the number of pairs of close elements

$$|X \bigcap_e Y| = |\{(x \in X, y \in Y): |x - y| \le e\} \quad (1)$$

The definition (1) simultaneously produces many potential quasi intersections of different sizes between sets X, Y since many combinations of mating pairs are possible.

Example 1. Let $X = \{1, 2\}$, $Y = \{2, 3\}$, $e = 1$. Here, there exist two quasi intersections:

- single pair (2, 2) quasi intersection is all there is since the pair (1, 3) does not qualify: $|1 - 3| > 1$;

- two pair quasi intersection is also possible since both pairs (1, 2), (2, 3) do qualify.

The specific intersection size emerges only after a specific pairing of elements of two sets is completed, that is, one combination of pairs is created. Here, the

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exclusion rule applies, which states that each element in one set cannot pair simultaneously with more than one element from the other set.

Another reason for introduction of the non-standard definition (1) is as follows. In machine learning applications, when patterns are classified in accordance with their similarities, the intersection sizes must be stable. The term e specifies acceptable variability of sets' elements. With classical intersections, even though the identity $|\{x\}\bigcap\{x\}| = |\{x\}|$ is true, a small noise e can render the classical intersection empty: $|\{x\}\bigcap\{x+e\}| = 0$. But, the quasi intersection $\{x\}\bigcap_e\{x+e\}$ can restore the original intersection size $|\{x\}\bigcap_e\{x+e\}| = |\{x\}|$ if elements are properly paired. Then, for instance, the Jaccard similarity $J(X,Y)=|X\bigcap Y|/|X\bigcup Y|$ (Jaccard, 1901) can be used for classification purposes.

Example 2. For intuitively similar sets $X=\{100,150,200\}$ and $Y = \{199, 151, 101\}$, classical intersection size is equal to 0. However the non-standard definition (1) restores the intersection size to $|X\bigcap_e Y| = 3$ at $e=1$.

The main results of the paper are provided in Section 2. The Section 3 describes basics of applications of quasi intersections to machine learning, which is illustrated by two examples. Section 3.3 presents an introductory example of classification of visual patterns and Section 3.4 considers using quasi intersections for 1100 commercial trademarks image classification. Finally, the Appendix Section discusses the basics of pattern inversion techniques that are employed together with the quasi intersections method.

2 RESULTS

a) A novel definition of the quasi intersection was introduced, which is a specific pairing of close elements.
b) At an arbitrarily large e, the number of quasi intersections grows exponentially as $n!$
c) Two pattern recognition experiments of quasi intersection-based machine learning show 100% accuracy achieved at testing (ref. to Section 3.3, Table 1, and Section 3.4, Table 2).

3 QUASI INTERSECTIONS IN MACHINE LEARNING

With the objective of pattern recognition by classification that is based on set similarities, it might seem reasonable to look for maximum or minimum quasi intersection sizes as a measure of set similarity. However, such approach is not feasible. Indeed, for two sets with N elements each, there exist $N!$ ways of pairing the sets' element. This number can easily grow bigger than the number of atoms in the Universe. But, importantly, experiments show that **as the feature set size goes beyond some 50 – 100 features, the quasi intersection-based classification accuracy may approach 100%**. It is a consequence of experimental results showing that the pairing of elements subject to condition $|x-y| \le e$ can often be chosen randomly.

3.1 Classification of Feature Sets

Let $X=\{x\}$ be a set and $X_n=\{x\}_n$, $n=1,...,N$, be a collection of sets, where n-th set represents the n-th pattern class. Also, let the following non-standard intersection Jaccard measure

$$J_e(X,X_n)=|X\bigcap_e X_n|/(|X|+|X_n|-|X\bigcap_e X_n|) \quad (2)$$

be used to evaluate the similarity between patterns X, X_n. Note that $0\le J_e(X,X_n)\le 1$. Then the unknown pattern X is assigned to the class n, such that $J_e(X,X_n)=\max_{i=1,...,N} J_e(X,X_i)$, if $J_e(X,X_n)\ge t$ where t is a threshold at training. Otherwise, a new class name $N=N+1$ is incrementally created and the unknown pattern X, which is now indexed as X_N, becomes its representative. Note that the above procedure amounts to an unsupervised learning.

