# Leverage Score Sampling

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#### Abstract

Given a matrix  $A \in \mathbb{R}^{m \times n}$ , the column subset selection task aims at the construction of a column submatrix  $C \in \mathbb{R}^{m \times k}$  of A that has favorable spectral properties, for a target rank k. In recent developments, randomized algorithms for the selection of the column subset  $J = [j_1, \ldots, j_k]$ , where each  $j_i \in [n]$  is sampled from a suitable probability distribution defined on the columns of A, has been shown to provably provide good low-rank approximations. In this work, we analyze the leverage scores associated to the columns of A as one such sampling distribution. We introduce the latter as a measure of the importance of each column in composing the range of A and we prove their main properties, including their coherence-revealing features. Motivated by their provable poor performance for full column rank matrices, we then follow the idea of Cohen et al. [2017] and we introduce the ridge leverage scores as a stable and natural procedure to filter out small principal components. To conclude our work, we provide experimental evidence for the superior behaviour of ridge leverage scores sampling with respect to other commonly used distributions on a benchmark Hilbert matrix. In the final section, we draw conclusions about the work and we discuss possible improvements. The code is available at https://github.com/federicobetti99/Low-Rank-Approximation-Techniques. The obtained results are completely reproducible by executing the scripts provided thereby.

#### 1 Introduction

#### Introduction to the problem

Let us consider an arbitrary matrix  $A \in \mathbb{R}^{m \times n}$  with columns  $[a_1, \ldots, a_n]$ . The problem of column subset selection requests a columns index set  $J = \{j_1, \cdots, j_k\}$  such that the column sub-matrix  $C = [a_{j_1}, \ldots, a_{j_k}] \in \mathbb{R}^{m \times k}$  has favorable spectral properties.

**Definition 1.1.** Let  $A \in \mathbb{R}^{m \times n}$  and  $C = [a_{j_1}, \dots, a_{j_k}] \in \mathbb{R}^{m \times k}$  a column submatrix of A. Moreover, let the columns of  $Q \in \mathbb{R}^{m \times k}$  be an orthogonal basis for the range of C. Then, we call the subset  $J = \{j_1, \dots, j_k\}$  a  $(1 + \epsilon)$  factor column subset if

 $||A - QQ^T A||_2^2 \le (1 + \epsilon)||A - \mathcal{T}_k(A)||_2^2, \tag{1}$ 

where  $\mathcal{T}_k(A)$  denotes the best rank-k approximation of A, obtained by projecting onto the k dominant left singular vectors of A.

According to the definition above, the goal is to find a  $(1 + \epsilon)$  factor column subset. We recall that for a generic matrix C (not necessarily a column subset of A), it holds that

$$||A - QQ^T A||_2^2 \le ||A - \mathcal{T}_k(A)||_2^2 + 2||AA^T - CC^T||_2.$$
(2)

Therefore, the task at hand for sketching can be reduced to finding a column subset J for which the corresponding matrix C approximates well enough the range of the original matrix A, which eventually means that the additional term  $||AA^T - CC^T||_2$  is not too large.

#### Literature review

In recent developments, random sampling has been shown to perform well for the column subset selection problem. The main task in this setting is the definition of a suitable probability distribution p on the columns of A such that  $p_j \geq 0$  for  $j \in [n]$  and  $\sum_{j=1}^n p_j = 1$ . Then, for a prescribed rank k, k columns  $a_{j_1}, \ldots, a_{j_k}$  are sampled from the defined distribution p independently and with replacement and rescaled according to  $a_{j_t} = a_{j_t} / \sqrt{kp_{j_t}}$  for  $t = 1, \ldots, k$ , as to guarantee the unbiased estimate  $\mathbb{E}\left[CC^T\right] = AA^T$ , where  $C = [a_{j_1}, \ldots, a_{j_k}] \in \mathbb{R}^{m \times k}$ . Note that the replacement is not needed, as picking the same column twice does not enhance the range of C, but it simplifies the analysis from a theoretical point of view. The rank-k approximation of A is then given by  $QQ^TA$ , where the

columns of  $Q \in \mathbb{R}^{m \times k}$  are an orthogonal basis for the column space of C. Many sampling probabilities have been studied in the literature. In our opinion, Tropp [2009] provides an exhaustive summary. A naive approach would be to pick uniformly at random columns of A, each of them with probability 1/n: this is clearly a sub-optimal choice, as one can diminish the importance sampling probability of the main principal components by simply adding an arbitrary number of zero columns to A which are not contributing in forming the column space. In this work, we mainly consider the latter as a baseline to gain further insight on the advantage of the proposed sampling distribution. On the other hand, standard results in low-rank approximation have shown that sampling proportionally to the column norms of A, namely

$$p_l = \frac{\|a_l\|_2^2}{\|A\|_F^2},\tag{3}$$

minimizes the quantity  $\mathbb{E}\left[\|AA^T - CC^T\|_F^2\right]$  over all possible choices of sampling probabilities p. Note that the distribution (3) remains unchanged under the addition of an arbitrary number of zero columns (because the corresponding column norm is zero), and thus improves with respect to uniform sampling. However, the main drawback for such a sampling strategy is that the importance of a column is based only on its vector norm, and this may be unnecessarily restrictive. As a consequence, one can easily construct counterexamples for which this sampling algorithm gets tricked into selecting a sub-optimal index set J: consider for example a  $n \times 2$  matrix  $A = [Q_1, Q_2\mathbbm{1}_{n-1}]$  where Q is the orthogonal factor of the QR decomposition of a random matrix and  $\mathbbm{1}_{n-1}$  denotes the vector of length n-1 with all components equal to one.

The disadvantage in sampling from the distribution (3) introduced in the previous paragraph leads to the main focus of this work: we discuss how sketching proportionally to the **leverage scores** Cohen et al. [2017] associated to the columns of A will mitigate this issue, while providing additional desirable properties for the underlying probability distribution. In Section 2, we define the **classical leverage scores** associated to the columns of A and we prove their main properties and advantages for their application to sketching algorithms: we show that maximum sampling probability is given to columns which are orthogonal to all the others and that the underlying sampling distribution is invariant to the addition of zero columns. While these are desirable properties, they imply that the leverage scores are all equal to one for a full column rank matrix, and hence their application in this setting reduces to uniform sampling, which performs poorly in practice. To mitigate this issue, in Section 3 we study instead the **ridge leverage scores** which arise as a stable and natural procedure to filter out small principal components by regularization. In Section 4, we show empirically the advantage of the ridge leverage scores with respect to the other sketching distributions introduced in the above by testing the different sampling strategies on a Hilbert matrix of reduced size. Finally, in Section 5 we draw conclusions about the work and we remark possible improvements of the presented work.

