Random Matrix-Improved Estimation of the Wasserstein Distance between two Centered Gaussian Distributions

Malik Tiomoko¹, Romain Couillet¹,²,*
¹CentraleSupélec, Université ParisSaclay, ²GIPSA-lab, Université Grenoble-Alpes

Abstract—This article proposes a method to consistently estimate functionals \( \frac{1}{p} \sum_{i=1}^{p} f(\lambda_i(C_1C_2)) \) of the eigenvalues of the product of two covariance matrices \( C_1, C_2 \in \mathbb{R}^{p \times p} \) based on the empirical estimates \( \lambda_i(C_1C_2) \) \( (\hat{C}_a = \frac{1}{n_a} \sum_{i=1}^{n_a} x_i^{(a)} x_i^{(a)\text{T}}) \), when the size \( p \) and number \( n_a \) of the (zero mean) samples \( x_i^{(a)} \) are similar. As a corollary, a consistent estimate of the Wasserstein distance (related to the case \( f(t) = \sqrt{t} \)) between centered Gaussian distributions is derived.

The new estimate is shown to largely outperform the classical sample covariance-based “plug-in” estimator. Based on this finding, a practical application to covariance estimation is then devised which demonstrates potentially significant performance gains with respect to state-of-the-art alternatives.

I. INTRODUCTION

Many machine learning and signal processing applications require an adequate framework to compare statistical objects, starting with probability distributions. The Wasserstein distance, initially inspired by Monge [1] and later by Kantorovich [2] in a transport theory analogy, provides a natural notion of dissimilarity for probability measures and finds a wide spectrum of applications in image analysis [3], shape matching [4], computer vision [5], etc.

However, computing the Wasserstein distance is expensive as it requires to minimize a cost function taking the form of an integral over the space of probability measures. Despite recent advances [6], where regularized approximations that reduce this numerical cost are proposed, the latter is still involved in general. Special cases exist for which the Wasserstein distance assumes a closed form, particularly when the underlying distributions are zero-mean Gaussian with covariance matrices \( C_1 \) and \( C_2 \). The closed-form formula however involves the eigenvalues of \( C_1C_2 \) and thus depends on the unknown population covariance matrices \( C_1 \) and \( C_2 \).

Assuming the observation of \( n_1, n_2 \gg p \) samples with covariances \( C_1, C_2 \), respectively, \( C_1C_2 \) is conveniently approximated by its empirical version \( \hat{C}_1\hat{C}_2 \). As we will show, this induces a dramatic estimation bias in practical applications where \( p \) is rather large or, equivalently, \( n_1, n_2 \) rather small, a standard assumption in big data applications.

Based on recent advances in random matrix theory, this article proposes a new consistent estimate for the Wasserstein distance between two centered Gaussian distributions when the dimension \( p \) of the samples is of the same order of magnitude as their numbers \( n_1, n_2 \). This work enters the scope of Mestre’s seminal ideas [7] on the estimation of functionals \( \frac{1}{p} \sum_{i=1}^{p} f(\lambda_i(C)) \) of the eigenvalue distribution of population covariance matrices \( C \), which can be related to the (limiting) eigenvalue distribution of the sample estimates \( C \) via a complex integration trick. We recently extended this work to the estimation of functionals of the eigenvalue distribution of F-matrices in [8], i.e., matrices of the form \( C_1^{-1}C_2 \), and applied to the estimation of the natural geodesic Fisher distance, Bhattacharyya distance, and Rényi/Kullback-Leibler divergences between Gaussian distributions.

Our main contribution is the extension of [7], [8] to functionals \( f \) of the eigenvalues of products \( C_1C_2 \) of population covariance matrices. The Wasserstein distance falls within this scope for \( f(t) = \sqrt{t} \). Unlike [8], where the functionals of interest \( f(t) = t, \log(t), \log^2(t) \) are amenable to explicit evaluations of the complex integrals, the present \( f(t) = \sqrt{t} \) scenario is more technically involved and gives rise to real non-explicit, yet numerically computable, integrals.

In the remainder of the article, Section II introduces the main model and assumptions, Section III provides our key technical result and its corollary to the Wasserstein distance estimation, and a practical application to covariance matrix estimation is finally proposed in Section IV.

Reproducibility. Matlab codes for the various estimators introduced and studied in this article are available at https://github.com/maliktiomoko/RMTWasserstein

II. MODEL AND MAIN OBJECTIVE

For \( a \in \{1, 2\} \), let \( X_a = [x_1^{(a)}, \ldots, x_{n_a}^{(a)}] \) be \( n_a \) independent and identically distributed random vectors with \( x_i^{(a)} = C_2^{1/2} \tilde{x}_i^{(a)} \), where \( \tilde{x}_i^{(a)} \in \mathbb{R}^p \) has zero mean, unit variance and finite fourth order moment entries. This holds in particular for \( x_i^{(a)} \sim \mathcal{N}(0, C_a) \). In order to control the growth rates of \( n_1, n_2, p \), we make the following assumption:

Assumption 1 (Growth Rates). As \( n_a \to \infty \), \( p/n_a \to c_a \in (0, 1) \) and \( \limsup_p \max\{\|C_a^{-1}\|, \|C_a\|\} < \infty \) for \( \|\cdot\| \) the operator norm.