3.2 Intersections vs. Frequencies

Direct matching of an input pattern X to all patterns X_n, $n=1,...,N$, is computationally expensive. Fortunately, pattern indexing by pattern inversion provides way of calculating occurrence frequencies f_n, $n=1,...,N$, of pattern classes in inverse data, e.g., inverse files, inverse patterns, inverse sets, etc. A strict formal definition of inverse data structures is provided in Appendix Section. The following identity

$$|X\bigcap_e X_n| = f_n,\ n=1,...,N \quad (3)$$

makes possible the replacement of quasi intersection sizes for frequency. The identity is illustrated by the example provided in Figure 1 that shows red (R), green (G) and blue (B) curves intersecting at 3 points, where columns emerge that contain names of these curves. Obviously, the intersection size of G, B and G, R patterns is $|G \cap B| = 2$ and $|G \cap R| = 1$.

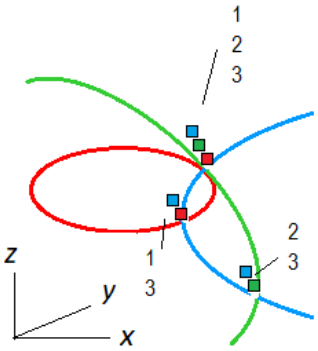

Figure 1: Red (1), green (2) and blue (3) intersecting curves on (x, y)- coordinate plane.

At the same time, a histogram h of names contained in the columns {2, 3} and {1, 2, 3], where the green curve intersects the other lines, has the samples $h(B) = 2$ and $h(R) = 1$.

Hence, $|G \cap B| = h(B)$ and $|G \cap R| = h(R)$,

as it is stated in (3). Importantly, when columns are available, all intersection sizes can be obtained by scanning only three columns (in this example), even though curves may be represented by thousands of dots. Note that the above approach is similar to TF-IDF method (Jones, 2004) developed for document search or information retrieval, where inverse document frequencies of the word across a set of documents are calculated. This is a typical method employed in search engines, where document name frequencies in fully inverted files are calculated.

However, search engines will not work for pattern recognition applications where textual inputs are replaced with noisy numerical data. That is, not just inverse document frequency of the word may change depending on the currently available document collection, but the word itself may completely disappear because of even a small noise.

In case of numeric patterns, the objective is to index each pattern by its features, for instance, the edge pattern on the plain – by (x, y) - coordinates of its dots. Then pattern identities would replace page numbers of the back-of-book-index, where each set $\{n\}_{x, y}$ contains identities of patterns that share the coordinate pair (x, y).

3.3 Introductory Example of Indexing-Based Pattern Recognition

Let consider a pattern recognition example where elements of feature sets are 2D-features, which are (x, y) - coordinates of dark (below a certain grey level) dots, that is, pixel coordinates are used as raw features of image objects. Figure 2 shows an example, where the shapes 1 - 8 are used for training and the distorted shapes 1a - 8a are used for testing. Note that for recognition of distorted objects 1a - 8a the classification system should be invariant with respect to scaling and orientation.