#### Notation and preliminary results

For the rest of the work we consider a matrix  $A \in \mathbb{R}^{m \times n} = [a_1, \dots, a_n]$  with  $m \geq n$ . Thus,  $a_i$  denotes the *i*-th column of A. We denote by  $A = U \Sigma V^T$  the singular value decomposition (SVD) of A and  $\sigma_1 \geq \sigma_2 \geq \dots \geq \sigma_n \geq 0$  the singular values of A. We always denote by  $r \leq n$  the rank of A, namely we assume  $\sigma_r > 0$  and  $\sigma_{r+1} = 0$ . For the sake of completeness, we now recall the main properties of the pseudo-inverse which we use in the elaborate. The Moore-Penrose inverse of A, which we always denote by  $A^{\dagger}$ , is defined as the unique matrix  $P \in \mathbb{R}^{n \times m}$  satisfying the Moore-Penrose conditions:

$$APA = A$$
,  $PAP = P$ ,  $(AP)^T = AP$ ,  $(PA)^T = PA$ . (4)

Given an SVD of A, one can show that  $A^{\dagger} = V \Sigma^{\dagger} U^{T}$ , where  $\Sigma^{\dagger} = \text{diag}(\sigma_{1}^{-1}, \dots, \sigma_{r}^{-1}, 0, \dots, 0)$ , always denoting r = rank(A). Moreover, the following useful properties hold:

$$(A^{\dagger})^T = (A^T)^{\dagger}, \quad (A^{\dagger})^{\dagger} = A, \quad (AA^T)^{\dagger} = (A^T)^{\dagger}A^{\dagger}, \quad A^{\dagger} = A^T(AA^T)^{\dagger}. \tag{5}$$

For space reasons, we refer to Barata and Hussein [2012] for an exhaustive summary with corresponding proofs of the results above and of the definitions. The best rank-k approximation of A given by  $U_kU_k^TA$ , which corresponds to the orthogonal projection of A onto the k dominant left singular vectors, is denoted by  $\mathcal{T}_k(A)$ , where  $\mathcal{T}$  denotes the truncation operator and  $U_k$  is the restriction of U to its first k columns. Finally, we denote by  $e_i$  the i-th canonical basis vector with  $(e_i)_j = 1$  if i = j, 0 otherwise.

In the context of randomized strategies for sketching algorithms, it is always implicitly assumed that the k columns of A contained in C  $\{a_{j_t}\}_{t=1}^k$ , sampled according to some probability distribution p defined for  $i \in [n]$ , are properly re-scaled according to  $a_{j_t} = a_{j_t} / \sqrt{kp_{j_t}}$ , as to guarantee the unbiased estimate  $\mathbb{E}\left[CC^T\right] = AA^T$ .

# 2 Leverage Scores

The leverage scores associated to the columns of A are used ubiquitously as importance sampling probabilities for matrix sketching and low-rank approximations. In the literature, they are often introduced as the minimum norm solution to a least squares problem. We find that this interpretation gives a nice insight, hence this is also our starting point.

**Definition 2.1.** (Leverage score). The leverage score associated with the i-th column of A, which we denote as  $l_i(A)$  for the rest of the work, is the solution to the problem

$$\min_{y} ||y||_2^2 \quad such \ that \ a_i = Ay. \tag{6}$$

**Lemma 2.1.** For  $i \in [n]$  it holds that

$$l_i(A) = a_i^T (AA^T)^{\dagger} a_i. (7)$$

*Proof.* The solution  $\hat{y}$  to problem (6) is the minimum norm solution to the least squares problem

$$\min_{y} \|Ay - a_i\|_2^2, \tag{8}$$

in the sense that for any  $\tilde{y} \in \operatorname{argmin}_y ||Ay - a_i||_2^2$ , we have  $||\hat{y}||_2 \leq ||\tilde{y}||_2$ . It is a well known fact, see for example Barata and Hussein [2012], that the minimum norm solution to (8) is given by  $\hat{y} = A^{\dagger}a_i$ , for which

$$\|\hat{y}\|_{2}^{2} = a_{i}^{T}(A^{T})^{\dagger}A^{\dagger}a_{i} = a_{i}^{T}(AA^{T})^{\dagger}a_{i}.$$
(9)

In the last equality of (9), we used the fact that for any matrix A,  $B = A^T$  is a sufficient condition for  $(AB)^{\dagger} = B^{\dagger}A^{\dagger}$  to hold true. Furthermore, the constraint  $A\hat{y} = a_i$  is satisfied because we have

$$A\hat{y} = AA^{\dagger}a_i = AA^{\dagger}Ae_i \stackrel{(*)}{=} Ae_i = a_i,$$

where  $e_i$  is the *i*-th canonical basis vector and at (\*) we used the fact that by definition  $AA^{\dagger}A = A$ .

**Remark 2.1.** An alternative way to see that the constraint is satisfied is to notice that the residual  $||Ay - a_i||_2 = 0$  when taking  $y = e_i$ , being  $e_i$  the i-th canonical basis vector. Hence, this must hold true also for any solution  $\hat{y}$  to (8), as  $||A\hat{y} - a_i||_2 \le ||Ae_i - a_i||_2 = 0$ .

**Definition 2.2.** Let  $A \in \mathbb{R}^{m \times n}$ . We define the **hat matrix** of A as  $H = A^T (AA^T)^{\dagger} A$ .

