# **Metric Incremental Clustering of Nominal Data**

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#### **Abstract**

We present an algorithm for clustering nominal data that is based on a metric on the set of partitions of a finite set of objects; this metric is defined starting from a lower valuation of the lattice of partitions. The proposed algorithm seeks to determine a clustering partition such that the total distance between this partition and the partitions determined by the attributes of the objects has a local minimum. The resulting clustering is quite stable relative to the ordering of the objects.

### 1 Introduction

Clustering is an unsupervised learning process that partitions data such that similar data items are grouped together in sets referred to as clusters. This activity is important for condensing and identifying patterns in data. Despite the substantial effort invested in researching clustering algorithms by the data mining community, there are still many difficulties to overcome in building clustering algorithms. Indeed, as pointed in [12] "there is no clustering technique that is universally applicable in uncovering the variety of structures present in multidimensional data sets".

In this paper we focus on an incremental clustering algorithm that can be applied to nominal data, that is, to data whose attributes have no particular natural ordering. In general clustering, objects to be clustered are represented as points in an n-dimensional space  $\mathbb{R}^n$  and standard distances, such as the Euclidean distance is used to evaluate similarity between objects. For objects whose attributes are

nominal (e.g., color, shape, diagnostic, etc.), no such natural representation of objects is possible, which leaves only the Hamming distance as a dissimilarity measure, a poor choice for discriminating among multi-valued attributes of objects.

Incremental clustering has attracted a substantial amount of attention starting with Hartigan's algorithm [11] implemented in [6]. A seminal paper by D. Fisher [10] contained COBWEB, an incremental clustering algorithm that involved restructurings of the clusters in addition to the incremental additions of objects. Incremental clustering related to dynamic aspects of databases were discussed in [4, 5]. It is also notable that incremental clustering has been used in a variety of applications [13, 14, 7, 9]. The interest in incremental clustering stems from the fact that the main memory usage is minimal since there is no need to keep in memory the mutual distances between objects and the algorithms are scalable with respect to the size of the set of objects and the number of attributes.

An object system is a pair S = (S, H), where S is set called the set of objects of S,  $H = \{A_1, \ldots, A_m\}$  is a set of mappings defined on S. We assume that for each mapping  $A_i$  (referred to as an attribute of S) there exists a nonempty set  $E_i$  called the domain of  $A_i$  such that  $A_i : S \longrightarrow E_i$  for  $1 \le i \le m$ . The value of an attribute  $A_i$  on an object t is denoted by  $t[A_i]$ . Our terminology is consistent with the terminology used in relational databases, where a table can be regarded as an object system; however, the notion of object system is more general because objects have an identity as members of the set S, instead of being regarded as just m-tuples of values. In this spirit, we shall refer to  $t[A_i]$  as projection of t on  $A_i$ .

Let S be a set. A partition on S is a non-empty collection of subsets of S indexed by a set I,  $\pi = \{B_i \mid i \in I\}$  such



that  $\bigcup_{i \in I} B_i = S$  and  $i \neq j$  implies  $B_i \cap B_j = \emptyset$ . The sets  $B_i$  are commonly referred to as the *blocks of the partition*  $\pi$ . The set of partitions on S is denoted by  $\mathsf{PART}(S)$ .

PART(S) can be naturally equipped with a partial order. For  $\pi, \sigma \in \mathsf{PART}(S)$  we write  $\pi \leq \sigma$  if every block B of  $\pi$  is included in a block of  $\sigma$ , or equivalently, if every block of  $\sigma$  is an exact union of blocks of  $\pi$ . This partial order generates a lattice structure; this means that for every  $\pi, \pi' \in \mathsf{PART}(S)$  there is a least partition  $\pi_1$  such that  $\pi \leq \pi_1$  and  $\pi' \leq \pi_1$  and there is a largest partition  $\pi_2$  such that  $\pi_2 \leq \pi$  and  $\pi_2 \leq \pi'$ . The first partition is denoted by  $\pi \vee \pi'$ , while the second is denoted by  $\pi \wedge \pi'$ .

