

SVD Factorization for Tall and Fat Matrices on Map/Reduce Architectures

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Abstract

We demonstrate an implementation for an approximate rank- k SVD factorization combining well-known randomized projection techniques with previously implemented map/reduce solutions in order to compute various steps of this procedure, such as local QR and local SVD implementations that can run on a single machine. We structure the problem in a way reduces to single machine Cholesky and SVD factorizations on $k \times k$ matrices, thereby greatly easing the computability of the problem.

1 Introduction

[1] presents many excellent techniques for utilizing map/reduce architectures to compute QR and SVD for the so-called tall-and-skinny matrices. The ideas are based on the fact that QR factorization can be turned into an $A^T A$ computation problem which is easy to compute using map/reduce. First idea is,

$$A^T A = (QR)^T (QR) = R^T Q^T QR = R^T R$$

Then, we take a look at Cholesky factorization of an $n \times n$ symmetric positive definite matrix which is

$$A = LL^T$$

where L is an $n \times n$ lower triangular matrix. R is upper triangular. Then if we factorize A into L and L^T , and $LL^T = RR^T$, we have a method of calculating QR using Cholesky factorization on $A^T A$. The key observation here is that after $A^T A$ computation is completed we will have an $n \times n$

matrix and if A is “skinny” then n is relatively small (in the thousands), and Cholesky decomposition can be executed on this small matrix on a single computer. We can calculate SVD based on QR. SVD decomposition is represented as

$$A = U\Sigma V^T$$

Expand it with $A = QR$

$$QR = U\Sigma V^T$$

$$R = Q^T U \Sigma V^T$$

Let's call $\tilde{U} = Q^T U$

$$R = \tilde{U} \Sigma V^T$$

This means if we run a local SVD on R (we just calculated above with Cholesky) which is an $n \times n$ matrix, we will have calculated \tilde{U} , and the real Σ , and real V^T . Hence we have a map/reduce way of calculating QR and SVD on $m \times n$ matrices where n is small.

1.1 Approximate rank-k SVD

Computing SVD with large n which are “fat” that might have columns in the billions would require reducing the dimensionality of the problem. According to [2], one way to achieve is through random projection. First we draw an $n \times k$ Gaussian random matrix Ω . Then we calculate

$$Y = A\Omega$$

We perform QR decomposition on Y

$$Y = QR$$

Then form $k \times n$ matrix

$$B = Q^T A$$

Then we can calculate SVD on this small matrix

$$B = \hat{U} \Sigma V^T$$

Then form the matrix

$$U = Q\hat{U}$$

The main idea based on

$$A = QQ^T A$$

if replace Q which comes from random projection Y ,

$$A \approx \tilde{Q}\tilde{Q}^T A$$

Q and R of the random projection are close to that of A . In the multiplication above R is called B where $B = \tilde{Q}^T A$, and,

$$A \approx \tilde{Q}B$$

then, as in [1], we can take SVD of B and apply the same transition rules to obtain an approximate U of A .

The reason the approximate works has to do with the fact that projecting points to a random subspace preserves distances between points, or in detail, projecting the n -point subset onto a random subspace of $O(\log n/\epsilon^2)$ dimensions only changes the interpoint distances by $(1 \pm \epsilon)$ with positive probability [3]. It is also said that Y is a good representation of the span of A .

References

- [1] Gleich, Benson, Demmel, *Direct QR factorizations for tall-and-skinny matrices in MapReduce architectures*
- [2] N. Halko, *Randomized methods for computing low-rank approximations of matrices*
- [3] S. Dangupta, A. Gupta *An Elementary Proof of a Theorem of Johnson and Lindenstrauss*