**Remark 2.2.** We can simplify the expression for the **hat matrix** by noticing that  $A^T(AA^T)^{\dagger} = A^{\dagger}$ . Indeed, by taking  $A = U\Sigma V^T$  an SVD of A, we have

$$A^T(AA^T)^\dagger = V\Sigma^T U^T (U\Sigma\Sigma^T U^T)^\dagger = V\Sigma^T (\Sigma\Sigma^T)^\dagger U^T = V\Sigma^T (\Sigma^T)^\dagger \Sigma^\dagger U^T = V\Sigma^\dagger U^T = A^\dagger U^T = V\Sigma^T U^T = V\Sigma^T$$

Throughout the remaining of the elaborate, we may prefer the former or the latter representation for the hat matrix, depending on convenience. Most importantly, we recognize the hat matrix  $H = A^{\dagger}A$  to be the orthogonal projector onto range( $A^{T}$ ), the row-space of A.

**Lemma 2.2.** The hat matrix H is idempotent, i.e.  $H^2 = H$ , and symmetric, i.e.  $H = H^T$ .

*Proof.* To show that H is idempotent, we compute directly that

$$H^2 = A^{\dagger}AA^{\dagger}A = A^{\dagger}A = H.$$

where we use the fact that  $A^{\dagger}AA^{\dagger}=A^{\dagger}$  by definition of the Moore-Penrose inverse. To show that H is symmetric, let us take  $A=U\Sigma V^T$  the singular value decomposition of A. Then

$$\begin{split} H^T &= (A^\dagger A)^T = (V \Sigma^\dagger U^T U \Sigma V^T)^T \\ &= (V \Sigma^\dagger \Sigma V^T)^T \\ &= V (\Sigma^\dagger \Sigma)^T V^T \\ &= V \Sigma^\dagger \Sigma V^T \\ &= V \Sigma^\dagger U^T U \Sigma V^T = A^\dagger A = H. \end{split} \qquad (U^T U = I_m)$$

Thus, the claim follows.

**Lemma 2.3.** For  $i \in [n]$ , it holds that  $l_i(A) \leq 1$ .

*Proof.* Note that the leverage scores  $\{l_i(A)\}_{i=1}^n$  are the diagonal entries of the hat matrix H, because we can write

$$l_i(A) = a_i^T (AA^T)^{\dagger} a_i = e_i^T A^T (AA^T)^{\dagger} A e_i = e_i^T H e_i = [H]_{ii}.$$

Therefore, we compute directly

$$l_i(A) = [H]_{ii} = e_i^T H e_i \overset{CS}{\leq} \|e_i\|_2 \|H e_i\|_2 \leq \|e_i\|_2 \|H\|_2 \|e_i\|_2 = \|H\|_2 \leq 1.$$

Thus,  $0 \le l_i(A) \le 1$  for  $i \in [n]$ . In the first inequality, we used the fact that  $l_i(A) \ge 0$  by construction of problem (6) and the Cauchy-Schwarz inequality. In the second inequality, we used the fact that  $||Hx||_2 \le ||H||_2 ||x||_2$  for any  $x \in \mathbb{R}^n$ . In the last inequality, we used that for any orthogonal projector P, it holds that  $||P||_2 \le 1$  Golub and Van Loan [2013].

**Remark 2.3.** By construction of the problem (6) which is solved by the i-th leverage score, we can alternatively prove that  $l_i(A) \leq 1$  by noting that

$$l_i(A) = \underset{y:Ay=a_i}{\operatorname{argmin}} \|y\|_2^2 \le \|e_i\|_2^2 = 1,$$

where  $e_i$  is the i-th canonical basis vector. However, we preferred the previous proof as the representation of the leverage scores through the diagonal elements of the hat matrix will often turn out to be useful throughout the work.

Similarly to the sampling distribution (3) for sketching, the sampling distribution induced by the leverage scores (7) remains unchanged in the case of addition of zero columns, because the leverage score for such columns will be trivially zero. This also implies that the normalization constant  $\sum_{i=1}^{n} l_i(A)$  should be invariant under this operation, as it happens for the Frobenius norm  $||A||_F^2$  for the sketching distribution given by (3). We now show that we can compute directly the normalization constant, exploiting again the fact that H is a projection matrix.

**Lemma 2.4.** It holds that  $\sum_{i=1}^{n} l_i(A) = rank(A) \leq n$ .

*Proof.* Let r denote the rank of A. We notice first that  $\sum_{i=1}^{n} l_i(A) = \text{Tr}(H)$ . Letting  $A = U\Sigma V^T$  be the SVD of A, this gives

$$\begin{aligned} \operatorname{Tr}(H) &= \operatorname{Tr}(A^{\dagger}A) \\ &= \operatorname{Tr}(V\Sigma^{\dagger}U^{T}U\Sigma V^{T}) \\ &= \operatorname{Tr}(V\Sigma^{\dagger}\Sigma V^{T}) \\ &= \operatorname{Tr}(V^{T}V\Sigma^{\dagger}\Sigma) \end{aligned} \qquad \text{(if } A = U\Sigma V^{T} \text{ then } A^{\dagger} = V\Sigma^{\dagger}U^{T}) \\ &= \operatorname{Tr}(V^{T}U = I_{m}) \\ &= \operatorname{Tr}(\Sigma^{\dagger}\Sigma) \\ &= \operatorname{Tr}(\Sigma^{\dagger}\Sigma) \end{aligned} \qquad \text{(trace invariance under cyclic permutations)} \\ &= \operatorname{Tr}(\Sigma^{\dagger}\Sigma) \\ &= \operatorname{rank}(A). \end{aligned} \qquad (\Sigma^{\dagger} = \operatorname{diag}(\sigma_{1}^{-1}, \dots, \sigma_{r}^{-1}, 0, \dots, 0))$$

Thus, the claim follows.

With a slight abuse of notation, for the rest of the work we implicitly assume the probability distribution induced by  $\{l_i(A)\}_{i=1}^n$  to be normalized by r = rank(A), unless specified otherwise.

We now show that we can simplify the expression for the leverage scores (7), assuming we have further information about the rank of the matrix A.

**Lemma 2.5.** Let  $A = U\Sigma V^T$  be the singular value decomposition of A. Furthermore, denote by  $V_k$  be the restriction of V to its first k columns. If A has rank k, then  $l_i(A) = \|(V_k^T)_i\|_2^2$ , where  $(V_k^T)_i$  denotes the i-th row of  $V_k$ .

