A Kernelized Stein Discrepancy for Goodness-of-fit Tests and Model Evaluation

Qiang Liu QLIU@CS.DARTMOUTH.EDU

Computer Science, Dartmouth College, NH, 03755

Jason D. Lee Michael I. Jordan JASONDLEE88@EECS.BERKELEY.EDU JORDAN@CS.BERKELEY.EDU

Department of Electrical Engineering and Computer Science University of California, Berkeley, CA 94709

Abstract

We derive a new discrepancy statistic for measuring differences between two probability distributions based on a novel combination of Stein's method and the reproducing kernel Hilbert space theory. We apply our result to test how well a probabilistic model fits a set of observations, and derive a new class of powerful goodness-of-fit tests that are widely applicable for complex and high dimensional distributions, even for those with computationally intractable normalization constants. Both theoretical and empirical properties of our methods are studied thoroughly.

1. Introduction

Evaluating the goodness-of-fit of models over observed data is a fundamental task in machine learning and statistics. Traditional approaches often involve calculating or comparing the likelihoods or cumulative distribution functions (CDF) of the models. Unfortunately, modern learning techniques increasingly involve complex probabilistic models with computationally intractable likelihoods or CDFs, such as large graphical models, hidden variables models and deep generative models (Koller & Friedman, 2009; Salakhutdinov, 2015). Although Markov chain Monte Carlo (MCMC) or variational methods can be used to approximate the likelihood, their approximation errors are often large and hard to estimate, making it difficult to give results with calibrated statistical significance. In fact, it is often a #P-complete problem to calculate or even approximate likelihoods for graphical models (e.g, Chandrasekaran et al., 2008), making likelihood-based approaches fundamentally infeasible.

We propose a likelihood-free approach for model evalua-

Preliminary work.

tion with guaranteed statistical significance. In particular, we consider the setting of goodness-of-fit testing, where we test whether a given sample $\{x_i\} \sim p(x)$ is drawn from a given distribution q(x), meaning $H_0: p=q$. Our method is based on a novel discrepancy measure between distributions that can be empirically estimated using U-statistics, and depends on q only through its score function $s_q = \nabla_x \log q(x)$; this score function does not depend on the normalization constant in q(x), and can often be calculated efficiently even when the likelihood is intractable. This allows us to apply our methods to complex and high dimensional models on which the likelihood-based methods, or other traditional goodness-of-fit tests, such as χ^2 -test and Kolmogorov-Smirnov test, can not be applied.

Main Idea Our method is motivated by a novel combination of Stein's method and the reproducing kernel Hilbert space (RKHS) theory. Stein's method (Stein, 1972) is a general theoretical tool for obtaining bounds on distances between distributions. It relies on the Stein's identity (Stein et al., 2004), which claims that two smooth densities p(x) and q(x) on $\mathbb R$ are identical if and only if

$$\mathbb{E}_p[\mathbf{s}_q(x)f(x) + \nabla_x f(x)] = 0 \tag{1}$$

for general smooth functions f(x), where $s_p(x) = \nabla_x \log p(x) = \nabla_x p(x)/p(x)$ is called the (Stein) score function of p(x); this identity can be proved by integration by parts. As a result, one can define a Stein discrepancy measure between p and q via

$$\mathbb{S}(p,q) = \max_{f \in \mathcal{H}} \mathbb{E}_p[\boldsymbol{s}_q(x)f(x) + \nabla_x f(x)],$$

where $\mathcal H$ is a set of smooth functions that satisfies (1) and is also rich enough to ensure $\mathbb S(p,q)>0$ whenever $p\neq q$. The problem, however, is that $\mathbb S(p,q)$ is often computationally intractable because it requires a difficult variational optimization. As a result, $\mathbb S(p,q)$ is rarely used in practical machine learning; perhaps the only exception is the recent work of Gorham & Mackey (2015) who obtained

a computationally tractable form by enforcing smoothness constraints only on a finite number of points, turning the optimization into a quadratic programming.

We propose a simpler method for obtaining computational tractable Stein discrepancy $\mathbb{S}(p,q)$ by taking \mathcal{H} to be the unit ball of a reproducing kernel Hilbert space (RKHS) associated with a smooth positive semi-definite kernel k(x,x'). In particular, we show that in this case

$$\mathbb{S}(p,q) = \mathbb{E}_{x,x' \sim p} [u_q(x,x')], \tag{2}$$

where x,x' are i.i.d. random variables drawn from p and $u_q(x,x')$ is a function (defined in Theorem 3.5) that depends on q only through the score function $\nabla_x \log q(x)$ which can be calculated efficiently even when q has an intractable normalization constant. Specifically, assuming q(x) = f(x)/Z with $Z = \int f(x)dx$ being the normalization constant, we have $s_q = \nabla_x \log f(x)$, independent of Z; calculating Z involves a high dimension integration, and has been the major challenge for likelihood-based and Bayesian methods for model evaluation (e.g., Chen et al., 2012).

With an i.i.d. sample $\{x_i\}$ drawn from the (unknown) p(x), the form (2) also enables efficient empirical estimation of $\mathbb{S}(p,q)$ via a U-statistic,

$$\hat{\mathbb{S}}(p,q) = \frac{1}{n(n-1)} \sum_{i \neq j} u_q(x_i, x_j).$$
 (3)

The distribution of $\hat{\mathbb{S}}(p,q)$ can be well characterized using the theory of U-statistics (Hoeffding, 1948; Serfling, 2009), allowing us to solve the goodness-of-fit test (p=q) by testing:

$$H_0: \mathbb{E}_p[u_q(x, x')] = 0$$
 vs. $H_1: \mathbb{E}_p[u_q(x, x')] > 0$.

Outline The rest of the paper is organized as follows. Section 2 introduces backgrounds on RKHS and Stein's method. Section 3 defines our kernelized Stein discrepancy (KSD) and studies its main properties, and Section 4 discusses the empirical estimation of KSD and its application in goodness-of-fit tests. We review several related methods in Section 5, present experiments in Section 6 and conclude the paper in Section 7.

Notations Throughout this paper, we denote by \mathcal{X} a subset of d-dimensional real space \mathbb{R}^d . For a vector-valued function $\boldsymbol{f}(x) = [f_1(x), \dots, f_{d'}(x)]$, its derivative $\nabla_x \boldsymbol{f}(x) = [\frac{\partial f_j(x)}{\partial x_i}]_{ij}$ is a $d \times d'$ matrix-valued function. For a two-variable function (kernel) k(x,x'), we use $k(\cdot,x')=k_{x'}(\cdot)$ to refer to a function of x indexed by fixed x'. For technical simplicity, we will assume all the functions we encounter are absolutely integrable, so that the Fubini-Tonelli theorem can be used to exchange the orders of integrals and infinite sums.

2. Backgrounds

We introduce backgrounds on positive semi-definite kernels and their related reproducing kernel Hilbert spaces (RKHS) in Section 2.1, and Stein's identity and operator in Section 2.2.

2.1. Kernels and Reproducing Kernel Hilbert Spaces

A symmetric kernel k(x, x') is said to be positive semidefinition, if for any function g

$$\int_{\mathcal{X}} g(x)k(x, x')g(x')dxdx' \ge 0.$$

The spectral decomposition of a positive semi-definite kernel k(x, x'), as implied by Mercer's theorem, is defined as

$$k(x, x') = \sum_{j} \lambda_j e_j(x) e_j(x'), \tag{4}$$

where $\{e_j\}$, $\{\lambda_j\}$ are the orthonormal eigenfunctions and positive eigenvalues of k(x,x'), respectively, satisfying $\int e_i(x)e_j(x)dx = \mathbb{I}[i=j]$, for $\forall i,j$.

