

# 8 Sparsification Algorithms

All low rank matrix approximation algorithms including the fundamental ones such as Power method or Orthogonal iterations, involve lots of matrix-matrix or matrix-vector multiplications. These basic operations require time proportional to number of non-zero entries in matrices, as one need to read the entire matrix into memory. Sparsifying a matrix, i.e. decreasing number of non-zeros, and quantizing it, i.e. rounding up entries to a constant, accelerate such computations as well as saving space in representation.

First sparsification algorithm was by Achlioptas and McSherry[1], where they sampled and quantized entries of a given matrix  $A \in \mathbb{R}^{n \times d}$  to lowered number of non-zeros and length of their representation. They observed acts of sampling and quantization can be viewed as adding a random noise matrix  $E \in \mathbb{R}^{n \times d}$  to  $A$ , whose entries are independent random variables with zero mean and bounded variance. Since with high probability a random matrix has a weak spectral structure, it does not alter the the main spectrum of input matrix. Below we first state a theorem on norm of random matrices, then describe their algorithms.

## 8.1 Spectral Structure of Random Matrices

Theorem below[2] shows a well constructed random matrix has a weak spectral structure.

**Theorem 8.1.1.** [2] Let  $E \in \mathbb{R}^{n \times d}$  be a random matrix such that entries  $E_{i,j} = r_{ij}$  are independent bounded random variables  $r_{ij} \in [-k, k]$ , with  $\mathbf{E}[r_{ij}] = 0$  and  $\mathbf{Var}(r_{ij}) \leq \sigma^2$ . For all  $\alpha \geq 1, \varepsilon > 0$ , and  $n + d \geq 20$ , if  $k \leq \left(\frac{4\varepsilon}{4+3\varepsilon}\right)^3 \frac{\sigma\sqrt{n+d}}{\log^3(n+d)}$  then

$$\Pr \left[ \|E\|_2 \geq (2 + \varepsilon + \alpha)\sigma\sqrt{n+d} \right] < (n+d)^{-\alpha^2}$$

## 8.2 Additive Error Sparsification Algorithms

Using theorem 8.1.1, Achlioptas and McSherry[1] showed a carefully constructed random matrix  $\hat{A} \in \mathbb{R}^{n \times d}$  can approximate spectral norm of  $A_k$ . Theorem 8.2.1 states their result.

**Theorem 8.2.1.** Let  $A \in \mathbb{R}^{n \times d}$  be an arbitrary matrix with  $b = \max_{i,j} |A_{i,j}|$  being the maximum entry in absolute value. Let  $\hat{A} \in \mathbb{R}^{n \times d}$  be a random matrix where entries  $\hat{A}_{i,j}$  are independent random variables with  $\mathbf{E}[\hat{A}_{i,j}] = A_{i,j}$ ,  $\mathbf{Var}(\hat{A}_{i,j}) = (\sigma b)^2$  and  $\|A_{i,j} - \hat{A}_{i,j}\|_2 \leq \frac{\sigma b\sqrt{n+d}}{2\log^3(n+d)}$ . Then for any  $\alpha \geq 1$ ,

$$\|A - \hat{A}_k\|_2 \leq \|A - A_k\|_2 + (8 + 2\alpha)\sigma b\sqrt{n+d}$$

holds with probability atleast  $1 - (n+d)^{-\alpha^2}$ .

*Proof.*

$$\begin{aligned} \|A - \hat{A}_k\|_2 &\leq \|A - \hat{A}\|_2 + \|\hat{A} - \hat{A}_k\|_2 && \text{triangle inequality} \\ &\leq \|A - \hat{A}\|_2 + \|\hat{A} - A_k\|_2 && \text{For any rank } k \text{ matrix } D: \|\hat{A} - \hat{A}_k\|_2 \leq \|\hat{A} - D\|_2 \\ &\leq \|A - \hat{A}\|_2 + \|\hat{A} - A\|_2 + \|A - A_k\|_2 && \text{triangle inequality} \\ &\leq 2\|A - \hat{A}\|_2 + \|A - A_k\|_2 \end{aligned}$$

Setting  $E = A - \hat{A}$  one can verify that  $E$  satisfies all conditions of theorem 8.1.1, as it has zero expectation  $\mathbf{E}[E_{i,j}] = A_{i,j} - \mathbf{E}[\hat{A}_{i,j}] = 0$ , bounded variance  $\mathbf{Var}(E_{i,j}) \leq (\sigma b)^2$ , and bounded entries  $E_{i,j} \in \left[-\frac{\sigma b\sqrt{n+d}}{2\log^3(n+d)}, \frac{\sigma b\sqrt{n+d}}{2\log^3(n+d)}\right]$ . Therefore taking  $\varepsilon = 2$ , the bound  $\|A - \hat{A}\|_2 \leq (4 + \alpha)\sigma b\sqrt{n+d}$  holds with probability atleast  $1 - (n+d)^{-\alpha^2}$ , and therefore  $\|A - \hat{A}_k\|_2 \leq (8 + 2\alpha)\sigma b\sqrt{n+d} + \|A - A_k\|_2$ .  $\square$

As theorem 8.2.1 holds for any random matrix  $\hat{A}$  with above conditions, authors of [1] proposed two concrete constructions. The first construction is based on sampling; matrix  $\hat{A}$  samples some entries of  $A$  and omits others, they show the stronger spectrum of input matrix is, the larger fraction of entries they can afford to lose. Theorem 8.2.2 states their sampling result.