*Proof.* If A has rank k it coincides with its k-truncated SVD, namely  $A = U_k \Sigma_k V_k^T$ , where  $U_k \in \mathbb{R}^{m \times k}$ ,  $\Sigma_k \in \mathbb{R}^{k \times k}$  and  $V_k \in \mathbb{R}^{n \times k}$ . This is because  $\sigma_{k+1} = \ldots = \sigma_n = 0$ . Thus, we compute

$$\begin{split} l_i(A) &= a_i^T (AA^T)^\dagger a_i \\ &= e_i^T A^\dagger A e_i \qquad \qquad \text{(because } A^T (AA^T)^\dagger = A^\dagger \text{)} \\ &= e_i^T V_k \Sigma_k^\dagger U_k^T U_k \Sigma_k V_k^T e_i \qquad \qquad \text{(if } A = U_k \Sigma_k V_k^T \text{ then } A^\dagger = V_k \Sigma_k^\dagger U_k^T \text{)} \\ &= e_i^T V_k \Sigma_k^\dagger \Sigma_k V_k^T e_i \qquad \qquad \text{(} U_k^T U_k = I_k \text{)} \\ &= (V_k^T e_i)^T (V_k^T e_i) = \|(V_k^T)_i\|_2^2. \qquad \qquad \text{(} \Sigma_k^\dagger = \Sigma_k^{-1} \text{ because } A \text{ has rank } k \text{ so } \sigma_k > 0 \text{)} \end{split}$$

Therefore, the claim follows.

**Remark 2.4.** Lemma 2.5 implies that, in the special case in which A has rank k, sketching proportionally to the leverage scores (7) coincides with using the distribution given by

$$p_i = \|(V_k^T)_i\|_2^2 / k, \quad i = 1, \dots, n.$$
 (10)

Drineas et al. [2008] have shown that sampling  $\mathcal{O}(k^2 \log(1/\delta)\epsilon^{-2})$  columns according to (10) gives

$$||A - QQ^T A||_F \le (1 + \epsilon)||A - \mathcal{T}_k(A)||_F \tag{11}$$

with probability  $1 - \delta$ , where  $\mathcal{T}_k(A)$  is the best rank-k approximation of A, obtained by projecting A onto the space spanned by its k dominant left singular vectors.

**Lemma 2.6.** If a column  $a_i$  is orthogonal to all the other columns then  $l_i(A) = 1$ .

*Proof.* If  $a_i = 0$ , we trivially have  $l_i(A) = 0$  which is inconsistent with the claim. Thus, let us assume that  $a_i \neq 0$ , so that  $||a_i||_2 \neq 0$ . Because  $a_i$  is orthogonal to all the other columns,  $u_i = \frac{a_i}{||a_i||_2}$  is a left singular vector of A. To see this, we show equivalently that  $u_i$  is an eigenvector of  $AA^T$  with eigenvalue  $\sigma_i^2 = ||a_i||_2^2$ . We have

$$AA^{T}u_{i} = AA^{T}\frac{a_{i}}{\|a_{i}\|_{2}} = A\begin{bmatrix} -a_{1} & -\\ & \ddots & \\ -a_{i} & -\\ & \ddots & \\ -a_{n} & -\end{bmatrix} \frac{a_{i}}{\|a_{i}\|_{2}} = A\begin{bmatrix} 0\\ & \ddots\\ & 0\\ & \frac{a_{i}^{T}a_{i}}{\|a_{i}\|_{2}}\\ 0\\ & \ddots\\ & 0\end{bmatrix} = A\begin{bmatrix} 0\\ & \ddots\\ & 0\\ & \|a_{i}\|_{2}\\ 0\\ & \ddots\\ & 0\end{bmatrix},$$

where in the third equality we used the fact that  $a_i$  is orthogonal to all the other columns of A (and hence, rows of  $A^T$ ). In the last two passages, the non-zero element is clearly in position i. We conclude that

$$AA^Tu_i = a_i ||a_i||_2 = ||a_i||_2^2 u_i,$$

which shows the claim. Up to permutations of the columns of A we can assume without loss of generality that i = 1. Thus, we write

$$A = \begin{bmatrix} | & \\ a_i & B \\ | & \end{bmatrix},$$

where  $B = [a_2, \ldots, a_n] \in \mathbb{R}^{m \times n - 1}$ . Now, if  $\tilde{U} \tilde{\Sigma} \tilde{V}^T$  denotes the economy-sized SVD of B, namely with  $\tilde{U} \in \mathbb{R}^{m \times (n - 1)}, \tilde{\Sigma} \in \mathbb{R}^{(n - 1) \times (n - 1)}$  and  $\tilde{V} \in \mathbb{R}^{(n - 1) \times (n - 1)}$ , we can derive an SVD for A as

$$A = \begin{bmatrix} | & \\ | & a_i & B \end{bmatrix} = \begin{bmatrix} | & \\ | & u_i || a_i ||_2 & \tilde{U} \tilde{\Sigma} \tilde{V}^T \end{bmatrix} = \begin{bmatrix} | & \\ \frac{a_i}{||a_i||_2} & \tilde{U} \end{bmatrix} \begin{bmatrix} ||a_i||_2 & 0 & \dots & 0 \\ 0 & & \\ \vdots & & \tilde{\Sigma} \tilde{V}^T \end{bmatrix}$$

$$= \begin{bmatrix} | & \\ u_i & \tilde{U} \end{bmatrix} \begin{bmatrix} ||a_i||_2 & 0 & \dots & 0 \\ 0 & & & \\ \vdots & & \tilde{\Sigma} & \end{bmatrix} \begin{bmatrix} 1 & 0 & \dots & 0 \\ 0 & & & \\ \vdots & & \tilde{V}^T & \\ 0 & & & \end{bmatrix}$$

$$= \begin{bmatrix} | & \\ u_i & \tilde{U} \\ | & & \end{bmatrix} \begin{bmatrix} ||a_i||_2 & 0 & \dots & 0 \\ 0 & & & & \\ \vdots & & \tilde{\Sigma} & & \\ 0 & & & & \end{bmatrix} \begin{bmatrix} 1 & 0 & \dots & 0 \\ 0 & & & & \\ \vdots & & \tilde{V} & & \\ 0 & & & & \end{bmatrix}^T := U \Sigma V^T.$$

