Nyström Kernel Stein Discrepancy*

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Quick Summary

- Kernel Stein discrepancy (KSD; [1, 2]): powerful goodness-of-fit measure and test.
- Applications: Assessing and improving sample quality, validating MCMC methods, comparing deep generative models,
- Limitations: Quadratic runtime complexity.
- Main contribution: Accelerated estimator with the same convergence rate as the quadratic-time estimator.

Kernel Stein Discrepancy

Goal: Test H₀ : P = Q vs. H₁ : P ≠ Q for fixed known target P and unknown sampling distribution Q, given samples Q̂_n := {**x**_i}ⁿ_{i=1} ⊂ R^d of Q.
We illustrate the method with the Langevin-Stein operator-based [3] KSD on R^d, which is

$$S_{p}(\mathbb{Q}) = \left\| \mathbb{E}_{X \sim \mathbb{Q}} h_{p}(\cdot, X) \right\|_{\mathcal{H}_{h_{p}}},$$
with kernel $(\mathbf{x}, \mathbf{y} \in \mathbb{R}^{d})$

$$h_{p}(\mathbf{x}, \mathbf{y}) := \left\langle \nabla_{\mathbf{x}} \log p(\mathbf{x}), \nabla_{\mathbf{y}} \log p(\mathbf{y}) \right\rangle_{\mathbb{R}^{d}} k(\mathbf{x}, \mathbf{y})$$

$$+ \left\langle \nabla_{\mathbf{y}} \log p(\mathbf{y}), \nabla_{\mathbf{x}} k(\mathbf{x}, \mathbf{y}) \right\rangle_{\mathbb{R}^{d}}$$

$$+ \left\langle \nabla_{\mathbf{x}} \log p(\mathbf{x}), \nabla_{\mathbf{y}} k(\mathbf{x}, \mathbf{y}) \right\rangle_{\mathbb{R}^{d}} + \sum_{i=1}^{d} \frac{\partial^{2} k(\mathbf{x}, \mathbf{y})}{\partial x_{i} \partial y_{i}}$$

Main Results

• \sqrt{n} -consistency of KSD estimator: If

$$\left\| \left\| h_p(\cdot, X) \right\|_{\mathcal{H}_{hp}} \right\|_{\psi_2} < \infty$$

holds (implied by (1)), then

$$\left|S_p(\mathbb{Q}) - S_p(\hat{\mathbb{Q}}_n)\right| = \mathcal{O}_P(n^{-1/2}).$$

• \sqrt{n} -consistency of N-KSD: If the sub-Gaussian property (1) holds, then $\left|S_p(\mathbb{Q}) - \tilde{S}_p(\hat{\mathbb{Q}}_n)\right| = O_P(n^{-1/2}),$

given that the effective dimension $\mathcal{N}_{\mathbb{Q},\bar{h}_p}(\lambda) := \operatorname{tr}\left(C_{\mathbb{Q},\bar{h}_p,\lambda}^{-1}C_{\mathbb{Q},\bar{h}_p}\right)(C_{\mathbb{Q},\bar{h}_p}:=\mathbb{E}_{X\sim\mathbb{Q}}\left[h_p\left(\cdot,X\right)\otimes h_p\left(\cdot,X\right)\right]; C_{\mathbb{Q},\bar{h}_p,\lambda} := C_{\mathbb{Q},\bar{h}_p} + \lambda I, \lambda > 0)$ either • decays polynomially:

$$\mathcal{N}_{\mathbb{Q},\bar{h}_p}(\lambda) \lesssim \lambda^{-\gamma}, \qquad m = \tilde{\Omega}\left(n^{1/(2-\gamma)}\right),$$

for $\gamma \in (0, 1]$ (computational savings if $\gamma < 1/2$), or

p the (Lebesgue) density of \mathbb{P} , and kernel $k : \mathbb{R}^d \times \mathbb{R}^d \to \mathbb{R}$.

- Note that knowledge of the derivative of the score function $\nabla_{\mathbf{x}} \log p(\mathbf{x})$ is enough, i.e., knowledge of p up to normalization suffices.
- Existing KSD estimators take (roughly) the form

• Runtime: $\mathcal{O}(mn+m^3)$, saving if $m = o(n^{2/3})$.

$$S_p^2(\hat{\mathbb{Q}}_n) = \frac{1}{n^2} \sum_{i,j=1}^n h_p(\mathbf{x}_i, \mathbf{x}_j),$$

and have a runtime requirement of $\mathcal{O}(n^2)$.