For a positive semi-definite kernel k(x,x'), its related RKHS $\mathcal H$ comprises of linear combinations of its eigenfunctions, i.e., $f(x) = \sum_j f_j e_j(x)$ with $\sum_j f_j^2/\lambda_j < \infty$, endowed with an inner product $\langle f,g\rangle_{\mathcal H} = \sum_j f_j g_j/\lambda_j$ between f(x) and $g(x) = \sum_j g_j e_j(x)$. Thus this Hilbert space is equipped with a norm $||f||_{\mathcal H}$ where $||f||_{\mathcal H} = \langle f,f\rangle_{\mathcal H} = \sum_j f_j^2/\lambda_j$. One can verify that $k(x,\cdot)$ is in $\mathcal H$ and satisfies the important "reproducing" property,

$$f(x) = \langle f, k(\cdot, x) \rangle_{\mathcal{H}}, \quad k(x, x') = \langle k(\cdot, x'), k(\cdot, x) \rangle_{\mathcal{H}}.$$

Every positive definite kernel k defines a unique RKHS for which k is a reproducing kernel.

We denote by $\mathcal{H}^d = \mathcal{H} \times \cdots \mathcal{H}$ the Hilbert space of $d \times 1$ vector-valued functions $\mathbf{f} = \{f_\ell \colon f_\ell \in \mathcal{H}\}_{\ell \in [d]}$, equipped with an inner product $\langle \mathbf{f}, \mathbf{g} \rangle_{\mathcal{H}^d} = \sum_{\ell \in [d]} \langle f_\ell, g_\ell \rangle_{\mathcal{H}}$ for \mathbf{f} and $\mathbf{g} = \{g_\ell\}_{\ell \in [d]}$, and hence norm $||\mathbf{f}||_{\mathcal{H}} = \sqrt{\sum_{\ell} ||f_\ell||_{\mathcal{H}}^2}$.

2.2. Stein's Identity and Operator

We introduce the basics of Stein's identity and operator, which play a central role in our method.

Definition 2.1. Assume that \mathcal{X} is a subset of \mathbb{R}^d and p(x) a continuous differentiable (smooth) density whose support is \mathcal{X} . The (Stein) score function of p is defined as

$$s_p = \nabla_x \log p(x) = \frac{\nabla_x p(x)}{p(x)}.$$

We say that a function $f: \mathcal{X} \to \mathbb{R}$ is in the **Stein class** of p

if f is smooth and satisfies

$$\int_{x \in \mathcal{X}} \nabla_x (f(x)p(x)) dx = 0.$$
 (5)

The **Stein's operator** of p is a linear operator acting on the Stein class of p, defined as

$$\mathcal{A}_p f(x) = \mathbf{s}_p(x) f(x) + \nabla_x f(x).$$

Note that both s_p and $A_p f$ are $d \times 1$ vector-valued functions mapping from \mathcal{X} to \mathbb{R}^d . A vector-valued function $\mathbf{f}(x) = [f_1(x), \dots, f_{d'}(x)]$ is said to be in the Stein class of p if all f_i , $\forall i \in [d']$ is in the Stein class of p. Applying A_p on a vector-valued $\mathbf{f}(x)$ results a $d \times d'$ matrix-valued function, $A_p \mathbf{f}(x) = \mathbf{s}_p(x) \mathbf{f}(x)^\top + \nabla_x \mathbf{f}(x)$.

Remark The condition (5) can be easily checked using integration by parts or divergence theorem; in particular, when $\mathcal{X} = \mathbb{R}^d$, (5) holds if

$$\lim_{x \to \infty} f(x)p(x) = 0,$$

which holds for regular p(x) if $\lim_{x\to\infty} f(x) = 0$. When $\mathcal X$ is a compact subset of $\mathbb R^d$ with piecewise smooth boundary $\partial \mathcal X$, then by divergence theorem (Marsden & Tromba, 2003), (5) holds if f(x)p(x) = 0 for $\forall x \in \partial \mathcal X$, or more generally if

$$\oint_{\partial \mathcal{X}} p(x)f(x) \cdot \vec{n}(x)dS(x) = 0,$$

where $\vec{n}(x)$ is the unit normal to the boundary $\partial \mathcal{X}$; $\oint_{\partial} dS(x)$ denotes the surface integral over $\partial \mathcal{X}$.

Lemma 2.2 (Stein Lemma, Stein et al. (2004)). Assume p(x) is a smooth density supported on \mathcal{X} , then

$$\mathbb{E}_p[\boldsymbol{\mathcal{A}}_p\boldsymbol{f}(x)] = \mathbb{E}_p[\boldsymbol{s}_p(x)\boldsymbol{f}(x)^\top + \nabla \boldsymbol{f}(x)] = 0.$$

for any f that is in the Stein class of p.

Proof. By the definition of the Stein class, simply note that
$$s_p(x) f(x)^\top + \nabla f(x) = \nabla_x (f(x)p(x))/p(x)$$
.

The following simple result provides a convenient tool for our derivation; it relates the expectation under p of Stein's operator $\mathcal{A}_q f$ with the difference of the score functions of p and q.

Lemma 2.3 (Ley & Swan (2013)). Assume p(x) and q(x) are smooth densities supported on \mathcal{X} and f(x) is in the Stein class of p, we have

$$\mathbb{E}_p[\mathcal{A}_q \boldsymbol{f}(x)] = \mathbb{E}_p[(\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x))\boldsymbol{f}(x)^\top].$$

Proof. Since
$$\mathbb{E}_p[\mathcal{A}_p \boldsymbol{f}(x)] = 0$$
 by Lemma 2.2, we have $\mathbb{E}_p[\mathcal{A}_q \boldsymbol{f}(x)] = \mathbb{E}_p[\mathcal{A}_q \boldsymbol{f}(x) - \mathcal{A}_p \boldsymbol{f}(x)] = \mathbb{E}_p[(\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x))\boldsymbol{f}(x)^\top].$

Therefore, $\mathbb{E}_p[\mathcal{A}_q \boldsymbol{f}(x)]$ is the $\boldsymbol{f}(x)$ -weighted expectation of the score function difference $(\boldsymbol{s}_q(x)-\boldsymbol{s}_p(x))$ under p. When $\boldsymbol{f}(x)$ is a $d\times 1$ vector-valued function, $\mathbb{E}_p[\mathcal{A}_q \boldsymbol{f}(x)]$ is a $d\times d$ matrix; taking its trace gives $\mathbb{E}_p[\mathrm{trace}(\mathcal{A}_q \boldsymbol{f}(x))] = \mathbb{E}_p[(\boldsymbol{s}_q(x)-\boldsymbol{s}_p(x))^{\top} \boldsymbol{f}(x)].$

3. Kernelized Stein Discrepancy

We introduce our kernelized Stein discrepancy (KSD) with an elementary definition motivated by Lemma 2.3, and then establish its connection with Stein's method and RKHS.

Definition 3.1. Given a positive semi-definite kernel $k: \mathcal{X} \times \mathcal{X} \to \mathbb{R}$, the **kernelized Stein discrepancy** (KSD) $\mathbb{S}(p,q)$ between distribution p and q is defined as

$$\mathbb{S}(p,q) = \mathbb{E}_{x,x' \sim p} [\boldsymbol{\delta}_{q,p}(x)^{\top} k(x,x') \boldsymbol{\delta}_{q,p}(x)], \quad (6)$$

where $\delta_{q,p}(x) = s_q(x) - s_p(x)$ is the score difference between p and q, and x, x' are i.i.d. draws from p(x).