Let us denote  $k = \operatorname{rank}(A)$ . To show the claim, we note that the elements of the matrix  $\Sigma$  are not yet necessarily in non-increasing order. However, because we assumed that  $a_i \neq 0$ , we have  $||a_i||_2 > 0$ , which equivalently means that if we order the elements of  $\Sigma$  in non-increasing order, such a singular value will belong in the top k diagonal entries of  $\Sigma$ . Hence, when considering the truncation  $\mathcal{T}_k(A) = U_k \Sigma_k V_k^T$ ,  $u_i$  belongs to the k dominant left singular vectors contained in  $U_k$ , the canonical vector  $e_1$  belongs to the columns of  $V_k$  and the corresponding row is a canonical vector in  $\mathbb{R}^k$ , thus with norm 1. As a consequence, by using Lemma 2.5, we conclude that  $l_i(A) = ||(V_k^T)_i||_2^2 = 1$ , and the claim follows.

**Remark 2.5.** In the proof of Lemma 2.6, we assumed without loss of generality and for simplicity of exposition that  $a_i$  was corresponding to the first column of A. However, as we claimed to conclude the proof, the crucial observation is that when reordering the singular values of A, the entry equal to one which we derived in the previous calculations will still be captured by the rank-k truncation for k = rank(A), and as a consequence the corresponding row of length k is always a canonical vector in  $\mathbb{R}^k$ . Thus, from Lemma 2.5 we conclude that  $l_i(A) = 1$ .

The presence of a column which is orthogonal to all the others (and not identically zero) means that removing such a column would decrease the column rank of A, completely changing its column space. As a consequence, we want to give very high probability to such a column. Lemma 2.6 shows that for such a column the leverage score (7) is maximum. Thus, the **classical leverage score** is a measure of the importance of  $a_i$  in the range of A

Remark 2.6. In the introduction, we claimed that the main drawback of sampling proportionally to the columns norm (3) is the fact that the importance sampling probability for a column is just given by its norm, while ideally the sampling probability of a column should also measure its contribute in the column space of A. The previous result shows precisely that the leverage scores do satisfy this property; namely, the i-th leverage score takes into account the importance of the column  $a_i$  in composing the range of A, for  $i \in [n]$ . As a consequence of this, we naturally expect sketching proportionally to (7) to be superior to sampling from (3).

**Definition 2.3.** The maximum leverage score  $c(A) = \max_{i \in [n]} l_i(A)$  over the columns of A is defined as the coherence of A.

Remark 2.7. Matrix coherence measures the extent to which the singular vectors of a matrix are correlated with the standard basis. If there is a column  $a_i$  which is orthogonal to all the others, then A contains at least one column whose removal would significantly affect the composition of A's column space, and hence has maximum coherence. Sampling accordingly to the column norms (3) may not capture such a feature of A, while the leverage scores carry useful information about the coherence of a matrix, and hence about the presence or not of columns which are much more important than others in composing the range of a matrix A. In this sense, the classical leverage scores are said to be coherence-revealing.

# 3 Ridge Leverage Scores

We concluded Section 2 by claiming that the classical leverage scores are able to reveal the most important columns of A in composing the column space. However, combining Lemma 2.3 and Lemma 2.4 from the above, we conclude that for a full column rank matrix A all the leverage scores are equal to one, because we have  $0 \le l_i(A) \le 1$  for  $i \in [n]$  and  $\sum_{i=1}^n l_i(A) = n$ . Therefore, in this setting, the leverage score sampling reduces to uniform sampling of the columns of A with probability 1/n, which performs poorly in practice. As a first solution, inspired by the result of Lemma 2.5, one can compute rank-k subspace leverage scores McCurdy [2018].

**Definition 3.1.** (Rank-k subspace leverage scores). Let  $A \in \mathbb{R}^{m \times n}$  and  $\mathcal{T}_k(A)$  its best rank k approximation obtained by projecting onto its k principal left singular vectors. The rank-k subspace leverage scores associated to the columns of A are defined for  $i \in [n]$  as

$$l_i^k(A) = a_i^T (\mathcal{T}_k(A)\mathcal{T}_k(A)^T)^{\dagger} a_i.$$
(12)

**Remark 3.1.** By substituting A with its truncated SVD  $\mathcal{T}_k(A)$  in (7) and repeating the same exact calculations carried out in Lemma (2.5), we recover an expression for the rank-k subspace leverage scores given by  $l_i^k(A) = \|(V_k^T)_i\|_2^2$ , i.e. the squared norm of the i-th row of  $V_k$ .

In (12), the classical leverage scores introduced in Section 2 are modified to only capture how important each column  $a_i$  is in composing the top k singular directions of the range of A. This procedure aims at mitigating the small singular values components of  $AA^T$  in the classical leverage scores, which are selected with probability 1/n otherwise, by completely omitting them. However, note that in principle the best rank-k approximation  $\mathcal{T}_k(A)$  may not be well defined. For example, the latter is not unique with respect to the Frobenius norm if  $\sigma_k = \sigma_{k+1}$ ; in the spectral norm, the best rank-k approximation is unique only if  $\sigma_{k+1} = 0$ . Furthermore, even if  $\mathcal{T}_k(A)$  is unique, Cohen et al. [2017] argue that it can be very sensitive to matrix perturbations, so the scores can drastically change when A is slightly modified or when the algorithm does not have access to all the entries of the matrix A, but has only a partial knowledge. As a consequence, they argue that the rank-k subspace leverage scores are not stable and, instead, propose to regularize the problem as a more natural solution.

