# Efficient Calculation of Interaction Features on Sparse Matrices

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#### Abstract

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### 1 Introduction

Introduction Interaction features are a way of capturing correlations between features in a machine learning setting. A feature vector  $\vec{x}$  of dimensionality D has second degree interaction features  $\{x_a \cdot x_b : a, b \in \{0, 1, ..., D-1\} \land a < b\}$ , so a D dimensional vector has  $\binom{D}{2} = \frac{D^2 - D}{2}$  second degree interaction features. A naive approach to calculating these features is to simply iterate through the combinations of the column indices. For a sparse vector, many of the resulting interaction features would be zero, and could therefore be ignored. This work describes a method to efficiently calculate second degree interaction features for a sparse matrix that has time and space complexities that decrease quadratically with the density of the input matrix with respect to the naive approach.

# 2 Approach

Let the list of nonzero columns for a given row  $\vec{x}$  be denoted by  $N_{zc}$ . The nonzero second degree interaction features are simply the products of all combinations of two elements whose columns are in  $N_{zc}$ . However, to properly place an interaction feature into the correct column, a mapping from the column index pairs of  $N_{zc}$  into the columns of the interaction matrix is

needed. The mapping is from the space (a, b) where a,b are in 1, 2, ..., D onto the space  $1, 2, ..., \frac{D^2 - D}{2}$ . This is isomorphic to mapping the coordinates of the upper triangle of a matrix onto a flat list. The following is a proof by construction for such a mapping.

#### INSERT JOHN'S PROOF HERE

With this mapping, an algorithm for generating second degree interaction features on a matrix A can be formulated as follows:

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Sparse Interaction(A)  \max(b,a) = \frac{2Da - a^2 + 2b - 3a - 2}{2}   B = \text{Compressed Sparse Row Matrix of size } N \times \frac{D^2 - D}{2}   \text{for } row \text{ in } A   N_{zc} = \text{nonzero columns of } row   \text{for each combination } a,b \text{ of elements of } N_{zc}   k = \max(b,a)   i = \text{index of } row   B[i,k] = row[a] \cdot row[b]
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## 3 Complexity Analysis

Assume that A is a matrix with sparsity 0 < d < 1, N rows, and D columns. Finding interaction features with the proposed algorithm has time and space complexity  $O(dND^2)$ , whereas a naive approach of using non-sparse matrices and multiplying all column combinations has time and space complexity  $O(ND^2)$ . The algorithm is therefore an improvement by a factor of the density factor of A.

This can represent a large gain in speed and time. For example, the 20 Newsgroups dataset has density d of 0.12 when its unigrams are represented in a vector space model. This means the proposed approach would take less than  $\frac{1}{8}$  time and memory.

The real benefit of this method is revealed when the average complexity is analysed. The number of interactions calculated for a given row are  $\binom{|N_{zc}|}{2}$ . If the matrix has density d, then on average,  $N_{zc} = Dd$ , so the number of interaction features calculated in total is

$$N\binom{dD}{2} = N\frac{(Dd)!}{2!(Dd-2)!}$$
$$= N\frac{(D^2d^2 - Dd)}{2}$$

This means that the average complexity decreases quadratically with the density.

### 4 Future Work

The approach for generating second degree interaction features required a mapping from combinations of two to the space  $1, 2, ..., \frac{D^2 - D}{2}$ , which is isomorphic to a mapping from the indices of an upper triangular matrix to the indices of a flat list of the same size. To generate third degree interaction features, a mapping from combinations of three (a, b, c) to the space  $1, 2, ... \frac{D^3 - 3D^2 + 2D}{6}$  (which is  $\binom{D}{3}$ ), or the upper 3-simplex of a tensor to a flat list of the same size  $\frac{D^3 - 3D^2 + 2D}{6}$  would be required. In general, for interaction features of degree k, the upper k-simplex of a k-dimensional tensor must be mapped to the space  $1, 2, ... \frac{D!}{k!(D-k)!}$ . A similar approach for finding these mappings could be taken as the one used here for k=2.

Motivation for deriving mapping functions for higher orders of interaction features is that the average complexity of generating degree k interaction features is  $N\binom{Dd}{k}$ , which decreases polynomially with respect to k compared to generating the features naively.