

**1** Let  $\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_n$  be independent Rademacher random variables (that is,  $\mathbf{x}_i = \pm 1$  with equal probability) and  $\mathbf{a} = (a_1, a_2, \dots, a_n)$  be a sequence of real numbers.

- (a) Show that  $\mathbb{E}e^{s\mathbf{x}_i} = \cosh(s)$ .
- (b) Prove that  $\cosh(s) \leq e^{s^2/2}$ .
- (c) Use (a), (b), and Markov's inequality to prove that

$$\mathbb{P}\left(\left|\sum_{i=1}^n a_i \mathbf{x}_i\right| \geq \epsilon\right) \leq e^{-\frac{\epsilon^2}{2\|\mathbf{a}\|_2^2}}$$

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**2<sup>a</sup>** In the  $k$ -means clustering problem we are given some input vectors  $\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_m \in \mathbb{R}^n$  and a positive integer  $k$ , and we'd like to output a partition  $P$  of  $\{1, 2, \dots, n\}$  into  $k$  disjoint subsets  $P_1, P_2, \dots, P_k$  and cluster centers  $\mathbf{y}_1, \mathbf{y}_2, \dots, \mathbf{y}_k \in \mathbb{R}^n$ . which minimize the function

$$f(\{P_i\}, \{\mathbf{y}_i\}; \{\mathbf{x}_i\}) = \sum_{j=1}^k \sum_{i \in P_j} \|\mathbf{x}_i - \mathbf{y}_j\|_2^2.$$

That is, the  $x_i$  are clustered into  $k$  clusters according to  $P$ . This problem is NP-hard, but good approximation algorithms exist which can return almost-optimal clusterings.

- (a) For a fixed partition  $P$ , show that the optimal  $\mathbf{y}_1, \mathbf{y}_2, \dots, \mathbf{y}_k$  is where for every nonempty  $P_i \in P$ ,

$$\mathbf{y}_i = \frac{1}{|P_i|} \sum_{j \in P_i} \mathbf{x}_j$$

is just the average of the points in  $P_i$ . Thus we can restrict ourselves to optimizing over partitions  $P$ .

- (b) Prove that for a given cluster  $P_i$ , the optimal cost is

$$\frac{1}{2|P_i|} \sum_{j, k \in P_i} \|\mathbf{x}_j - \mathbf{x}_k\|_2^2$$

- (c) Using the Johnson-Lindenstrauss lemma, show that for any  $0 < \epsilon < \frac{1}{2}$  there is a linear map  $\mathbf{S} \in \mathbb{R}^{m \times n}$ ,  $m = \mathcal{O}(\epsilon^{-2} \log m)$  such that for all partitions  $P$  simultaneously

$$(1 - \epsilon)f(\{P_i\}; \{\mathbf{x}_i\}) \leq f(\{P_i\}; \{\mathbf{S}\mathbf{x}_i\}) \leq (1 + \epsilon)f(\{P_i\}; \{\mathbf{x}_i\})$$

and where  $\mathbf{S}$  can be found efficiently with a randomized algorithm with small failure probability. Thus if one does not mind worsening the quality of our clusters by a factor  $1 + \epsilon$ , without loss of generality one can assume that the input vectors  $\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_m \in \mathbb{R}^n$  are in dimension  $n = \mathcal{O}(\epsilon^{-2} \log m)$ , which can be *much* smaller than the original dimension  $n$ .

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<sup>a</sup>Jelani Nelson, 2013

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