Nyström-based Estimator (N-KSD)

• Denote by $\tilde{\mathbb{Q}}_m := \{\{\tilde{\mathbf{x}}_i\}\}_{i=1}^m$ a subsample of $\hat{\mathbb{Q}}_n$. The Nyström estimator is $\tilde{S}_p^2(\hat{\mathbb{Q}}_n) = \boldsymbol{\beta}_p^\mathsf{T} \mathbf{K}_{h_p,m,m}^- \boldsymbol{\beta}_p$, with $\boldsymbol{\beta}_p = \frac{1}{n} \mathbf{K}_{h_p,m,n} \mathbf{1}_n \in \mathbb{R}^m$, matrices $\mathbf{K}_{h_p,m,n} = [h_p(\tilde{\mathbf{x}}_i, \tilde{\mathbf{x}}_j)]_{i,j=1}^m \in \mathbb{R}^{m \times m}$, and $\mathbf{K}_{h_p,m,n} = [h_p(\tilde{\mathbf{x}}_i, \mathbf{x}_j)]_{i,j=1}^{m,n} \in \mathbb{R}^{m \times n}$, and \mathbf{A}^- denoting the (Moore-Penrose) pseudo-inverse of a matrix \mathbf{A} . • decays exponentially:

$$\mathcal{N}_{\mathbb{Q},\bar{h}_p}(\lambda) \lesssim \log(1+c_1/\lambda), \qquad m = \tilde{\Omega}\left(n^{1/2}\right),$$

for some $c_1 > 0$ (computational savings if n is large enough).

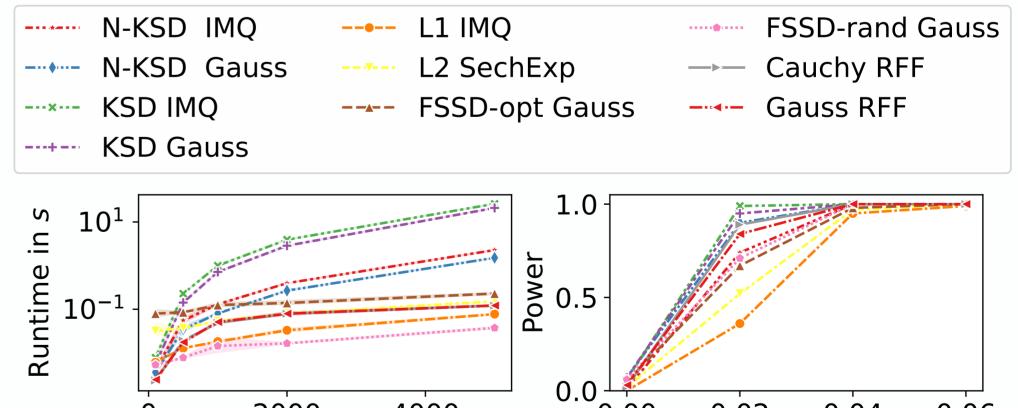
• The decay of the effective dimension can be linked to the decay of the eigenvalues of the covariance operator $C_{\mathbb{Q},\bar{h}_p}$ [4, Proposition 4, 5].

Discussion

- Unboundedness of the feature map handled by sub-Gaussian assumption.
- The quadratic-time and the N-KSD estimator both have \sqrt{n} -consistency, i.e., computational gain with no loss in statistical accuracy.
- Our results apply in the general KSD framework [5].
- Open: Weaker assumption for the Nyström case.

Goodness-of-fit Benchmark

• Runtime and power of Nyström KSD (N-KSD) and competitors on the restricted Boltzmann machine (RBM) goodness-of-fit benchmark.



Sub-Gaussian Assumption

 \bullet Existing Nyström analysis considers bounded kernels only. In practice, h_p is usually unbounded and existing results do not apply.

• Example: Consider d = 1, standard normal $p(x) \propto \exp(-x^2/2)$, and the RBF kernel $k(x, y) = \exp(-\gamma(x - y)^2)$ ($\gamma > 0$). Then

$$h_p(x,x) = x^2 + 2\gamma \stackrel{x \to \infty}{\to} \infty.$$

Similarly, for the IMQ kernel $k(x, y) = (c^2 + (x - y)^2)^{-\beta} (\beta, c > 0).$

• For the Nyström analysis, assume that $\bar{h}_p(\cdot, X) := h_p(\cdot, X) - \mathbb{E}_{X \sim \mathbb{Q}} h_p(\cdot, X)$ with the sampling distribution \mathbb{Q} is sub-Gaussian, that is,

$$\left| \left\langle \bar{h}_{p}\left(\cdot, X \right), u \right\rangle_{\mathcal{H}_{h_{p}}} \right\|_{\psi_{2}} \lesssim \left\| \left\langle \bar{h}_{p}\left(\cdot, X \right), u \right\rangle_{\mathcal{H}_{h_{p}}} \right\|_{L_{2}(\mathbb{Q})} < \infty$$

$$(1)$$

holds for all $u \in \mathcal{H}_{h_p}$, with a *u*-independent absolute constant in \leq , and $\|\cdot\|_{\psi_2}$ denoting the sub-Gaussian norm.

• Code: https://github.com/flopska/nystroem-ksd

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