**Definition 3.2.** (Ridge leverage scores). Let  $A \in \mathbb{R}^{m \times n}$  and let  $\lambda > 0$  be a suitable regularization parameter. The ridge leverage scores associated to the columns of A are defined for  $i \in [n]$  as

$$l_{i,\lambda}(A) = a_i^T (AA^T + \lambda^2 I)^{-1} a_i.$$
(13)

Remark 3.2. Note that in (13), we can take the inverse of  $(AA^T + \lambda^2 I)$  instead of the Moore-Penrose pseudoinverse of the latter. This is because by regularizing the problem, all the singular values of  $AA^T + \lambda^2 I$  are at least  $\lambda^2$ . Furthermore, we recover the orthogonal projector onto range $(A^T)$  given by  $A^{\dagger}A$  in the limit as  $\lambda \to 0$ . Indeed note that, by definition

$$A^T(AA^T)^{\dagger} = A^{\dagger} := \lim_{\lambda \to 0} A^T(AA^T + \lambda^2 I)^{-1}.$$

In particular, if we always denote r = rank(A), it holds that Golub and Van Loan [2013]

$$||A^{\dagger} - A^{T} (AA^{T} + \lambda^{2} I)^{-1}||_{2} \le \frac{\lambda^{2}}{\sigma_{r}(\sigma_{r}^{2} + \lambda^{2})}.$$
 (14)

Hence, we also have

$$|l_{i,\lambda}(A) - l_{i}(A)| = |e_{i}^{T} A^{T} (AA^{T} + \lambda^{2} I)^{-1} A e_{i} - e_{i}^{T} A^{\dagger} A e_{i}|$$

$$= |e_{i}^{T} (A^{T} (AA^{T} + \lambda^{2} I)^{-1} A - A^{\dagger} A) e_{i}| \stackrel{CS}{\leq} ||e_{i}||_{2} ||A^{T} (AA^{T} + \lambda^{2} I)^{-1} A - A^{\dagger} A||_{2} ||e_{i}||_{2}$$

$$= ||A^{T} (AA^{T} + \lambda^{2} I)^{-1} A - A^{\dagger} A||_{2} \leq ||A^{T} (AA^{T} + \lambda^{2} I)^{-1} - A^{\dagger} ||_{2} ||A||_{2}$$

$$\leq \frac{\sigma_{1} \lambda^{2}}{\sigma_{r} (\sigma_{r}^{2} + \lambda^{2})} \rightarrow 0 \quad \text{as } \lambda \rightarrow 0,$$

where in the first inequality we used the Cauchy-Schwarz inequality and in the last inequality we used (14). Thus, the classical leverage scores (7) are recovered in the limit as  $\lambda \to 0$  of the ridge leverage scores (13). Note that, the factor  $\frac{\sigma_1}{\sigma_r}$  corresponds to the conditioning number of the matrix A. Thus, while the result holds true in general for badly conditioned matrices the convergence may be slower.

Remark 3.3. Intuitively, sampling by ridge leverage scores (13) is equivalent to sampling by classical leverage scores (7) for the matrix  $[A, \lambda I_n]$ . However, samples are only taken among the columns of A and not among the ones of the identity matrix. One can show a derivation for (13) similar to the one proved for the classical leverage scores in Lemma 2.1 and involving a regularized version of the least squares problem (8). For space reasons, we omit such a discussion, referring the interested reader to El Alaoui and Mahoney [2014].

Ridge leverage scores take a different approach to mitigate the small singular values components of  $AA^T$  compared to the rank-k leverage scores (12): they only diminish the importance of small principle components through regularization, instead of completely omitting them. We give an intuition for this claim by deriving a simpler expression for the ridge leverage scores.

**Lemma 3.1.** Let  $A = U\Sigma V^T$  be the singular value decomposition of A. For  $i \in [n]$  it holds that

$$l_{i,\lambda}(A) = (V^T e_i)^T \operatorname{diag}\left(\frac{\sigma_1^2}{\sigma_1^2 + \lambda^2}, \dots, \frac{\sigma_n^2}{\sigma_n^2 + \lambda^2}\right) V^T e_i.$$
(15)

*Proof.* We compute

$$\begin{split} l_{i,\lambda}(A) &= a_i^T (AA^T + \lambda^2 I)^{-1} a_i \\ &= e_i^T V \Sigma^T U^T (U \Sigma \Sigma^T U^T + \lambda^2 I)^{-1} U \Sigma V^T e_i \\ &= e_i^T V \Sigma^T U^T (U \Sigma \Sigma^T U^T + U \lambda^2 I U^T)^{-1} U \Sigma V^T e_i \\ &= e_i^T V \Sigma^T U^T U (\Sigma \Sigma^T + \lambda^2 I)^{-1} U^T U \Sigma V^T e_i \\ &= e_i^T V \Sigma^T (\Sigma \Sigma^T + \lambda^2 I)^{-1} \Sigma V^T e_i \\ &= (V^T e_i)^T \operatorname{diag} \left( \frac{\sigma_1^2}{\sigma_1^2 + \lambda^2}, \dots, \frac{\sigma_n^2}{\sigma_n^2 + \lambda^2} \right) V^T e_i. \end{split}$$

The perturbation of  $AA^T$  with  $\lambda^2 I$  allows to regularize the problem and can mitigate the issue presented at the beginning of the section for full column rank matrices. Given the expression worked out in Lemma 3.1 for the ridge leverage scores (13), conside the case of a  $m \times n$  matrix A (always with  $m \ge n$ ) with a large gap between the k-th and (k+1)-th singular values, and suppose to choose the regularization parameter  $\lambda$  such that

$$\sigma_1 \ge \sigma_2 \ge \ldots \ge \sigma_k \gg \lambda \gg \sigma_{k+1} \ge \ldots \ge \sigma_n > 0.$$

Because  $\sigma_n > 0$ , A has full column rank so we already argued previously that  $l_i(A) = 1$  for all  $i \in [n]$ . Hence, sampling from the classical leverage scores (7) is equivalent to uniformly sample columns of A, which performs poorly. On the other hand, for such a regularization parameter  $\lambda$ , the diagonal entries of the matrix

$$\Sigma^{2,\lambda} = \operatorname{diag}\left(\frac{\sigma_1^2}{\sigma_1^2 + \lambda^2}, \dots, \frac{\sigma_n^2}{\sigma_n^2 + \lambda^2}\right)$$
(16)

are satisfying  $\Sigma_{ii}^{2,\lambda} \approx 0$  for  $i \geq k+1$  and  $\Sigma_{ii}^{2,\lambda} \approx 1$  for  $i \leq k$ . Therefore

$$l_{i,\lambda}(A) = \sum_{j=1}^{n} \frac{\sigma_j^2}{\sigma_j^2 + \lambda^2} v_{ij}^2 \approx \sum_{j=1}^{k} v_{ij}^2 = \|(V_k^T)_i\|_2^2,$$
(17)

where  $V_k$  denotes the restriction of V to its first k columns. Therefore, sampling proportionally to the ridge leverage scores  $l_{i,\lambda}(A)$  is approximately equal to using the distribution (10).

