#### Stochastic Optimization: Complexity-Based Analysis and Development Engineering Applications

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#### <span id="page-5-0"></span>Sample Average Approximation (SAA)

Sample Average Approximation (SAA) is a commonly-used procedure for approximating solutions to stochastic optimization problems of the form

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The idea of SAA is to first generate an i.i.d. sample  $\xi_1, \ldots, \xi_n$  of the random variable *ξ*, and then approximate the expectation E*ξf*(*x, ξ*) using its sample average

$$
\min_{x \in \mathcal{X}} \{ F_n(x) := \frac{1}{n} \sum_{i=1}^n f(x, \xi_i) \}.
$$
 (2)



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Towards this goal, a now classical analysis [\[Kleywegt et al., 2002,](#page-53-0) [Shapiro, 2003,](#page-54-0) [Shapiro et al., 2009](#page-54-1)] showed that in order to ensure

$$
\mathbb{P}\big(\mathbf{F}(\hat{\mathbf{x}}_n) - \mathbf{F}(\mathbf{x}^*) \le \delta\big) \ge 1 - \alpha \tag{3}
$$

for any  $\delta \in (0,1]$  and  $\alpha \in (0,1]$ , the number of samples *n* should satisfy

$$
n \gtrsim \frac{\rho}{\delta^2} \log \frac{1}{\delta} + \frac{1}{\delta^2} \log \frac{1}{\alpha}.
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These bounds depend polynomially on problem dimension *p*.





#### <span id="page-11-0"></span>Rademacher Complexity

Analysis depends on stochastic process theory.



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Let  $\epsilon_1,\ldots,\epsilon_n$  be i.i.d. Rademacher random variables, where  $\mathbb{P}(\epsilon=\pm 1)=\frac{1}{2}$ ; and let *f*(*x, ξ*) be the function from the objective of the SAA problem. We define the *Rademacher complexity* of the function set  $\mathcal{F} := \{f(x, \xi) : x \in \mathcal{X}\}\)$  to be

$$
\mathcal{R}_n[f] = \mathbb{E}_{\xi} \Big( \sup_{x \in \mathcal{X}} \Big| \frac{1}{n} \sum_{i=1}^n \epsilon_i f(x, \xi_i) \Big| \Big). \tag{5}
$$



#### Rademacher Complexity

With an assumption that  $-\Delta/2 \le f(x,\xi) \le \Delta/2$  for all  $(x,\xi) \in \mathcal{X} \times \Xi$ , for some finite constant  $\Delta \in \mathbb{R}_+$ , we give a concentration bound of the form: Proposition 1

$$
\mathbb{P}\Big(\sup_{x\in\mathcal{X}}\Big|F_n(x)-F(x)\Big|>t\Big)\leq \exp\Big(-2n\Big(\frac{t-2\mathcal{R}_n[f]}{\Delta}\Big)^2\Big).
$$
 (6)

The proof involves use of Jensen and McDiarmid's inequalities, a symmetrization argument, and an application of the triangle inequality.



#### <span id="page-14-0"></span>Improved Sample Bounds

#### Proposition 2

*Let*  $g : \mathbb{R} \to \mathbb{R}$  *be Lipschitz with constant L, and consider the stochastic optimization problem*

$$
\min_{x \in \mathcal{S}} \left\{ \mathbb{E}_{\xi} \left( g(\xi^{\mathsf{T}} x) \right) \middle| ||x||_1 \leq \lambda \right\}
$$
\n(7)

*where S ⊆* R *p and* max*<sup>ξ</sup>∈*<sup>Ξ</sup> *∥ξ∥<sup>∞</sup> ≤ <sup>C</sup> <* +*∞. Then the Rademacher complexity of the*  $a$ bove problem is bounded by  $\mathcal{R}_n[f] \leq \lambda$ LC $\sqrt{2\log 2p/n}$ , and we need

$$
n \ge \left(\frac{3\lambda\mathcal{L}\mathcal{C}}{\delta}\right)^2 \cdot \left(2\log\left(\frac{2}{\alpha}\right) + 2\log 2\rho\right) \tag{8}
$$