An attribute A of an object system  $\mathbb{S}=(S,H)$  generates a partition  $\pi^A$  of the set of objects S, where two objects belong to the same block of  $\pi^A$  if they have the same projection on A. We denote by  $B_a^A$  the block of  $\pi^A$  that consists of all tuples of S whose A-component is a. Note that for relational databases,  $\pi^A$  is the partition of the set of rows of a table that is obtained by using the **group by** A option of **select** in standard SOL.

A clustering of an object system S = (S, H) is defined as a partition  $\kappa$  of S. We seek to find clusterings starting from their relationships with partitions induced by attributes. As we shall see, this is a natural approach for nominal data.

The mapping  $v: \mathsf{PART}(S) \longrightarrow \mathbb{R}$  by  $v(\pi) = \sum_{i=1}^n |B_i|^2$ , where  $\pi = \{B_1, \dots, B_n\}$  is a lower valuation on  $\mathsf{PART}(S)$ , that is,

$$v(\pi \vee \sigma) + v(\pi \wedge \sigma) > v(\pi) + v(\sigma) \tag{1}$$

for  $\pi, \sigma \in \mathsf{PART}(S)$ . For every lower valuation v the mapping  $d: (\mathsf{PART}(S))^2 \longrightarrow \mathbb{R}$  defined by  $d(\pi, \sigma) = v(\pi) + v(\sigma) - 2v(\pi \wedge \sigma)$  is a metric on  $\mathsf{PART}(S)$  (see [2, 1, 15]). A special property of this metric allows the formulation of an incremental clustering algorithm.

# 2 AMICA - A Metric Incremental Clustering Algorithm

Let S=(S,H) be an object system. We seek a clustering  $\kappa=\{C_1,\ldots,C_n\}\in\mathsf{PART}(S)$  such that the total distance from  $\kappa$  to the partitions of the attributes:  $D(\kappa)=\sum_{i=1}^n d(\kappa,\pi^{A_i})$  is minimal. The definition of d allows us to write:

$$d(\kappa,\pi^A) = \sum_{i=1}^n |C_i|^2 + \sum_{j=1}^{m_A} |B_{a_j}^A|^2 - 2\sum_{i=1}^n \sum_{j=1}^{m_A} |C_i \cap B_{a_j}^A|^2,$$

Suppose now that t is a new object,  $t \notin S$ , and let  $Z = S \cup \{t\}$ . The following cases may occur:

- 1. the object t is added to an existing cluster  $C_k$ ;
- 2. a new cluster,  $C_{n+1}$  is created that consists only of t.

Also, from the point of view of partition  $\pi^A$ , t is added to the block  $B_{t[A]}^A$ , which corresponds to the value t[A] of the A-component of t.

In the first case let:

$$\kappa_{(k)} = \{C_1, \dots, C_{k-1}, C_k \cup \{t\}, C_{k+1}, \dots, C_n\}$$
  
$$\pi^{A'} = \{B_{a_1}^A, \dots, B_{t[A]}^A \cup \{t\}, \dots, B_{a_{m-k}}^A\}$$

be the partitions of Z. Now, we have:

$$\begin{split} d(\kappa_{(k)}, \pi^{A'}) - d(\kappa, \pi^A) \\ &= (|C_k| + 1)^2 - |C_k|^2 + (|B_{t[A]}^A| + 1)^2 \\ &- |B_{t[A]}^A|^2 - 2(2|C_k \cap B_{t[A]}^A| + 1) \\ &= 2|C_k| + 1 + 2|B_{t[A]}^A| + 1 - 4|C_k \cap B_{t[A]}^A| - 2 \\ &= 2|C_k \oplus B_{t[A]}^A|. \end{split}$$

The minimal increase of  $d(\kappa_{(k)}, \pi^{A'})$  is given by:

$$\min_{k} \sum_{A} 2|C_k \oplus B_{t[A]}^A|.$$

In the second case we deal with the partitions:

$$\kappa' = \{C_1, \dots, C_n, \{t\}\}\$$

$$\pi^{A'} = \{B_{a_1}^A, \dots, B_{t[A]}^A \cup \{t\}, \dots, B_{a_{m_A}}^A\}$$

and we have  $d(\kappa',\pi^{A'})-d(\kappa,\pi^{A})=2|B_{t[A]}^{A}|.$  Consequently,

$$D(\kappa') - D(\kappa) = \begin{cases} 2 \cdot \sum_A |C_k \oplus B_{t[A]}^A| & \text{in Case 1} \\ 2 \cdot \sum_A |B_{t[A]}^A| & \text{in Case 2.} \end{cases}$$

Thus, if  $\min_k \sum_A |C_k \oplus B^A_{t[A]}| < \sum_A |B^A_{t[A]}|$  we add t to a cluster  $C_k$  for which  $\sum_A |C_k \oplus B^A_{t[A]}|$  is minimal; otherwise, we create a new one-object cluster.