**Remark 3.4.** Note that, as we discussed at the beginning of the section, for a similar matrix using the rank-k subspace leverage scores (12) will omit completely the small principal components successive to  $\sigma_k$ . Instead, with ridge leverage scores sampling we assign to these principal components a non-zero probability, which is decreasing with  $\lambda$ .

From this example, we see that the regularization parameter becomes to all effects a hyper-parameter which needs to be tuned, depending on the desired cutoff effect on the columns of V. A too large value of  $\lambda$  may drastically change the spectral properties of the middle factor  $AA^T$ , while a too small  $\lambda$  may not regularize sufficiently the problem, thus recovering the rank-k subspace leverage scores (12). To our knowledge, for a general matrix A, the choice of  $\lambda$  is still an unresolved task. In the pioneering work on ridge leverage scores sampling, Cohen et al. [2015] propose a wise choice of  $\lambda$  given by  $\lambda^2 = \frac{\|A - \mathcal{T}_k(A)\|_F^2}{k}$ .

Remark 3.5. Note that while  $\mathcal{T}_k(A)$ , used to compute the rank-k subspace leverage scores (12), may not be uniquely determined,  $\lambda = \frac{\|A - \mathcal{T}_k(A)\|_F}{\sqrt{k}}$  is uniquely determined because by the Von-Neumann trace inequality we can lower bound the best squared rank-k approximation error in the Frobenius norm with  $\sum_{j=k+1}^n \sigma_j^2$ . In particular, while the ridge scores (13) with this choice of regularization parameter depend on the value of  $\|A - \mathcal{T}_k(A)\|_F^2$ , they do not depend on a specific low-rank approximation. This is sufficient for stability since by the Weyl's inequality Marcus  $[1965] \|A - \mathcal{T}_k(A)\|_F^2$  changes predictably under matrix perturbations even when  $\mathcal{T}_k(A)$  itself does not.

Note that computing both the rank-k subspace leverage scores (12) and the ridge scores (13) with  $\lambda^2 = \frac{\|A - \mathcal{T}_k(A)\|_F^2}{k}$  may be computationally too expensive or even simply impossible in practice. For the former case, having access to the best rank-k approximation of A  $\mathcal{T}_k(A)$  is more costly than the rest of the procedure. For the latter case, computing such a value of  $\lambda$  requires knowledge (or an estimation) of the singular values of A  $\sigma_{k+1}, \ldots, \sigma_n$ , which may not be feasible in many practical situations. Even for a generic choice of  $\lambda$ , computing the ridge leverage scores for all the columns of A may be unfeasible in some applications in which the algorithm does not have full access to the entries of A. By this point of view, many strategies have been studied in the literature to approximate the ridge leverage scores associated to the columns of a matrix by recursive algorithms. For example, Cohen et al. [2017] discuss an iterative algorithm to compute overestimates of the true ridge leverage scores by uniformly sampling at random columns of A. We now conclude our survey on the ridge leverage scores by showing that under the aforementioned choice of  $\lambda$ , the sum of the corresponding ridge scores is not too large.

**Lemma 3.2.** Let  $\lambda^2 = \frac{\|A - \mathcal{T}_k(A)\|_F^2}{k}$  in the ridge scores (13) of the columns of A. Then  $\sum_{i=1}^n l_{i,\lambda}(A) \leq 2k$ .

*Proof.* We rewrite (13) using the SVD decomposition of  $A = U\Sigma V^T$ . This gives

$$l_{i,\lambda}(A) = a_i^T (U\bar{\Sigma}^{-2}U^T)a_i,$$

where  $\bar{\Sigma}_{ii}^{-2} = (\sigma_i^2(A) + \lambda^2)^{-1}$ . Then

$$\sum_{i=1}^{n} l_{i,\lambda}(A) = \text{Tr}(\Sigma^{2}\bar{\Sigma}^{-2}) = \sum_{i=1}^{n} \frac{\sigma_{i}^{2}(A)}{\sigma_{i}^{2}(A) + \frac{\|A - \mathcal{T}_{k}(A)\|_{F}^{2}}{L}}.$$

Now, for  $i \leq k$  we can bound the terms in the previous sum by 1. This gives eventually

$$\sum_{i=1}^{n} l_{i,\lambda}(A) \le k + \sum_{i=k+1}^{n} \frac{\sigma_i^2(A)}{\sigma_i^2(A) + \frac{\|A - \mathcal{T}_k(A)\|_F^2}{k}} \le k + \frac{1}{\frac{\|A - \mathcal{T}_k(A)\|_F^2}{k}} \sum_{i=k+1}^{n} \sigma_i^2(A) = 2k.$$

### 4 Numerical results

We conclude this work on the classical and ridge leverage scores associated to the columns of a matrix by presenting numerical evidence of their superior behaviour in sketching algorithms. We start this section by considering a Hilbert matrix of size n = 100. Recall that the Hilbert matrix is defined as follows:

$$A_{ij} = \frac{1}{i+j-1}, \quad i, j = 1, \dots n.$$
 (18)

It is a well-known fact that the Hilbert matrix have exponentially fast decaying of the singular values, which implies that it is not a full column rank matrix and the leverage scores are not all trivially equal. Moreover, these feature makes it suitable for numerical experiments on sketching algorithms. In the sequel, we compare the performance of the three main sampling probabilities which we discussed in the work: uniform sampling, columns norm sampling (3) and ridge leverage scores (13) with  $\lambda = 1e - 4$ . While such a value of  $\lambda$  was tuned by experimental evidence, the choice of the regularization parameter  $\lambda$  had a little influence on the performance.