**Theorem 8.2.2.** Let  $A \in \mathbb{R}^{n \times d}$  be the input matrix and  $b = \max_{i,j} |A_{i,j}|$  be the maximum entry in absolute value. Define matrix  $\hat{A} \in \mathbb{R}^{n \times d}$  as

$$\hat{A}_{i,j} = \begin{cases} 0 & \text{w.p. } 1 - \frac{1}{s} \\ sA_{i,j} & \text{w.p. } \frac{1}{s} \end{cases}$$

where  $1 \leq s \leq \frac{n+d}{4 \log^6(n+d)}$ . Then with probability atleast  $1 - 1/(n+d)$  the following error bound holds

$$\|A - \hat{A}\|_2 \leq \|A - A_k\|_2 + 10b\sqrt{s(n+d)}$$

*Proof.* It is easy to verify that matrix  $\hat{A}$  satisfies all conditions of theorem 8.2.1:

- $\mathbf{E}[\hat{A}_{i,j}] = 0(1 - 1/s) + sA_{i,j}(1/s) = A_{i,j}$
- $\mathbf{Var}(\hat{A}_{i,j}) = \mathbf{E}[\hat{A}_{i,j}^2] - \mathbf{E}[\hat{A}_{i,j}]^2 = 1/s(sA_{i,j})^2 - A_{i,j}^2 = (s-1)A_{i,j}^2 \leq (\sqrt{s}b)^2$  therefore  $\sigma = \sqrt{s} \leq \frac{\sqrt{n+d}}{2 \log^3(n+d)}$
- $\forall i \in [1, n], j \in [1, d] : |A_{i,j} - \hat{A}_{i,j}| \in \{A_{i,j}, sA_{i,j}\}$ , and in both cases it is upper bounded by  $|A_{i,j} - \hat{A}_{i,j}| \leq sA_{i,j} \leq \frac{(n+d)b}{4 \log^6(n+d)}$

Fitting conditions of theorem 8.2.1, and using  $\alpha = 1$ , we obtain  $\|A - \hat{A}_k\|_2 \leq 10b\sqrt{s(n+d)} + \|A - A_k\|_2$ .  $\square$

In their second construction, they randomly quantize entries of  $A$ , and shorten the representation, this allows them to store each entry in one bit. Theorem 8.2.3 explains their result.

**Theorem 8.2.3.** Let  $A \in \mathbb{R}^{n \times d}$  be the input matrix and  $b = \max_{i,j} |A_{i,j}|$  be the maximum entry in absolute value. Define matrix  $\hat{A} \in \mathbb{R}^{n \times d}$  as

$$\hat{A}_{i,j} = \begin{cases} +b & \text{w.p. } \frac{1}{2} + \frac{A_{i,j}}{2b} \\ -b & \text{w.p. } \frac{1}{2} - \frac{A_{i,j}}{2b} \end{cases}$$

Then  $\|A - \hat{A}\|_2 \leq \|A - A_k\|_2 + 10b\sqrt{(n+d)}$  with probability atleast  $1 - 1/(n+d)$ .

*Proof.* Again it's easy to see that matrix  $\hat{A}$  satisfies all conditions of theorem 8.2.1:

- $\mathbf{E}[\hat{A}_{i,j}] = b(\frac{1}{2} + \frac{A_{i,j}}{2b}) - b(\frac{1}{2} - \frac{A_{i,j}}{2b}) = A_{i,j}$
- $\mathbf{Var}(\hat{A}_{i,j}) = \mathbf{E}[\hat{A}_{i,j}^2] - \mathbf{E}[\hat{A}_{i,j}]^2 = b^2(\frac{1}{2} + \frac{A_{i,j}}{2b}) + b^2(\frac{1}{2} - \frac{A_{i,j}}{2b}) - A_{i,j}^2 \leq b^2$  therefore  $\sigma = 1$
- $\forall i \in [1, n], j \in [1, d] : |A_{i,j} - \hat{A}_{i,j}| = |A_{i,j} \pm b| \leq 2b$

Fitting conditions of theorem 8.2.1, and using  $\alpha = 1$  completes the proof

$$\|A - \hat{A}_k\|_2 \leq 10b\sqrt{(n+d)} + \|A - A_k\|_2$$

$\square$

### 8.3 Relative Error Sparsification Algorithm

Latest result in using sparsification for low-rank approximation [3] takes advantage of a popular technique in matrix completion line of work, called as *alternating minimization*. We first give a brief review of this technique, then elaborate the main algorithm.