*samples to ensure that*

$$
\mathbb{P}\big(F(\hat{x}_n) - F(x^*) \le \delta\big) \ge 1 - \alpha \tag{9}
$$

*for any*  $\delta \in (0,1]$  *and*  $\alpha \in (0,1]$ *, holds.* 



#### Improved Sample Bounds

Proposition 3

*Let*  $g : \mathbb{R} \to \mathbb{R}$  *be Lipschitz with constant L, and consider the stochastic optimization problem*

$$
\min_{X \in \mathcal{S}} \left\{ \mathbb{E}_{\xi} \big( g(tr(\xi^T X)) \big) \middle| ||X||_* \leq \lambda \right\}
$$

 $\mathcal{S} \subseteq \mathbb{R}^{p \times q}$  and  $\max_{\xi \in \Xi} \|\xi\|_2 \leq \zeta + \infty$ . Then the Rademacher complexity of *the above stochastic optimization problem is bounded by*  $\mathcal{R}_n[f] \leq \lambda$ LC $\sqrt{3\log(\min\{\rho, q\})/n}$ , and we need

$$
n \geq \left(\frac{3\lambda\iota\zeta}{\delta}\right)^2 \cdot \left(2\log\left(\frac{2}{\alpha}\right) + 3\log\left(\min\{p,q\}\right)\right)
$$

samples to ensure that for any  $\delta \in (0,1]$  and  $\alpha \in (0,1]$ , (17) holds.



<span id="page-16-0"></span>Consider a scenario where we would like to choose a portfolio that allocates investments into some combination of *p* risky assets and 1 risk-free asset, while considering a tradeoff between maximizing the expected return of the portfolio and the risk tolerance of the investor.





Consider a scenario where we would like to choose a portfolio that allocates investments into some combination of *p* risky assets and 1 risk-free asset, while considering a tradeoff between maximizing the expected return of the portfolio and the risk tolerance of the investor.



The Markowitz portfolio selection model [\[Markowitz, 1952,](#page-53-1) [Bruder et al., 2013](#page-52-0)] is a simple framework to pose such a problem.



Let  $\xi \in \mathbb{R}^p$  is a random variable of the returns from the  $p$  risky assets, and define  $\mu = \mathbb{E}_{\xi}\xi$  and  $\Sigma = \mathbb{E}_{\xi}((\xi - \mu)(\xi - \mu)^{\intercal}).$  Then one formulation of the problem involves solving a convex quadratic program

$$
\min_{x \in \mathbb{R}^p} \left\{ x^T \Sigma x - \gamma \cdot x^T (\mu - r\mathbf{1}) \middle| x \ge 0, ||x||_1 \le 1 \right\}
$$
\n(10)

where:

*r* is the rate of return for the risk-free asset,



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Each entry of the vector *x* gives the fraction of the portfolio allocated to the *p* risky assets; hence  $1 - \sum_{i=1}^p x_i$  is the fraction of the portfolio allocated to the risk-free asset.





Figure 1: Comparison of 95% upper confidence bound of SAA solution gap (solid blue) with bounds on 95% upper confidence bound gap predicted classically (dash-dotted red), our Proposition (dashed orange), and our Corollary (dotted green).

In both plots, the x-axis is the dimension *p* of the decision variable, and the y-axis is the 95% upper confidence bound gap.



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• Tensors generalize matrices.



#### Examples

Vectors are 1D tensors, matrices 2D, and so on...



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- **Tensor Rank and Decomposition**: Though related, many problems that are polynomial-time solvable for matrices are NP-hard for tensors.



- Tensors generalize matrices.
- **Tensor Rank and Decomposition**: Though related, many problems that are polynomial-time solvable for matrices are NP-hard for tensors.
- It is NP-hard to compute the rank of a tensor[[Hillar and Lim, 2013](#page-53-2)], & tensor versions of the spectral norm, nuclear norm, and matrix singular value decomposition are also NP-hard to compute. [[Hillar and Lim, 2013](#page-53-2), [Friedland and Lim, 2014\]](#page-52-1).



<span id="page-27-0"></span>Tensor completion is the problem of observing (possibly with noise) a subset of entries of a tensor and then estimating the remaining entries based on an assumption of low-rankness.