Incremental clustering algorithms are affected, in general, by the order in which objects are processed by the clustering algorithm. Moreover, as pointed in [8], each such algorithm proceeds typically in a hill-climbing fashion that yields local minima rather than global ones. For some incremental clustering algorithms certain object orderings may result in rather poor clusterings. To diminish the ordering effect problem we expand the initial algorithm by adopting the "not-yet" technique introduced by Roure and Talavera in [16]. The basic idea is that a new cluster is created only when the inequality:

$$r(t) = \frac{\sum_{A} |B_{t[A]}^{A}|}{\min_{k} \sum_{A} |C_{k} \oplus B_{t[A]}^{A}|} < \alpha,$$

is satisfied, that is, only when the effect r(t) of adding the object t on the total distance is significant enough. Here  $\alpha$ 



is a parameter provided by the user, such that  $\alpha <= 1$ . Note that if  $\alpha = 1$ , we make no use of the NOT-YET buffer.

We formulate now a metric incremental clustering algorithm (referred to as AMICA – an acronym of the previous five words) that is using the properties of distance d. The variable nc denotes the current number of clusters. If  $\alpha < r(t) \le 1$ , then we place the object t in a NOT-YET buffer. If  $r(t) \le \alpha$  a new cluster that consists of the object t is created. Otherwise, that is if r(t) > 1, the object t is placed in an existing cluster  $C_k$  that minimizes  $\sum_A |C_k \oplus B_{t[A]}^A|$ ; this limits the number of new singleton clusters that would be otherwise created. After all objects of the set S have been examined, the objects contained by the NOT-YET buffer are processed with  $\alpha=1$ . This prevents new insertions in the buffer and results in either placing these objects in existing clusters or in creating new clusters. The pseudocode of the algorithm is given next:

```
Input:
            data set S and threshold \alpha
Output:
              clustering C_1, \ldots, C_{\rm nc}
Method:
nc = 0;
\ell=1;
while S \neq \emptyset do
         select an object t;
         S = S - \{t\};
         if \sum_{t \in A} |B_{t[A]}^A| \le \alpha \min_{1 \le k \le \text{nc}} \sum_{A} |C_k \oplus B_{t[A]}^A|
                  create a new single-object
              cluster C_{nc} = \{t\};
                  r(t) = \frac{\sum_{A} |B_{t[A]}^A|}{\min_{1 \le k \le \text{nc}} \sum_{A} |C_k \oplus B_{t[A]}^A|}
         if r(t) > 1
              then
                  k = \arg\min_k \sum_A |C_k \oplus B_{t[A]}^A|
                  add t to cluster C_k;
              else /* this means \alpha < r(t) \le 1 */
                  place t in NOT-YET buffer;
         end if;
endwhile;
process objects in the NOT-YET buffer
as above with \alpha = 1;
```

## 3 Experimental Results

We applied AMICA to synthetic data sets produced by an algorithm that generates clusters of objects having realnumbered components grouped around a specified number of centroids. The resulting tuples were discretized using a specified number of discretization intervals which allowed us to treat the data as nominal. The experiments were applied to several data sets with an increasing number of tuples and increased dimensionality and using several permutations of the set of objects. All experiments describe in this paper used  $\alpha=0.95$ .

The stability of the obtained clusterings is quite remarkable. For example, in an experiment applied to a set that consists of 10,000 objects (grouped by the synthetic data algorithm around 6 centroids) a first pass of the algorithm produced 11 clusters; however, most objects (9895) are concentrated in the top 6 clusters, which approximate very well the "natural" clusters produced by the synthetic algorithm.