Often a target matrix  $A$  can be represented in a bi-linear form as  $A = UV$  (matrices  $U, V$  are not necessarily orthonormal). Having this parametrization, the task of approximating  $A$  reduces to finding  $U$  and  $V$  that minimize an error metric, for example  $\|A - UV\|_F$ . The *alternating minimization* technique starts with some initial guess for  $U$  and  $V$  (say  $U^{(0)}, V^{(0)}$ ), iteratively keep one of  $U, V$  fixed and optimize over the other, that is  $V^{(i+1)} = \arg \min_V \|A - U^{(i)}V\|_F$ , then switch and repeat until it converges.

In order to use this technique in matrix approximation, algorithm[3] samples some entries of matrix  $A$ , partition them into multiple subsets and iterates over those subsets to refine the approximation it obtained from first subset. The full method is described in algorithm 8.3.1 and 8.3.2.

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#### Algorithm 8.3.1 Leverage Element Low Rank Approximation (LELA)

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- 1: **Input:**  $A \in \mathbb{R}^{d \times n}$ , rank  $r$ , number of samples  $m$ , number of iterations  $T$
  - 2: **Output:**  $P_\Omega(A), \Omega, r, \hat{q}, T$
  - 3:  $\Omega \subset [n] \times [d] \leftarrow$  indices of  $m$  independently sampled entries with probability  $\hat{q}_{i,j} = \min\{1, q_{i,j}\}$  with  $q_{i,j} = m \cdot \left( \frac{\|A_{i,:}\|^2 + \|A_{:,j}\|^2}{2(n+d)\|A\|_F^2} + \frac{|A_{i,j}|}{2\|A\|_1} \right)$
  - 4: obtain  $P_\Omega(A) \subset A$  as the matrix of sampled entries, using another pass over  $A$
  - 5:  $\hat{A}_r = W \text{AltMin}(P_\Omega(A), \Omega, r, \hat{q}, T)$
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#### Algorithm 8.3.2 Weighted Alternative Minimization

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- 1: **Input:**  $P_\Omega(A), \Omega, r, \hat{q}, T$
  - 2: **Output:**  $\hat{A}_r \in \mathbb{R}^{n \times d}$
  - 3: For all  $i, j \in [n] \times [d]$  set  $w_{i,j} = 1/\hat{q}_{i,j}$  if  $\hat{q}_{i,j} > 0$ , otherwise  $w_{i,j} = 0$
  - 4: Divide  $\Omega$  into  $2T + 1$  equal uniformly random subsets  $\Omega = \{\Omega_0, \dots, \Omega_{2T}\}$
  - 5:  $R_{\Omega_0}(A) \leftarrow w \cdot * P_{\Omega_0}(A)$
  - 6: Set  $U^{(0)}\Sigma^{(0)}(V^{(0)})^T = \text{svd}(R_{\Omega_0}(A), r)$
  - 7: **for**  $t = 0$  to  $T - 1$  **do**
  - 8:  $\hat{V}^{(t+1)} = \arg \min_{V \in \mathbb{R}^{d \times r}} \|R_{\Omega_{2t+1}}^{1/2} (A - \hat{U}^{(t)}V^T)\|_F^2$
  - 9:  $\hat{U}^{(t+1)} = \arg \min_{U \in \mathbb{R}^{n \times r}} \|R_{\Omega_{2t+2}}^{1/2} (A - U(\hat{V}^{(t+1)})^T)\|_F^2$
  - 10: **return**  $\hat{A}_r = \hat{U}^{(T)}(\hat{V}^{(T)})^T$
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In the sampling phase, whose aim is to sparsify the matrix, each entry  $A_{i,j}$  is sampled with a defined probability  $q_{i,j}$  and weighted as  $A_{i,j}/q_{i,j}$ , so that sampled matrix  $\hat{A} \in \mathbb{R}^{n \times d}$  has same frobenious norm as  $A$  in expectation. As decomposing a matrix takes time inversly proportional to the sparsity of the matrix authors spread non-zero entries of  $\hat{A}$  equally and randomly amongst some fixed numbers of matrices  $\hat{A}^{(j)} \in \mathbb{R}^{n \times d}$ , therefore  $\sum_{j=1} \hat{A}^{(j)} = \hat{A}$ . Now that each  $\hat{A}^{(j)}$  is a sparse random sample of  $\hat{A}$ , they take **svd** decomposition of  $\hat{A}^{(1)}$  explicitly, i.e  $[U, S, V] = \hat{A}^{(1)}$ . Considering  $\hat{A}^{(1)}$  in bi-linear form  $\hat{A}^{(1)} = U(SV^T)$ , they iterate over further matrices  $\{A^{(j)}\}$  and minimize the fronbenious error of approximation.

They show that 8.3.1 needs  $T = O(\log(\frac{\|A\|_2}{\epsilon \|A - A_r\|_F}))$  iterations, runs in time  $O(\text{nnz}(A) + \frac{nr^5}{\epsilon^2} \kappa^2 \log n)$  where  $\kappa = \sigma_1/\sigma_r$  is the condition number of  $A$ , and achieves the relative error bound

$$\|A - \hat{A}_r\|_2 \leq \|A - A_r\|_2 + 2\epsilon \|A - A_r\|_F$$



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# Bibliography

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