The following result shows that $\mathbb{S}(p,q)$ is a valid discrepancy measure between p and q when $\boldsymbol{\delta}_{p,q} = \boldsymbol{s}_q - \boldsymbol{s}_p \in \mathcal{H}^d$.

Proposition 3.2. For any semi-definite kernel k(x, x') associated with a reproducing kernel Hilbert space (RHKS) \mathcal{H} , we have $\mathbb{S}(p,q) \geq 0$ for any p, q with smooth densities. In addition, when $\boldsymbol{\delta}_{q,p} = \boldsymbol{s}_q - \boldsymbol{s}_p \in \mathcal{H}^d$, we have $\mathbb{S}(p,q) = 0$ if and only if p = q.

Proof. $\mathbb{S}(p,q) \geq 0$ is a direct result of the positive semi-definiteness of k(x,x'). Because $\pmb{\delta}_{q,p} = \{\pmb{\delta}_{p,q}^{\ell}\}_{\ell \in [d]} \in \mathcal{H}^d$, we have $\pmb{\delta}_{q,p}^{\ell}(x) = \sum_j \alpha_{j\ell} e_j(x), \, \forall \ell \in [d], \, \text{and} \, \mathbb{S}(p,q) = \sum_\ell \sum_j \lambda_j \alpha_{j\ell}^2$ by the spectral representation (4). Therefore, $\mathbb{S}(p,q) = 0$ implies $\alpha_{j\ell} = 0$ for $\forall j,\ell$, which then implies $\pmb{s}_p = \pmb{s}_q$ and hence p = q.

Proposition 3.2 requires $(s_q - s_p) \in \mathcal{H}^d$ to make $\mathbb{S}(p,q)$ a valid discrepancy measure. Therefore, we would need to take \mathcal{H} to be rich enough to capture a broad class of score differences. This can be achieved by taking k(x,x') to be a universal kernel, whose related RKHS \mathcal{H} is a dense subset of continuous functions on compact subsets of \mathbb{R}^d under L_∞ norm (Steinwart & Christmann, 2008). Many commonly used kernels, such as the RBF kernel, have been shown to be universal (Steinwart & Christmann, 2008).

The $\mathbb{S}(p,q)$ as defined in (6) requires to know both s_p and s_q ; we now apply Stein's identity to derive the more convenient form (2) that only requires s_q .

Definition 3.3. A kernel k(x, x') is said to be in the Stein class of p if k(x, x') has continuous second order partial derivatives, and both $k(x, \cdot)$ and $k(\cdot, x)$ are in the Stein class of p for any fixed x.

It is easy to check that the RBF kernel $k(x,x') = \exp(-\frac{1}{2h^2}||x-x'||_2^2)$ is in the Stein class for regular smooth densities.

Proposition 3.4. If k(x, x') is in the Stein class of p, so is any $f \in \mathcal{H}$.

Theorem 3.5. Assume p and q are smooth densities and k(x, x') is in the Stein class of p. Define

$$u_q(x, x') = \boldsymbol{s}_q(x)^{\top} k(x, x') \boldsymbol{s}_q(x') + \boldsymbol{s}_q(x)^{\top} \nabla_{x'} k(x, x') + \nabla_x k(x, x')^{\top} \boldsymbol{s}_q(x') + \operatorname{trace}(\nabla_{x, x'} k(x, x')).$$

then

$$\mathbb{S}(p,q) = \mathbb{E}_{x,x' \sim p}[u_q(x,x')]. \tag{7}$$

Proof. Apply Lemma 2.3 twice, first on $k(\cdot, x')$ for fixed x', and then with fixed x. See the Appendix.

The representation in (7) is of central importance for our framework, since it provides a tractable formula for empirical evaluation of $\mathbb{S}(p,q)$ and its confidence interval based on the sample $\{x_i\} \sim p$ and score function s_q ; see Section 4 for further discussion.

Using the spectral decomposition of k(x, x'), we can show that $\mathbb{S}(p, q)$ is effectively applying Stein's operator simultaneously on all the eigenfunctions $e_i(x)$ of k(x, x').

Theorem 3.6. Assume k(x,x') is a semi-definite kernel in the Stein class of p, with positive eigenvalues $\{\lambda_j\}$ and eigenfunctions $\{e_j(x)\}$, then $u_q(x,x')$ is also a positive semi-definite kernel, and can be rewritten into

$$u_q(x, x') = \sum_{j} \lambda_j \left[\mathcal{A}_q e_j(x) \right]^\top \left[\mathcal{A}_q e_j(x') \right], \quad (8)$$

where $A_q e_j(x) = s_q(x)e_j(x) + \nabla_x e_j(x)$ is the Stein's operator acted on e_j . In addition,

$$\mathbb{S}(p,q) = \sum_{j} \lambda_j ||\mathbb{E}_{x \sim p}[\mathcal{A}_q e_j(x)]||_2^2. \tag{9}$$

Note that although $\{e_j\}$ are orthonormal, the $\{A_q e_j(x)\}$ are no longer orthonormal in general.

Finally, we are ready to establish the variational interpretation of $\mathbb{S}(p,q)$ that motivated this work, that is, it can be treated as the maximum of $\mathbb{E}_{x \sim p}[\mathcal{A}_q f(x)]$ when optimizing f in the unit ball of RKHS \mathcal{H} related to kernel k(x,x').

Theorem 3.7. Let \mathcal{H} be the RKHS related to a positive semi-definite kernel k(x, x') in the Stein class of p. Denote by $\beta(x') = \mathbb{E}_{x \sim p}[A_q k_{x'}(x)]$, then

$$\mathbb{S}(p,q) = ||\boldsymbol{\beta}||_{\mathcal{H}^d}^2. \tag{10}$$

Further, we have $\langle \boldsymbol{f}, \boldsymbol{\beta} \rangle_{\mathcal{H}^d} = \mathbb{E}_x[\operatorname{trace}(\mathcal{A}_q \boldsymbol{f})]$ for $\boldsymbol{f} \in \mathcal{H}^d$, and hence

$$\sqrt{\mathbb{S}(p,q)} = \max_{\boldsymbol{f} \in \mathcal{H}^d} \left\{ \mathbb{E}_x[\operatorname{trace}(\mathcal{A}_q \boldsymbol{f})] \quad s.t. \quad ||\boldsymbol{f}||_{\mathcal{H}^d} \le 1 \right\}$$
(11)

where the maximum is achieved when $f = \beta$.

4. Goodness-of-fit Testing Based on KSD

The form in (7) allows efficient estimation of $\mathbb{S}(p,q)$ in practice. Given i.i.d. sample $\{x_i\}$ drawn from an unknown p and the score function $s_q(x)$, we can estimate $\mathbb{S}(p,q)$ by

$$\hat{\mathbb{S}}_{u}(p,q) = \frac{1}{n(n-1)} \sum_{1 < i \neq j < n} u_{q}(x_{i}, x_{j}), \quad (12)$$

where $\hat{\mathbb{S}}_u(p,q)$ is a form of U-statistics ("U" stands for unbiasedness), which provides a minimum-variance unbiased estimator for $\mathbb{S}(p,q)$ (Hoeffding, 1948; Serfling, 2009). We can also estimate $\mathbb{S}(p,q)$ using a V-statistic of form $\frac{1}{n^2}\sum_{i,j=1}^n u_q(x_i,x_j)$, which provides a biased estimator, but has the advantage of always being nonnegative since $u_q(x,x')$ is positive semi-definite. We will focus on the U-statistic in this work because of its unbiasedness.