The next table compares the clusters produced by the first run of the algorithm with the cluster produced from a data set obtained by applying a random permutation.

Initial Run		Random Permutation				
Cluster	Size	Cluster Size Distribution				
				(Original cluster)		
1	1548	1	1692	1692 (2)		
2	1693	2	1552	1548 (1), 3 (3), 1 (2)		
3	1655	3	1672	1672 (5)		
4	1711	4	1711	1711 (4)		
5	1672	5	1652	1652 (3)		
6	1616	6	1616	1616 (6)		
7	1	7	85	85 (8)		
8	85	8	10	10 (9)		
9	10	9	8	8 (10)		
10	8	10	1	1 (11)		
11	1	11	1	1 (7)		

Note that the clusters are stable; they remain almost invariant with the exception of their numbering. Similar results were obtained for other random permutations and collections of objects.

As expected with incremental clustering algorithms, the time requirements scale up very well with the number of tuples. On an IBM T20 system equipped with a 700 MHz Pentium III and with a 256 MB RAM, we obtained the following results for three randomly chosen permutations of each set of objects.

Number of	T	Average					
objects	perm	time (ms)					
2000	131	140	154	141.7			
5000	410	381	432	407.7			
10000	782	761	831	794.7			
20000	1103	1148	1061	1104			

Another series of experiments involved the application of the algorithm to databases that contain nominal data. We applied AMICA to the mushroom data set from the standard UCI data mining collection (see [3]). The data set contains 8124 mushroom records and is typically used as



test set for classification algorithms. In classification experiments the task is to construct a classifier that is able to predict the poisonous/edible character of the mushrooms based on the values of the attributes of the mushrooms. We discarded the class attribute (poisonous/edible) and applied AMICA to the remaining data set. Then, we identified the edible/poisonous character of mushrooms that are grouped together in the same cluster. This yields the clusters  $C_1, \ldots, C_9$ :

Cl.	Poisonous/Edible	Total	Percentage of
num.			dominant group
1	825/2752	3577	76.9%
2	8/1050	1058	99.2%
3	1304/0	1304	100%
4	0/163	163	100%
5	1735/28	1763	98.4%
6	0/7	7	100%
7	0/192	192	100%
8	36/16	52	69%
9	8/0	8	100%

Note that in almost all resulting clusters there is a dominant character, and for five out of the total of nine clusters there is complete homogeneity.

A study of the stability of the clusters similar to the one performed for synthetic data shows the same stability relative to input orderings as follows from the next table that describe a clustering obtained under a randomly chosen permutation of the set of objects:

$C_i$	Computed Clusters									
	First Random Permutation									
	$C'_1$	$C_2'$	$C_3'$	$C_4'$	$C_5'$	$C_6'$	$C_7'$	$C_8'$	$C_9'$	$C'_{10}$
	3540	1797	1095	192	1296	8	36	7	137	16
3577	3540	0	37	0	0	0	0	0	0	0
1058	0	0	1058	0	0	0	0	0	0	0
1304	0	8	0	0	1296	0	0	0	0	0
163	0	26	0	0	0	0	0	0	137	0
1763	0	1763	0	0	0	0	0	0	0	0
7	0	0	0	0	0	0	0	7	0	0
192	0	0	0	192	0	0	0	0	0	0
52	0	0	0	0	0	0	36	0	0	16
8	0	0	0	0	0	8	0	0	0	0

Note that the previous table contains mostly zeros. This shows that the clusters remain essentially stable under input data permutations (with the exception of the order in which they are created).

## 4 Conclusion and Future Work

AMICA provides good quality, stable clusterings for nominal data, an area of clustering that is less explored than the standard clustering algorithms that act on ordinal data. Clusterings produced by the algorithm show a rather low sensitivity to input orderings.

Further investigations in the behavior of the algorithm are warranted. For example, we ran AMICA with a rather high value of the threshold  $\alpha=0.95$ . Future work will

include an examination of the dependency of the maximal size of the NOT-YET buffer for various values of  $\alpha$ .

AMICA could be combined with special discretization algorithms such as metric discretization [17] to obtain a more general incremental clustering algorithm applicable to mixed data, that is, to data having both nominal and ordinal attributes. This is currently work in progress.

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