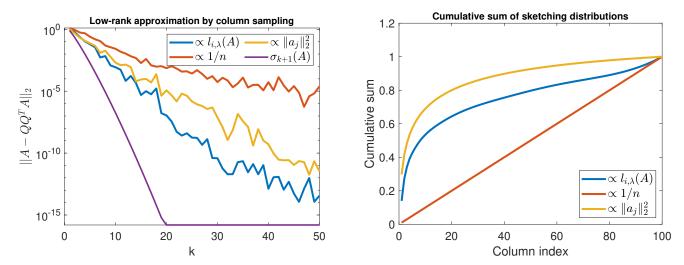


Figure 1: The figure on the left shows the performance of different column sampling distributions for the low-rank approximation of a Hilbert matrix of size n = 100. Approximations errors for all the sampling strategies (uniform sampling, columns norm sampling and ridge leverage score sampling with  $\lambda = 1e - 4$ ) are computed as  $||A - QQ^TA||_2$ , where Q is the orthogonal factor of the reduced QR decomposition of the sampled columns matrix C, and compared with the gold standard for a rank-k approximation given by  $\sigma_{k+1}(A)$ . As these strategies are randomized, results are averaged over 20 runs for all the considered ranks  $k = 1, \ldots, 50$ . The figure on the right shows the cumulative sum of the sketching distributions: by construction, the Frobenius norm of A is concentrated in the very first columns, while the ridge leverage scores assign higher probabilities to the last columns. Both distributions diverge clearly from the uniform column scores.

Figure 1 (left) shows the low-rank approximation error using the three different strategies for ranks k = 1, ..., 50, and their comparison to the gold standard given by  $\sigma_{k+1}(A)$ . As these strategies are randomized, the results are averaged over 20 runs for all considered ranks. The projection error in the spectral norm is measured as  $||A - QQ^T A||_2$ , where Q is computed from the reduced QR factorization of C and hence  $QQ^T$  is the orthogonal projector onto the column space of C. We see that ridge leverage score sampling (13) is outperforming both columns norm sampling (3) and uniform column sampling and the hierarchy in the quality of the approximation follows the outline of the work: uniform sampling performs poorly and remains far away from the gold standard, because all the columns of A are given the same probability. Columns norm sampling improves majorly the results, but sampling proportionally to the ridge leverage scores consistently yields a better column subset for the approximation of the range of A.

To understand the gap in the performance between columns norm sampling and ridge leverage scores, we show in Figure 1 (right) the cumulative sum of the scores obtained with the two strategies. First, we note that they both deviate quite clearly from the uniform distribution. This explains the poor performance of uniform sampling and suggests that the Hilbert matrix does not have low coherence, meaning that the importance of all the columns is not the same. On the other hand, coherently with the structure of (18), the cumulative sum of the column norms is saturated in the very first columns, and almost zero sampling probability is given to the last columns of A. On the other hand, the ridge leverage scores are assigning non-negligible sampling probabilities

to such columns. This difference in the sketching distributions gives an additional insight on the difference in performance which we observe in Figure 1 (left).

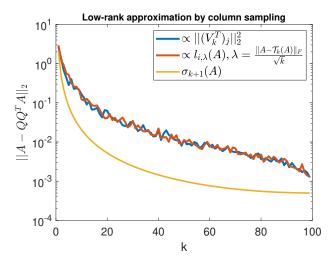


Figure 2: Comparison between column subset selection using as sampling probabilities the rank-k subspace leverage scores (12) and the ridge scores (13) with adaptive regularization parameter  $\lambda^2 = \frac{\|A - \mathcal{T}_k(A)\|_F^2}{k}$  on the matrix (19) of size n = 100. Approximations errors for the two sampling strategies are computed as  $\|A - QQ^TA\|_2$ , where Q is the orthogonal factor of the reduced QR decomposition of the sampled columns matrix C, and compared with the gold standard for a rank k approximation given by  $\sigma_{k+1}(A)$ . Results are averaged over 20 runs. We observe, coherently with what we expected, that the approximation error induced by the two strategies is on average the same.

Figure 2 presents evidence for the similarity between the truncated rank-k subspace leverage scores (12) and the ridge leverage scores (13) with adaptive regularization parameter  $\lambda^2 = \frac{\|A - \mathcal{T}_k(A)\|_F^2}{k}$ . For this experiment, to avoid numerical instability, we picked a matrix A of size 100 with entries defined by

$$A_{ij} = \exp(-|i - j|/1000), \tag{19}$$

which has a very slow singular value decay. As this matrix has full column rank, all its columns have leverage score equal to one. The figure shows that the approximation error induced by the two sampling strategies is on average the same for all ranks k = 1, ..., 100.

#### 5 Conclusions

In this work, we started from the main purpose of defining a powerful sampling probability for the random column subset selection task of a generic matrix  $A \in \mathbb{R}^{m \times n}$ . To this aim, we introduced in Section 2 the classical leverage scores (7) associated to each column of A as the minimum norm solution to a constrained least squares problem. Exploiting this interpretation, we showed their main properties and we argued that they represent a measure of the importance of each column in composing the column space of A. In particular, we proved that the leverage score is maximum for a column which is orthogonal to all the others, and we argued consequently that the leverage scores are superior with respect to the columns norms for column sampling because they possess coherence-revealing features. In Section 3, motivated by their provable poor performance for full column rank matrices (as they reduce to uniform scores for the columns of A), we followed the work presented in Cohen et al. [2015] and we introduced the ridge leverage scores (13) as a stable and natural procedure to filter out small principal components. We derived a simplified expression for the ridge leverage scores which gave insight on their superior behaviour for full column rank matrices. However, we clarified in the end our awareness of their highly expensive (or even simply unfeasible) computation. In this direction, we may follow as an improvement of the current work the recursive algorithm presented in Cohen et al. [2017], which computes overestimates of the ridge leverage scores by random column sampling of A at computation time. Finally, to conclude our work, we presented in Section 4 numerical evidence for the superior behaviour of ridge leverage score sampling with respect to columns norm sampling and uniform sampling in the low-rank approximation of a Hilbert matrix of size n = 100. Moreover, we presented a numerical example for the similarity between ridge leverage scores and scoring proportionally to the row norms of  $V_k$  on a full column rank matrix with slow singular value decay.

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