Theorem 4.1. Let k(x, x') be a positive semi-definite kernel in the Stein class of p and q. Assume $s_q - s_p \in \mathcal{H}^d$ and $\mathbb{E}_{x,x' \sim p}[u_q(x,x')^2] < \infty$, we have

1) If $p \neq q$, then $\hat{\mathbb{S}}_{u}(p,q)$ is asymptotically normal with

$$\sqrt{n}(\hat{\mathbb{S}}_u(p,q) - \mathbb{S}(p,q)) \xrightarrow{d} \mathcal{N}(0,\sigma_u^2),$$

where $\sigma_u^2 = \operatorname{var}_x(\mathbb{E}_{x'}[u_q(x, x')])$ and $\sigma_u^2 \neq 0$.

2) If p=q, then we have $\sigma_u^2=0$ (the U-statistics is degenerate) and

$$n\hat{\mathbb{S}}_u(p,q) \stackrel{d}{\to} \sum_{i=1}^{\infty} c_j(Z_j^2 - 1),$$
 (13)

where $\{Z_j\}$ are i.i.d. standard Gaussian random variables, and $\{c_j\}$ are the eigenvalues of kernel $u_q(x,x')$ under p(x), that is, they are the solutions of $c_j\phi_j(x) = \int_{x'} u_q(x,x')\phi_j(x')p(x')dx'$ for non-zero ϕ_j .

Proof. Using the standard asymptotic results of *U*-statistics in Serfling (2009, Section 5.5), we just need to check that $\sigma_u^2 \neq 0$ when $p \neq q$ and $\sigma_u^2 = 0$ when p = q. See Appendix for details.

Algorithm 1 Bootstrap Goodness-of-fit Test based on KSD

Input: Sample $\{x_i\}$ and score function $s_q(x) = \nabla_x \log q(x)$. Bootstrap sample size m.

Test: H_0 : $\{x_i\}$ is drawn from q v.s H_1 : $\{x_i\}$ is not drawn from q.

- 1. Compute $\hat{\mathbb{S}}_u$ by (12) and $u_q(x, x')$ as defined in Theorem 3.5. Generate m bootstrap sample $\hat{\mathbb{S}}_u^*$ by (14).
- 2. Reject H_0 with significance level α if the percentage of $\hat{\mathbb{S}}_u^*$ that satisfies $\hat{\mathbb{S}}_u^* > \hat{\mathbb{S}}_u$ is less than α .

Theorem 4.1 suggests that $n\hat{\mathbb{S}}_u(p,q)$ has a well defined limit distribution under the null p=q, that is, $n\hat{\mathbb{S}}_u(p,q)<\infty$ with probability one, but grows to ∞ at a \sqrt{n} -rate under any fixed alternative hypothesis $q\neq p$. This suggests a straightforward goodness-of-fit testing procedure: Denote by $F_{\hat{\mathbb{S}}_u}$ the CDF of $n\hat{\mathbb{S}}_u$ under the null p=q, and set $\gamma_{1-\alpha}$ the $1-\alpha$ quantile of $F_{\hat{\mathbb{S}}_u}$, i.e., $\gamma_{1-\alpha}=\inf\{s\colon F_{\hat{\mathbb{S}}_u}(s)\geq 1-\alpha\}$, then we reject the null with significant level α if $n\hat{\mathbb{S}}_u\geq \gamma_{1-\alpha}$.

Proposition 4.2. Assume the conditions in Theorem 4.1. For any fixed $q \neq p$, the limiting power of the test that rejects the null p = q when $n\hat{\mathbb{S}}_u(p,q) > \gamma_{1-\alpha}$ is one, that is, the test is consistent in power against any fixed $q \neq p$.

One difficulty in implementing this test is that the limit distribution in (13) and its α -quantile does not have analytic form unless $c_j=0$, or 1. Fortunately, the same type of asymptotics appears in many other classical goodness-of-fit tests, such as Cramer-von Mises test, Anderson-Darling test, as well as two-sample tests (Gretton et al., 2012). As a consequence, a line of work has been devoted to approximating the critical values of (13), including bootstrap methods (Arcones & Gine, 1992; Huskova & Janssen, 1993; Chwialkowski et al., 2014) and eigenvalue approximation (Gretton et al., 2009).

In this work, we adopt the bootstrap method suggested in Huskova & Janssen (1993); Arcones & Gine (1992): We repeatedly draw multinomial random weights $(w_1, \ldots w_n) \sim \operatorname{Mult}(n \; ; \; \frac{1}{n}, \ldots, \frac{1}{n})$, and calculate bootstrap sample

$$\hat{\mathbb{S}}_{u}^{*}(p,q) = \sum_{i \neq j} (w_{i} - \frac{1}{n})(w_{j} - \frac{1}{n})u_{q}(x_{i}, x_{j}), \quad (14)$$

and then calculate the empirical quantile $\hat{\gamma}_{1-\alpha}$ of $n\hat{\mathbb{S}}_u^*(p,q)$. The consistency of $\hat{\gamma}_{1-\alpha}$ for degenerate U-statistics has been established in Arcones & Gine (1992); Huskova & Janssen (1993).

Theorem 4.3 (Huskova & Janssen (1993)). Assume the conditions in Theorem 4.1. If p = q, then as the bootstrap

sample size $m \to \infty$,

$$\sup_{s \in \mathbb{R}} \left| \Pr(n \hat{\mathbb{S}}_u^* \le s \mid \{x_i\}_{i=1}^n) - \Pr(n \hat{\mathbb{S}}_u \le s) \right| \to 0,$$

that is, the bootstrap test attains the correct significance level asymptotically (consistent in level).

It is important to note, on the other hand, that the more usual bootstrap, such as $\sum_{i\neq j} w_i w_j u_q(x_i, x_j)$, may not work for degenerate U-statistics as discussed in Arcones & Gine (1992).

This bootstrap test is summarized in Algorithm 1; its cost is $O(mn^2)$ where n is the size of the sample $\{x_i\}$ and m the bootstrap sample size. A more computationally efficient, but less statistically powerful, method can be constructed based on the following linear estimator:

$$\hat{\mathbb{S}}_{lin} = \frac{1}{\lfloor n/2 \rfloor} \sum_{i=1}^{\lfloor n/2 \rfloor} u_q(x_{2i-1}, x_{2i}), \tag{15}$$

which has a zero-mean Gaussian limit under the null. This gives a test with only O(n) time complexity: reject the null if $\hat{\mathbb{S}}_{lin} > \hat{\sigma}z_{1-\alpha}$, where $z_{1-\alpha}$ is the $1-\alpha$ quantile of the standard Gaussian distribution, and $\hat{\sigma}$ the standard deviation of $\{u_q(x_{2i-1},x_{2i})\}$. This test, however, tends to perform much worse than the U-statistic based test as we show in our experiments. Further computation-efficiency tradeoff between the linear- and U-statistic can be obtained by block-wise averaging; see Ho & Shieh (2006); Zaremba et al. (2013) for details.

5. Related Methods

KSD has close connections with Fisher divergence and maximum discrepancy measure, which we discuss in Section 5.1 and Section 5.2, respectively.