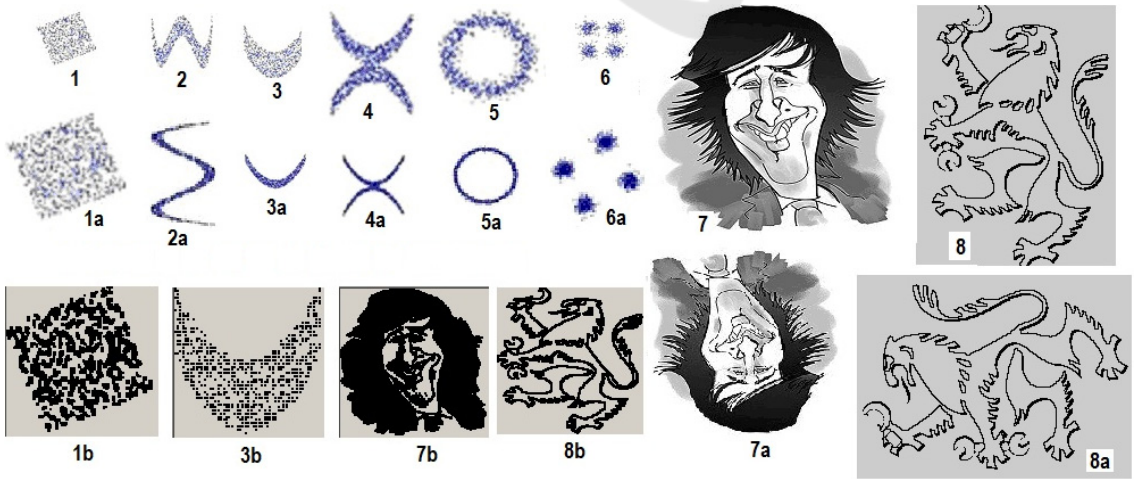

Figure 2: Pictures 1 – 8 (Wikipedia, 2022) were used for learning (single image per pattern class), pictures 1a – 8a (Wikipedia, 2022) are used for testing. Pictures 1b, 3b, 7b, 8b show outputs of a scale normalization algorithm when it takes in pictures 1, 3, 7, 8, respectively.

Invariance with respect to scaling was achieved by the normalization, whereby left-hand, right-hand, upper and lower boundary coordinates $x_{\min}, x_{\max}, y_{\min}, y_{\max}$ were found and (x, y) - coordinates of object's dots were normalized to the range $D = 128$ as follows

$$x = D*(x - x_{\min})/(x_{\max} - x_{\min})$$

$$y = D*(y - y_{\min})/(y_{\max} - y_{\min})$$

Invariance with respect to orientation was achieved by learning the objects 1 – 8 at a single view angle and multiple recognition of objects 1a – 8a at angles 0º through 360°, step 1º.

A notation for inverse patterns was introduced, which is an indexed set $\{n\}_x$ or $\{n\}_{x,y}$, where the set element n is the name of a pattern class and the indexes x, y are the feature values. The notation $\{n\}_{x,y}$ helps to conveniently represent the algorithm that calculates occurrence frequencies f_n, $n = 1,...,N$, of names in columns. Let x_l, y_l, $l = 1,...,L$ be the coordinates of dark dots, where L is the number of dark dots in the current image (image 1 through 8). Then name occurrence frequencies are best obtained as samples of the histogram

Algorithm 1: Name histogram evaluation.

$$l = 1,...,L,\ \forall r,s \in [-e,e],\ \forall n \in \{n\}_{x_l+r,y_l+s}:$$

$$f_n = f_n + 1$$

If classification of the current input image fails, that is, $\max_{i=1,...,N} (f_i)_n < t$, then the new image class is introduced, $N = N + 1$, and the columns are updated using the Algorithm 2.

Algorithm 2: Updating of columns.

$$\{n\}_{x_l,y_l} = \{n\}_{x_l,y_l} \bigcup N,\ l = 1,...,L$$

Here the initial conditions are

$$N = 1,\ \{n\}_{x_l,y_l} = \{1\}_{x_l,y_l},\ l = 1,...,L$$

But, the algorithm 1 cannot work properly as it violates the exclusion rule from Section 1. That is, the element x_l, y_l of the current input pattern may be accidentally compared within the loop $\forall r,s \in [-e,e]$ to more than one element of a previously stored pattern (and visa versa). This is why columns and class names must be inhibited once they have been accessed, which is achieved by replacing histogram algorithm 1 with the algorithm that inhibits patterns by flagging dirty names $flag_n = 0 \rightarrow flag_n = 1$ and columns $col_{x,y} = 0 \rightarrow col_{x,y} = 1$. Then the algorithm 1 can be re-written as Algorithm 3.