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Suppose we have data  $(x\langle i \rangle, y\langle i \rangle) \in \mathcal{R} \times \mathbb{R}$  for  $i = 1, \ldots, n$ . Let  $I = \{i_1, \ldots, i_u\} \subseteq [n]$  be any set of points that specify all the unique  $x \langle i \rangle$  for  $i = 1, \ldots, n$ .



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The nonnegative tensor completion problem is given by

$$
\widehat{\psi} \in \arg\min_{\psi} \frac{1}{n} \sum_{i=1}^{n} (y \langle i \rangle - \psi_{x \langle i \rangle})^{2}
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\n
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\text{s.t. } \text{rank}(\psi) \le \lambda
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The "state of the art" computational methods use decomposition and an alternating minimization procedure to solve.



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Completion of nonnegative rank-1 tensors can be written as a convex optimization problem [\[Aswani, 2016\]](#page-52-2).

For symmetric orthogonal tensors, a variant of the Frank-Wolfe algorithm has been proposed[[Rao et al., 2015](#page-54-2)], which can be shown to achieve the information-theoretic rate.



To date, no tensor completion algorithm has been shown to achieve the information-theoretic sample complexity rate, while guaranteeing convergence.



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Namely, for a tensor completion problem on a rank *k* tensor with sample size *n*, the information theoretic rate for estimation error is

$$
\sqrt{k\cdot\sum_i r_i/n}
$$

[\[Gandy et al., 2011\]](#page-52-3).



#### <span id="page-36-0"></span>A New Norm for Nonnegative Tensors

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It also depends on concepts and sets from the tensor world we cannot introduce here.



## A New Norm for Nonnegative Tensors

Proposition 4

*The function defined as*

$$
\|\psi\|_{+} := \inf\{\lambda \ge 0 \mid \psi \in \lambda C_1\} \tag{12}
$$

*is a norm for nonnegative tensors*  $\psi \in \mathbb{R}_+^{r_1 \times \cdots \times r_p}.$ 

We will call the set  $C_{\lambda}$  the nonnegative tensor polytope. A useful observation is that the following relationships hold:  $B_{\lambda} = \lambda B_1$ ,  $S_{\lambda} = \lambda S_1$ , and  $C_{\lambda} = \lambda C_1$ .



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In our case  $C_{\lambda}$  is not symmetric about the origin, and so without proof we do not *a priori* know whether scaling  $C_1$  eventually includes the entire space of nonnegative tensors. Thus we have to explicitly prove the gauge is a norm.



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Since the set of nonnegative tensors forms a cone [\[Qi et al., 2014\]](#page-53-3), we must prove our norm using a modified definition of a norm (Proof omitted).



#### <span id="page-41-0"></span>Results: Order 3 Tensors (r = dimensions)



Figure 2: Results for order-3 nonnegative tensors with size  $r \times r \times r$  and  $n = 500$  samples.



#### Results: Increasing Tensor Order (p)



Figure 3: Results for increasing order nonnegative tensors with size  $10^{\times p}$  and  $n=10,000$ samples.



#### Results:  $10^6$  entries and Increasing Sample **Size**



Figure 4: Results for nonnegative tensors with size 10*×*<sup>6</sup> and increasing *n* samples.



#### Results:  $10^7$  entries and Increasing Sample **Size**



Figure 5: Results for nonnegative tensors with size 10*×*<sup>7</sup> and increasing *n* samples.



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#### UN SDGs





### <span id="page-48-0"></span>Optimal Intervention Theory (OIT)

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Optimal Intervention Theory (OIT) is a method for improving human systems based on Statistical Learning Theory (SLT), which is the basis for learning in machines.

**Seeking a Productive Balance** 

# **Exploitation Persistence Exploration**



#### A New Solution Framework: Optimal Intervention Theory (OIT)

- OIT considers the classic trade-off between statistically rigorous methods and large-scale methods
- The Dynamic Programming algorithm, which solves problems with well-defined end-goals in stages, is the basis of the SLT applicable to OIT.





<span id="page-51-0"></span>

The full papers are available on my website at www.calebxb.com





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