5.1. Connection with Fisher Divergence

Fisher divergence, also known as Fisher information distance (Johnson, 2004), is defined as

$$\mathbb{F}(p,q) = \mathbb{E}_x \left[||\nabla_x \log p(x) - \nabla_x \log q(x)||_2^2 \right], \quad (16)$$

that is, it is the $\mathcal{L}^2(p)$ norm of $s_q(x)-s_p(x)$. An immediate connection is made by noting that $\mathbb{F}(p,q)$ can be treated as a special case of $\mathbb{S}(p,q)$ defined in (6) with $k(x,x')=\mathbb{I}[x=x']$, or a RBF kernel with bandwidth $h\to 0$; in this sense, we can also think KSD as a kernelized version of Fisher divergence. We can establish the follow inequalities between $\mathbb{F}(p,q)$ and $\mathbb{S}(p,q)$:

Theorem 5.1. 1) Following Definition (6) and (16), we have

$$|\mathbb{S}(p,q)| \le \sqrt{\mathbb{E}_{x,x' \sim p}[k(x,x')^2]} \cdot \mathbb{F}(p,q). \tag{17}$$

2) In addition, if k(x, x') is positive semi-definite and in the Stein class of p, and $s_q - s_p \in \mathcal{H}^d$, we have, for $p \neq q$,

$$\sqrt{\mathbb{S}(p,q)} \ge \mathbb{F}(p,q) / ||\boldsymbol{s}_q - \boldsymbol{s}_p||_{\mathcal{H}^d}. \tag{18}$$

Proof. (17) is a simple result of Cauchy-Schwarz inequality, and (18) can be obtained by taking $f=(s_q-s_p)/||s_q(x)-s_p(x)||_{\mathcal{H}^d}$ in (11). See Appendix.

(17) suggests that the convergence in Fisher divergence is stronger than that in KSD. In fact, using Stein's method, Ley & Swan (2013) showed that Fisher divergence is stronger than most other divergences, including KL, total variation and Hellinger distances.

In addition, we can also represent $\mathbb{F}(p,q)$ in a variational form similar to (11) but with f optimized over the unit ball of the intersection of $\mathcal{L}^2(p)$ space and the Stein class of p.

Proposition 5.2. Let $\mathcal{F}(p) = \mathcal{L}^2(p) \cap \mathcal{S}(p)$, where $\mathcal{S}(p)$ represents the Stein class of p, then we have

$$\sqrt{\mathbb{F}(p,q)} \ge \max_{\boldsymbol{f} \in \mathcal{F}(p)^d} \left\{ \mathbb{E}_p[\operatorname{trace}(\mathcal{A}_q \boldsymbol{f}(x))] \right.$$
s.t. $\mathbb{E}_p[||\boldsymbol{f}(x)||_2^2] \le 1$.

and the equality holds when $s_q - s_p \in \mathcal{F}(p)^d$.

Note that $\mathcal{L}^2(p)$ is larger than the Stein class and RKHS, and includes discontinuous, non-smooth functions, and hence we need to ensure f is in the Stein class explicitly.

Despite the connections, the critical disadvantage of Fisher divergence compared to KSD is that the computationally convenient representation (7) no longer holds for Fisher divergence, because its corresponding kernel $\mathbb{I}[x=x']$ is not differentiable. Therefore, we can not estimate $\mathbb{F}(p,q)$ using the U-statistic in (12). Instead, estimating $\mathbb{F}(p,q)$ seems to be substantially more difficult. To see this, note that

$$\mathbb{F}(p,q) = \mathbb{E}_{x \sim p}[||s_q(x)||_2^2 - 2s_p(x)^{\top} s_q(x) + ||s_p(x)||_2^2]$$

$$= \mathbb{E}_{x \sim p}[\phi_q(x)] + \mathbb{E}_{x \sim p}||s_p(x)||_2^2, \tag{19}$$

where $\phi_q(x) = ||s_q(x)||_2^2 + 2 \mathrm{trace}(\nabla_x s_q(x))$ and is obtained by applying Stein's identity on the cross term. Note that although the first term $\mathbb{E}_{x \sim p}[\phi_q(x)]$ in (19) can be estimated by the empirical mean of $\phi_q(x)$ (which only depends on s_q) under sample $\{x_i\} \sim p$, the second term $\mathbb{E}_{x \sim p}||s_p(x)||_2^2$ is more difficult to estimate, since it depends on the score function $s_p(x)$ of the unknown p(x), and hence requires a kernel density estimator for p(x); see Hall & Marron (1987); Birge & Massart (1995). We should point out that similar difficult "constant" terms appear in other common discrepancy measures such as KL divergence and α -divergence (e.g., Krishnamurthy et al., 2014).

For this reason, KSD provides a much more convenient tool for goodness-of-fit tests than the other discrepancies.

Meanwhile, Fisher divergence still has the advantage of being independent of the normalization constants of p and q, and provides a useful tool in cases when it does not require evaluating the term $s_p(x)$. For example, Fisher divergence has been widely used for parameter estimation, finding the optimal q(x) that best fits a sample $\{x_i\}$ by minimizing $\mathbb{F}(p,q)$; this yields the score matching methods developed in both parametric (Hyvärinen, 2005; Lyu, 2009) and non-parametric (Sriperumbudur et al., 2013) settings. We also believe that Fisher divergence can provide a promising tool for model comparison or selection, that is, given two candidate models q_1 and q_2 , decide which one fits the data $\{x_i\}$ better; this seems to be a natural idea, while we are not aware of any existing work in this direction.

5.2. Maximum Mean Discrepancy & Two-sample Tests

Closely related to goodness-of-fit tests are two sample tests, which test whether two i.i.d. samples $\{x_i\}$ and $\{y_i\}$ are drawn from the same distribution. In principle, one can turn a goodness-of-fit test into a two sample test by drawing $\{y_i\}$ from q(x). However, it is often difficult to draw exact i.i.d. samples for practical models, and furthermore MCMC sampling may be computationally expensive, suffer from the convergence problems, and introduce undesired correlations. When the MCMC approximation is poor, the two sample test would reject the null even when p=q (inconsistent in level).

Maximum Mean Discrepancy (Gretton et al., 2012) is a nonparametric distance measure widely used for two sample tests, defined as

$$\mathbb{M}(p,q) = \max_{h \in \mathcal{H}} \left\{ \mathbb{E}_p[h(x)] - \mathbb{E}_q[h(x)] \quad s.t. \quad ||h||_{\mathcal{H}} \le 1 \right\},$$

where \mathcal{H} is the RKHS of kernel k(x,x'). Gretton et al. (2012) showed that $\mathbb{M}(p,q)$ can be rewritten into

$$\mathbb{M}(p,q) = \mathbb{E}_{x,x';y,y'}[k(x,x') + k(y,y') - k(x,y') - k(x',y)],$$

where x,x' and y,y' are i.i.d. draws from p and q, respectively. Therefore, $\mathbb{M}(p,q)$ can be empirically estimated based on sample $x_i \sim p$ and $y_i \sim q$ using U- or V- statistics, making it a useful tool for two sample tests. Our KSD, on the other hand, is better estimated with sample $x_i \sim p$ and the score function s_q and hence suitable for goodness-of-fit tests. Finally, we remark that MMD is symmetric, while KSD is asymmetric.