Algorithm 3: Histogram evaluation with inhibition.

$$l = 1,...,L,\ \forall n, flag_n = 0, \forall r,s \in [-e,e],$$
$$\forall n \in \{n\}_{x_l+r,y_l+s}$$

if $col_{x_l+r,y_l+s} = 0$ and $flag_n = 0$

$$f_n = f_n + 1,\ col_{x_l+r,y_l+s} = 1,\ flag_n = 1$$

Recognition accuracy (%) for distorted patterns 1a - 8a (Figure 2) at different neighborhood radiuses *e* is provided in Table 1

Table 1: Recognition accuracy %.

e (pixels)	0	1	2	3	4
%	50	100	100	100	87.5

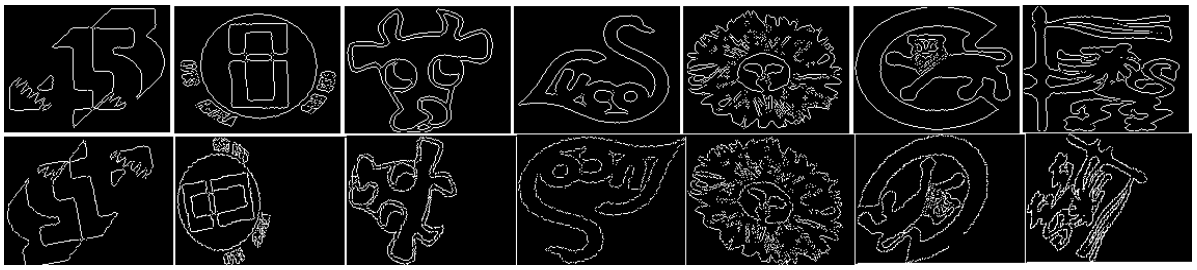

Figure 3: Some typical examples of commercial trademarks. Upper row images were used for training and bottom row images were used for testing.

Table 2: Recognition accuracy at testing of 1100 distorted commercial trademarks retrieved from (GoogleDrive, 2022).

e	0	1	2	3	4
%	98	100	100	100	99.2

3.4 Using Quasi-Intersections for Commercial Trademark Image Database

The same system was also trained on a database of commercial trademarks, retrieved from (GoogleDrive, 2022), which contains 1100 edge images, each comprising 160 x 120 pixels (upper row on the Figure 3 shows a few typical images out of 1100 images from this database). The training was performed at $e = 0$. Next, 1100 distorted images were used for testing (bottom row on the Figure 3 shows typical examples of distorted images and Table 2 shows the recognition accuracy.

The training time was shown to be about 3 seconds per 1000 images (pre-stored in RAM) on a single core 1.6 GHz Intel Pentium CPU, which amounts to 3 milliseconds per image. Although this indexing-based non-iterative learning is impressively fast, the burden of rotational invariance slows down the recognition time by the factor of 360 amounting to about 1 second per image).Pattern clustering may emerge at learning if $e > 0$. But, the discussion of it is beyond the scope of this paper.

4 DISCUSSION

Whereas a classic intersection of two sets always produces a single set, the quasi intersection definition produces a multitude of possible intersections. But, there is no way of knowing in advance which one of them will emerge. Analogously, in quantum mechanic, it is the quantum measurement that localizes a particle, whose potential positions are described by the wave function. And it is impossible to predict, which slit the photon will go through until a detector tells which way the particle had chosen. Besides, the exclusion rule, whereby no two elements of one set can simultaneously pair with one element of the other set, reminds of Pauli Exclusion Principle according to which no two identical fermions in any quantum system can be in the same quantum state.

The discussed instant learning, like deep learning, can work on raw features, which are image pixels, meaning that no feature engineering is needed.

However, the scope of possible applications of instant learning is not known as yet.

Implementation of quasi intersections implies inhibition. Without inhibition the pattern scores easily overflow expected levels, which resemble brain circuitry organization, where excitatory neurons are often accompanied by inhibitory neurons.