6. Experiments

We present empirical results in this section. We start with a toy case of 1D Gaussian mixture on which we can compare with the classical goodness-of-fit tests that only work

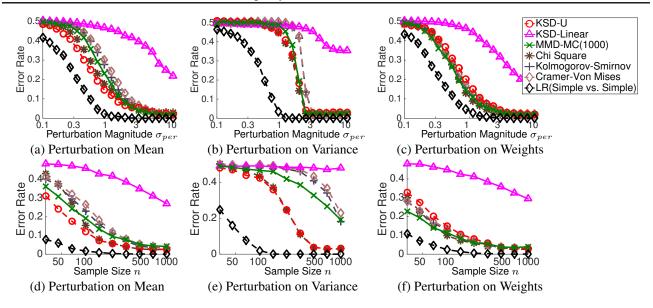


Figure 1. Results on 1D Gaussian mixture. (a)-(c) The error rates of different methods vs. the perturbation magnitude σ_{per} when perturbing the mean, variance and mixture weights, respectively; we use a fixed sample size of n=100. (d)-(f) the error rates vs. the sample size n, with fixed perturbation magnitude $\sigma_{per}=1$.

for univariate distributions, and then proceed to Gaussian-Bernoulli restricted Boltzmann machine (RBM), a graphical model widely used in deep learning (Welling et al., 2004; Hinton & Salakhutdinov, 2006). The following methods are evaluated, all with a significance level of 0.05:

- 1) KSD-U. The KSD-based bootstrap test using U-statistics as shown in Algorithm 1. We use the RBF kernel with bandwidth chosen to be median of the data distances.
- 2) KSD-Linear. The KSD test based on the linear estimator in (15) with asymptotically normal null distribution.
- 3) Classical goodness-of-fit tests, including χ^2 test, Kolmogorov-Smirnov test and Cramer-von Mises test (Lehmann & Romano, 2006); they are evaluated on only the 1D Gaussian mixture.
- 4) MMD-MC (n'). Draw exact sample $\{y_i\}$ of size n' from q(x) and perform two sample MMD test over $\{x_i\}$ and $\{y_i\}$ using bootstrap¹; see Gretton et al. (2012).
- 5) MMD-MCMC (n'). Draw approximate sample $\{y_i\}$ of size n' from q(x) using Gibbs sampler and perform MMD test on $\{x_i\}$ and $\{y_i\}$; we use 1000 burn-in steps.
- 6) LR (simple vs. simple). We evaluate the exact log-likelihood ratio $2\log(q(x)/p(x))$ and use it to test whether $\{x_i\}$ is drawn from p(x) or q(x). This approach is an oracle test in that it knows it exactly calculates the likelihood, and assumes we know p(x) and tests a much easier null hypothesis of simple vs. simple.

7) Likelihood Ratio (AIS). We approximately evaluate the likelihood ratio using annealed importance sampling (AIS), which is one of the most widely used algorithm for approximating likelihood (Neal, 2001; Salakhutdinov & Murray, 2008). Our AIS implementation uses a Gibbs sampler transition with a linear temperature grid of size 1000. We do not perform a test based on the AIS result because it is hard to know the approximation error.

1D Gaussian Mixture We draw i.i.d. sample $\{x_i\}_{i=1}^n$ from $p(x) = \sum_{k=1}^5 w_k \mathcal{N}(x; \mu_k, \sigma^2)$ with $w_k = 1/5$, $\sigma = 1$ and μ_k randomly drawn from Uniform[0, 10]. We then generate q(x) by adding Gaussian noise on μ_k , $\log w_k$, or $\log \sigma^2$, leading to three different ways for perturbation; the perturbation magnitude is controlled by the variance σ_{per}^2 of the Gaussian noise. In our experiment, we set q(x) randomly with equal probability to be either the true model p(x) ($H_0: p=q$), or the perturbed version ($H_1: p \neq q$), and then use different methods to test H_0 vs. H_1 . We repeat 1000 random trials, and report the average error rate in Figure 1.

We find from Figure 1 that the oracle LR (simple vs. simple) performs the best as expected. Otherwise, our KSD-U performs comparably with, or better than, the classical tests (χ^2 , Kolmogorov-Smirnov and Cramer-Von Mises) as well as MMD-MC (1000). KSD-Linear tends to perform the worst, suggesting it is not useful in this simple setting. However, it can serve as a computationally efficient alternative of KSD-U for more complex models on which the other tests are not practical. Note that because both the cases of p=q and $p\neq q$ happen with 0.5 prob-

¹We use the mmdTestBoot.m under http://www.gatsby.ucl.ac.uk/%7Egretton/mmd/mmd.htm

ability in our simulation, the error rate in the hardest case when p is close q is 0.5.

Gaussian-Bernoulli Restricted Boltzmann Machine (RBM) Gaussian-Bernoulli RBM is a hidden variable graphical models consist of a continuous observable variable $x \in \mathbb{R}^d$ and a binary hidden variable $h \in \{\pm 1\}^{d'}$, with joint probability

$$p(x,h) = \frac{1}{Z} \exp(\frac{1}{2} x^{\top} B h + b^{\top} x + c^{\top} h - \frac{1}{2} ||x||_2^2),$$

where Z is the normalization constant. The probability of the observable variable x is $p(x) = \sum_{h \in \{\pm 1\}^{d'}} p(x,h)$, which is intractable to calculate due to the difficult constant term Z. Nevertheless, one can show that its score function s_p can be easily calculated in a closed form,

$$s_p(x) = b - x + B \ \phi(B^{\top}x + c), \qquad \phi(y) = \frac{e^{2y} - 1}{e^{2y} + 1}.$$

In our experiment, we simulate a true model p(x) by drawing b and c from standard Gaussian and select B uniformly randomly from $\{\pm 1\}$; we use d=50 observable variables and d'=10 hidden variables, so that it remains possible to exactly calculate p(x) and draw exact samples using the brute-force algorithm. Similar to the case of 1D Gaussian mixture, we set q(x) randomly with equal probability to be equal to either p(x) or a perturbed version by adding Gaussian noise to B with variance σ^2_{per} . We report the the error rates of different tests in Figure 2; the results are averaged on 1000 random trials.

Figure 2(a) shows that the oracle LR (simple vs. simple) performs the best again as expected, followed by our KSD-U method. The MMD-MCMC breaks down because the MCMC sample is not representative of q, while the performance of MMD-MC depends on the size of the exact sample: it performs worse than KSD-U with MMD-MC (100), and is almost as good with MMD-MC (1000); see also Figure 2(b). Again, we find that KSD-linear generally performs much worse than KSD-U, but it provides a computationally efficient O(n) alternative to KSD-U which has a $O(mn^2)$ complexity and MMD which costs O(mnn'). A trade-off between linear and quadratic complexity can be achieved using block averaging; see Zaremba et al. (2013).

Figure 2(c)-(h) shows the different discrepancy measures under the case p=q and $p\neq q$, respectively. Again, we can find that the exact likelihood ratio provides the best discrimination, while MMD-MCMC fails to distinguish the two cases at all. The AIS approximation performs reasonably well, but is worse than KSD-U and MMD-MC (1000) in this particular case.

7. Conclusion and Future Directions

We propose a new computationally tractable discrepancy measure between complex probability models, and use it to derive a novel class of goodness-of-fit tests. We believe our discrepancy measure provides a new fundamental tool for analyzing and using complex probability models in statistics and machine learning. Future directions include extending our method to composite goodness-of-fit tests, in which we want to test if the observed data follows a given class of distributions, as well as understanding the theoretical discrimination power of KSD compared to the other classical goodness-of-fit tests, two sample tests (e.g., MMD with infinite exact Monte Carlo sample), and the method in Gorham & Mackey (2015).