The objective of this paper is not a discussion of a best way of implementing an image learning system, but a proof-of-the-concept of quasi intersections method. A better way of image recognition would be a two or more level approach, where level 1 classifies local features and level 2 classifies histograms of classes of local features. Then subsets of histogram samples will represent parts of input objects, which resembles the activation pattern of the inferotemporal brain region (Tsunoda et al., 2001). However, when patterns are represented by feature vectors, rather than feature sets, inhibition is not needed. The basics of feature vectors' indexing are provided in the Appendix section.

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APPENDIX

A.1 Patterns are Represented by Feature Vectors

For feature vectors, there is no need to inhibit accessed columns, even though the non-standard intersection definition (1) is used. Figure A.1 illustrates the above statement.

In Figure A.1 example, *x*-coordinates of colored marks represent component values of red, green and blue 11-dimensional vectors, and *k*-coordinates represent vectors' component indexes. For instance, the cell (15, 1) contains two vector names $\{n\}_{x,k} = \{1,3\}_{15,1}$. The cell (14, 4) contains the name of the 2nd vector and the cell (7, 3) contains names of 1st and 2nd vectors.

The vector coordinate values are

$$x_{red} = (15,13,7,1,17,6,18,3,15,20,11) \quad (1^{st})$$

$$x_{green} = (6,13,7,14,7,17,21,6,20,12,21) \quad (2^{nd})$$

$$x_{blue} = (15,13,3,7,7,7,21,8,4,8,14) \quad (3^{rd})$$

$$x_{black} = (6,13,3,13,7,17,21,8,4,9,13)$$

In K -dimensional space, the similarity between vectors $\vec{x}$, $\vec{x}_n$ can be evaluated as the intersection with the size $|X \bigcap_e X_n| = f_n$, $n = 1,...,N$. For vectors, this is always a single intersection because there exists only one way of counting coordinate pairs. It is an implication of the following two properties of distinct vectors with integer components (D is the range of component values, any cell and its content is referred to as a column). Given N memorized vectors

(a) For each dimension k (for each horizontal line), any two columns do not intersect

$$\{n\}_{x,k} \bigcap \{n\}_{y,k} = \varnothing, \ \forall x, y < D$$

(b) For each dimension k, the sum of column heights is equal to the number N of memorized vectors

$$\sum_{x=0}^{D-1} |\{n\}_{x,k}| = N$$

It follows from (a), (b) that intersection sizes $|\vec{x} \bigcap \vec{x}_n|$ of the input vector $\vec{x} = (x_1,...,x_k,...,x_K)$ with memorized vectors $\vec{x}_n$, $n = 1,...,N$, are equal to frequencies f_n of vector indexes in corresponding columns

$$|\vec{x} \bigcap \vec{x}_n| = f_n, \ n = 1,...,N$$

Computation of frequencies requires a definition of inverse patterns that contain names of pattern classes, feature sets or feature vectors.

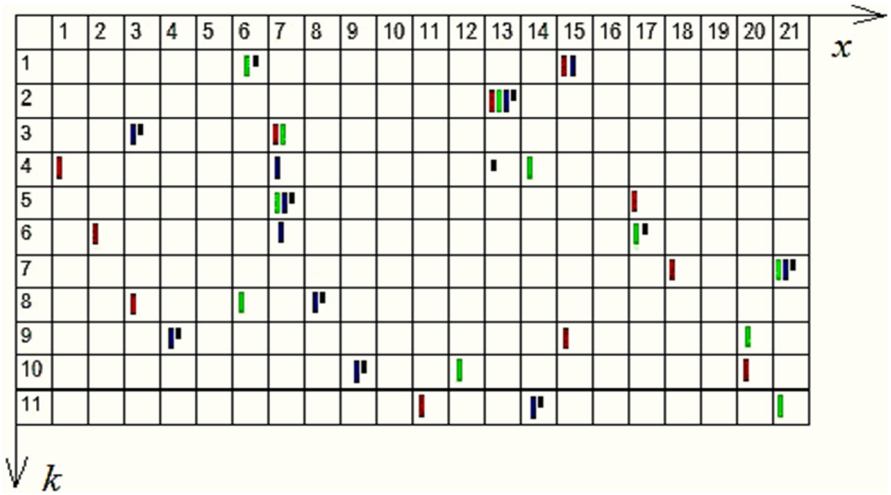

Figure A.1: Feature vectors plain.