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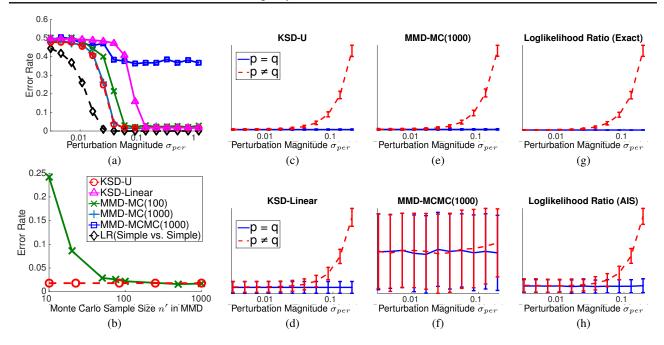


Figure 2. Results on Gaussian-Bernoulli RBM. (a) The error rate vs. the perturbation magnitude σ_{per} . (b) The error rate of MMD-MC vs. the size of the exact sample used. (c) Different discrepancy measures between p and q under the null p=q (blue solid lines) and the alternatives $p \neq q$ (dashed red lines); the x-axes are the deviation σ_{per} between p and q when $p \neq q$. We set n=100 in all the cases.

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Appendix for "A Kernelized Stein Discrepancy for Goodness-of-fit Tests and Model Evaluation"

Proof of Theorem 3.5. 1) Denote by $\mathbf{v}(x,x') = k(x,x')\mathbf{s}_q(x') + \nabla_{x'}k(x,x') = \mathcal{A}_qk_x(x')$; applying Lemma 2.3 on $k(x,\cdot)$ with fixed x,

$$\mathbb{S}(p,q) = \mathbb{E}_{x,x' \sim p}[(\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x))^{\top} k(x,x') (\boldsymbol{s}_q(x') - \boldsymbol{s}_p(x'))]$$
$$= \mathbb{E}_{x,x' \sim p}[(\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x))^{\top} \boldsymbol{v}(x,x')]$$

Because $k(\cdot, x')$ is in the Stein class of p for any x', we can show that $\nabla_{x'}k(\cdot, x')$ is also in the Stein class, because

$$\int_{x} \nabla_{x}(p(x)\nabla_{x'}k(x,x'))dx = \nabla_{x'} \int_{x} \nabla_{x}(p(x)k(x,x'))dx = 0,$$

and hence $v(\cdot, x')$ is also in the Stein class; apply Lemma 2.3 on $v(\cdot, x')$ with fixed x' gives

$$S(p,q) = \mathbb{E}_{x,x' \sim p}[(\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x))^{\top} \boldsymbol{v}(x,x')]$$
$$= \mathbb{E}_{x,x' \sim p}[\boldsymbol{s}_q(x)^{\top} \boldsymbol{v}(x,x') + \operatorname{trace}(\nabla_x \boldsymbol{v}(x,x'))]$$

The result then follows by noting that $\nabla_x \boldsymbol{v}(x,x') = \nabla_x k(x,x') \boldsymbol{s}_q(x')^\top + \nabla_{x'x'} k(x,x')$.

Proof of Theorem 3.6. Note that

$$\nabla_x k(x, x') = \sum_j \lambda_j \nabla_x e_j(x) \ e_j(x'), \qquad \qquad \nabla_{x, x'} k(x, x') = \sum_j \lambda_j \nabla_x e_j(x) \ \nabla_{x'} e_j(x')^\top,$$

and hence

$$\begin{aligned} u_{q}(x, x') &= \boldsymbol{s}_{q}(x)^{\top} k(x, x') \boldsymbol{s}_{q}(x') + \boldsymbol{s}_{q}(x)^{\top} \nabla_{x}' k(x, x') + \boldsymbol{s}_{q}(x')^{\top} \nabla_{x} k(x, x') + \operatorname{trace}(\nabla_{x, x'} k(x, x') \\ &= \sum_{j} \lambda_{j} \left[\boldsymbol{s}_{q}(x)^{\top} e_{j}(x) e_{j}(x') \boldsymbol{s}_{q}(x') + \boldsymbol{s}_{q}(x)^{\top} e_{j}(x) \nabla_{x'} e_{j}(x') + \boldsymbol{s}_{q}(x')^{\top} \nabla_{x} e_{j}(x) e_{j}(x') + \nabla_{x} e_{j}(x)^{\top} \nabla_{x'} e_{j}(x') \right] \\ &= \sum_{j} \lambda_{j} \left[\boldsymbol{s}_{q}(x) e_{j}(x) + \nabla_{x} e_{j}(x) \right]^{\top} \left[\boldsymbol{s}_{q}(x') e_{j}(x') + \nabla_{x'} e_{j}(x') \right] \\ &= \sum_{j} \lambda_{j} [\mathcal{A}_{q} e_{j}(x)]^{\top} [\mathcal{A}_{q} e_{j}(x')]. \end{aligned}$$

Therefore, $u_q(x, x')$ is positive semi-definite because $\lambda_j > 0$. In addition,

$$\begin{split} \mathbb{S}(p,q) &= \mathbb{E}_{x,x'}[u_q(x,x')] \\ &= \sum_j \lambda_j \mathbb{E}_x[\mathcal{A}_q e_j(x)]^\top \mathbb{E}_{x'}[\mathcal{A}_q e_j(x')] \\ &= \sum_j \lambda_j ||\mathbb{E}_x[\mathcal{A}_q e_j(x)]||_2^2. \end{split}$$

Proof of Theorem 3.7. We first prove (10) by applying the reproducing property $k(x, x') = \langle k(x, \cdot), k(x', \cdot) \rangle_{\mathcal{H}}$ on (6):

$$\begin{split} \mathbb{S}(p,q) &= \mathbb{E}_{x,x' \sim p}[(\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x))^\top k(x,x') \left(\boldsymbol{s}_q(x') - \boldsymbol{s}_p(x')\right)] \\ &= \mathbb{E}_{x,x' \sim p}[(\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x))^\top \left\langle k(x,\cdot), \ k(x,\cdot) \right\rangle_{\mathcal{H}} \left(\boldsymbol{s}_q(x') - \boldsymbol{s}_p(x')\right)] \\ &= \sum_{\ell=1}^d \left\langle \mathbb{E}_x[(\boldsymbol{s}_q^\ell(x) - \boldsymbol{s}_p^\ell(x))k(x,\cdot)], \ \mathbb{E}_{x'}[k(x,\cdot)(\boldsymbol{s}_q^\ell(x) - \boldsymbol{s}_p^\ell(x))] \right\rangle_{\mathcal{H}} \\ &= \sum_{\ell=1}^d \left\langle \boldsymbol{\beta}_\ell, \boldsymbol{\beta}_\ell \right\rangle_{\mathcal{H}} \\ &= ||\boldsymbol{\beta}||_{\mathcal{H}^d}^2 \end{split}$$

where we used the fact that $\boldsymbol{\beta}(x') = \mathbb{E}_{x \sim p}[\mathcal{A}_q k_{x'}(x)] = \mathbb{E}_{x \sim p}[(\boldsymbol{s}_q(x)k(x,x') + \nabla_x k(x,x')] = \mathbb{E}_x[(\boldsymbol{s}_q^\ell(x) - \boldsymbol{s}_p^\ell(x))k(x,x')]$. In addition,