A.2 Inverse Sets Definition

Given a collection of distinct sets $X_n = \{x\}_n,\ n = 1,...,N$, inverse set $\{n\}_x$ contains indexes of sets that share the element x

$$\{n\}_x = \{n : x \in X_n\},\ x \in \bigcup_{n=1}^{N} X_n$$

The sets/inverse sets transform

$$\{x\}_n,\ n = 1,...,N \ \leftrightarrow\ \{n\}_x,\ x \in \bigcup_{n=1}^{N} X_n$$

is a one-to-one correspondence.

Example. Given a collection of two sets $\{a,b,c\}_1,\ \{b,c,d\}_2$, inverse sets are

$$\{1\}_a,\ \{1,2\}_b,\ \{1,2\}_c,\ \{2\}_d$$

A.3 Inverse Patterns Definition

Let vectors in a K-dimensional space $x_{n,k},\ k = 1,...,K,\ n = 1,...,N$, differ in, at least, one component $\forall n,m\ \exists k :\ |x_{n,k} - x_{m,k}| > e$. Given N such vectors, the inverse patterns are defined as

$$\{n\}_{x,k} = \{n :\ |x - x_k| \le e\},\ x \in D,\ k = 1,...,K$$

Here D is the range of vector components' values. Calculation of frequencies can be done by scanning the content of K columns only, – even though millions of base vectors may be given. This replacement of similarities for frequencies minimizes computational complexity of classification. Frequencies are best obtained as samples of the histogram of names found in inverse sets. In vector case

$$\forall k,\ \forall r \in [-e,e],\ \forall n \in \{n\}_{x_k + r,\,k} : f_n = f_n + 1 \quad \text{(A.1)}$$

Here $\vec{x} = (x_k,\ k = 1,...,K)$ is the input vector. The input vector $\vec{x}$ is assigned to the most similar, that is, most frequent class

$$n : f_n = \max_{i=1,...,N}(f_i), \quad \text{if } f_n > t \quad \text{(A.2)}$$

where t is the recognition threshold. Properties (a), (b) imply that, in vector case, maximal frequency cannot exceed K.

A.4 Unsupervised Learning

Unsupervised learning starts with recognition. If input vector $\vec{x}$ is successfully recognized using the algorithm (A.1) and the expression (A.2), then proceed to next input vector. If recognition fails, then the number of classes is increased by 1 and the collection of inverse patterns is updated by inserting into the collection the name of the new class

1) $N = N + 1$

2) $\{n\}_{x_k,\,k} = \{n\}_{x_k,\,k} \cup N,\ k = 1,...,K$

Initial conditions are: $N = 1$,

$\{n\}_{x_k,\,k} = \{1\}_{x_k,\,k}$, $k = 1,...,K$

Once the learning is done, the columns are no longer updated. In case of supervised learning, the teacher can introduce a look-up table $m = class(n),\ n = 1,...,N$ that relates created and given classes.

A.5 Computational Complexity

Computational complexity of the frequency algorithm (A.1) is proportional to the average height of non-empty columns

$$h = \frac{1}{|C|}\sum_{(x,k)\in C} |\{n\}_{x,k}|,\ C = \{(x,k) : \{n\}_{x,k} \neq \varnothing\}$$

In turn, column height is inverse proportional to variability radius e. The greater the radius e, the bigger the generalization ability. As a result, less classes N will be created, the average column height will be reduced, which, in turn, decreases the computational time. But, on the other hand, this positive trend is limited by the declining discriminating power of the system.