$$\begin{split} \langle \boldsymbol{f}, \boldsymbol{\beta} \rangle_{\mathcal{H}^d} &= \sum_{\ell=1}^d \langle f_\ell, \ \mathbb{E}_{x \sim p}[(\boldsymbol{s}_q^\ell(x) k(x, \cdot) + \nabla_{x_\ell} k(x, \cdot)] \rangle_{\mathcal{H}} \\ &= \sum_{\ell=1}^d \mathbb{E}_{x \sim p}[(\boldsymbol{s}_q^\ell(x) \langle f_\ell, k(x, \cdot) \rangle_{\mathcal{H}} + \langle f_\ell, \nabla_{x_\ell} k(x, \cdot) \rangle_{\mathcal{H}}] \\ &= \sum_{\ell=1}^d \mathbb{E}_{x \sim p}[(\boldsymbol{s}_q^\ell(x) f_\ell(x) + \nabla_{x_\ell} f_\ell(x)] \\ &= \mathbb{E}_{x \sim p}[\text{trace}(\mathcal{A}_q \boldsymbol{f}(x))], \end{split}$$

where we used the fact that $\nabla_x f(x) = \langle f(\cdot), \nabla_x k(x, \cdot) \rangle_{\mathcal{H}}$; see (Zhou, 2008; Steinwart & Christmann, 2008). The variational form (11) then follows the fact that $||\beta||_{\mathcal{H}^d} = \max_{f \in \mathcal{H}^d} \left\{ \langle f, \beta \rangle_{\mathcal{H}^d}, ...s.t. ||f||_{\mathcal{H}^d} \leq 1 \right\}$.

Finally, the $\beta(\cdot) = \mathbb{E}_{x \sim p}[(s_q^{\ell}(x)k(x,\cdot) + \nabla_x k(x,\cdot)]]$ is in the Stein class of p because $k(x,\cdot)$ and $\nabla_x k(x,\cdot)$ are in the Stein class of p for any fixed x (see the proof of Theorem 3.5).

Proof Proposition 3.4. For any $f \in \mathcal{H}$ with kernel k(x, x'), we have $f = \langle f, k(\cdot, x) \rangle_{\mathcal{H}}$ and $\nabla_x f = \langle f, \nabla_x k(x, \cdot) \rangle_{\mathcal{H}}$. Therefore,

$$\mathbb{E}_{x \sim p}[\mathbf{s}_{p}(x)f(x) + \nabla_{x}f(x)] = \mathbb{E}_{x \sim p}[\mathbf{s}_{p}(x)\langle f, k(x, \cdot)\rangle_{\mathcal{H}} + \langle f, \nabla_{x}k(x, \cdot)\rangle_{\mathcal{H}}]$$

$$= \langle f, \mathbb{E}_{x \sim p}[\mathbf{s}_{p}(x)k(x, \cdot) + \nabla_{x}k(x, \cdot)]\rangle_{\mathcal{H}}$$

$$= \langle f, \mathbb{E}_{x \sim p}[\mathcal{A}_{p}k_{x}(\cdot)]\rangle_{\mathcal{H}}$$

$$= 0.$$

where the last step used the fact that $\mathbb{E}_{x \sim p}[A_p k_x(\cdot)]$ because $k_x(\cdot) = k(\cdot, x)$ is in the Stein class of p for any fixed x. \square

Proof of Theorem 4.1. Applying the standard asymptotic results of *U*-statistics in Serfling (2009, Section 5.5), we just need to check that $\sigma_u^2 \neq 0$ when $p \neq q$ and $\sigma_u^2 = 0$ when p = q.

We first note that we have $\mathbb{E}_{x'\sim p}[u_q(x,x')] = \sum_{\ell} (\mathbf{s}_q^{\ell}(x)\mathbf{g}_{\ell}(x) + \nabla_{x_{\ell}}\mathbf{g}_{\ell}(x))$, where $\mathbf{g}(x) = \mathbb{E}_{x'\sim p}[\mathcal{A}_q k_x(x')]$ and is in the Stein class of p. Therefore, when p=q, we have $g(x)\equiv 0$ by Stein's identity, and hence $\sigma_u^2=0$.

Assume $\sigma_u^2 = 0$ when $p \neq q$, we must have $\mathbb{E}_{x' \sim p}[u_q(x, x')] = c$, where c is a constant. Because $s_q^{\ell}(x) = \nabla_{x_{\ell}} q(x)/q(x)$, we have

$$\sum_{\ell} \nabla_{x_{\ell}}[q(x)\boldsymbol{g}_{\ell}(x)] = q(x) \sum_{\ell} (\boldsymbol{s}_{q}^{\ell}(x)\boldsymbol{g}_{\ell}(x) + \nabla_{x_{\ell}}\boldsymbol{g}_{\ell}(x)) = q(x) \mathbb{E}_{x' \sim p}[u_{q}(x, x')] = cq(x),$$

and therefore $\int \sum_{\ell} \nabla_{x_{\ell}}[q(x)g_{\ell}(x)]dx = c \int q(x)dx = c$. Since g(x) is in the Stein class of q by assumption, we should have c=0 and hence $\mathbb{S}(p,q)=c=0$, which contradicts with $p\neq q$.

Proof of Theorem 5.1. (17) is obtained by applying Cauchy-Schwarz inequality on (6),

$$\begin{split} \mathbb{S}(p,q)^2 &= |\mathbb{E}_{xx'}[(\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x))^\top k(x,x')(\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x))]|^2 \\ &\leq \mathbb{E}_{xx'}[k(x,x')^2] \cdot \mathbb{E}_{x,x'}[[(\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x))^\top (\boldsymbol{s}_q(x') - \boldsymbol{s}_p(x'))]^2] \\ &\leq \mathbb{E}_{xx'}[k(x,x')^2] \cdot \mathbb{E}_{x,x'}[||\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x)||_2^2 \cdot ||\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x)||_2^2] \\ &= \mathbb{E}_{xx'}[k(x,x')^2] \cdot \mathbb{F}(p,q)^2. \end{split}$$

To prove (18), we simply note that (11) is equivalent to

$$\sqrt{\mathbb{S}(p,q)} = \max_{\boldsymbol{f} \in \mathcal{H}^d} \bigg\{ \mathbb{E}_p[(\boldsymbol{s}_q(x) - \boldsymbol{s}_p(x))^\top \boldsymbol{f}(x)] \quad s.t. \quad ||\boldsymbol{f}||_{\mathcal{H}^d} \leq 1 \bigg\}.$$

Taking $f = (s_q - s_p)/||s_q(x) - s_p(x)||_{\mathcal{H}^d}$ then gives (18).

Proof of Proposition 5.2. Denote by $(\mathcal{L}^2(p))^d = \mathcal{L}^2(p) \times \cdots \times \mathcal{L}^2(p)$, note that by the definition of $\mathbb{F}(p,q)$, we have

$$\sqrt{\mathbb{F}(p,q)} = \max_{\mathbf{f} \in (\mathcal{L}^2(p))^d} \left\{ \sum_{\ell=1}^d \mathbb{E}_p[f_\ell(x)(\mathbf{s}_q^\ell(x) - \mathbf{s}_p^\ell(x))] \quad s.t. \quad \mathbb{E}_p[||\mathbf{f}(x)||_2^2] \le 1 \right\}.$$
 (7.20)

Restricting the maximizing to $\mathcal{F}(p)^d$ and applying Lemma 2.3 would give the result.