We define the sample covariance estimate \( \hat{C}_a \) of \( C_a \) as

\[
\hat{C}_a = \frac{1}{n_a} X_aX_a^T = \frac{1}{n_a} \sum_{i=1}^{n_a} x_i^{(a)} x_i^{(a)\text{T}}.
\]

The Wasserstein distance \( D_W(C_1, C_2) \) between two zero-mean Gaussian distributions with covariances \( C_1 \) and \( C_2 \), respectively, assumes the form [9, Remark 2.31]:

\[
D_W(C_1, C_2) = \text{tr}(C_1) + \text{tr}(C_2) - 2\text{tr}\left[\left(C_2^{1/2}C_1^{-1}C_2^{1/2}\right)^{1/2}\right].
\]
It is easily shown that, under Assumption \[1\]
\[
\frac{1}{p} \text{tr} \tilde{C}_n - \frac{1}{p} \text{tr} C_n \to 0
\]
almost surely. But estimating \(\frac{1}{p} \text{tr}(C_1^2 C_2 C_2^2)^\frac{1}{2}\) is more involved: this is the focus of the article. Up to a normalization by \(p\), this term can be written under the functional form:
\[
\frac{1}{p} \text{tr}(C_1^2 C_2 C_2^2)^\frac{1}{2} = \frac{1}{p} \sum_{i=1}^n \sqrt{\lambda_i(C_1 C_2)} \equiv D(C_1, C_2; \sqrt{\gamma})
\]
with \(\lambda_i(X)\) the \(i\)-th smallest eigenvalue of \(X\).

Our objective is to estimate the more generic form
\[
D(C_1, C_2; f) \equiv \frac{1}{p} \sum_{i=1}^n f(\lambda_i(C_1 C_2))
\]
for \(f : \mathbb{R} \to \mathbb{R}\) a real function admitting a complex-analytic extension. To this end, we shall relate the eigenvalues \(\lambda_i(C_1 C_2)\) to \(\lambda_i(\hat{C}_1 \hat{C}_2)\) through the Stieltjes transform \((m_\theta(z) \equiv \int \frac{d\theta(x)}{x-z})\) for measure \(\theta\) and \(z \in \mathfrak{C}\) of their associated normalized counting measures
\[
\mu_p = \frac{1}{p} \sum_{i=1}^p \delta_{\lambda_i(\hat{C}_1 \hat{C}_2)}, \quad \nu_p = \frac{1}{p} \sum_{i=1}^p \delta_{\lambda_i(C_1 C_2)}.
\]
In particular, \(m_{\nu_p}(z) = \frac{1}{p} \sum_{i=1}^p \frac{1}{\lambda_i - z} - \lambda_i(\hat{C}_1 \hat{C}_2)\).

With these notations, we are in position to introduce our main results.

III. MAIN RESULTS

The following theorem provides a consistent estimate for the metric \(D(C_1, C_2; f)\) defined in \[4\].

**Theorem 1.** Let \(\Gamma \subset \{ z \in \mathfrak{C}, \text{real}[z] > 0 \}\) be a contour surrounding \(\cup_{p=1}^\infty \text{Supp}(\mu_p)\). Then, under Assumption \[7\]
\[
D(C_1, C_2; f) - \hat{D}(X_1, X_2; f) \overset{a.s.}{\longrightarrow} 0
\]
where
\[
\hat{D}(X_1, X_2; f) = \frac{n_2}{2\pi ip} \int_{\Gamma} f \left( \frac{\varphi_p(z)}{\psi_p(z)} \right) \left[ \frac{\varphi_p'(z)}{\varphi_p(z)} - \frac{\psi_p'(z)}{\psi_p(z)} \right] \psi_p(z)dz
\]
and, recalling \(m_{\mu_p}(z) = \frac{1}{p} \sum_{i=1}^p \frac{1}{\lambda_i - z} - \lambda_i(\hat{C}_1 \hat{C}_2), \varphi_p(z) = 1 - \frac{\pi}{n_1} \log \left| m_{\mu_p}(z) \right|, \psi_p(z) = 1 - \frac{\pi}{n_2} m_{\mu_p}(z)\).

The result of Theorem \[1\] is very similar to [8, Theorem 1] established for functionals of the eigenvalues of \(C_1^{-1} C_2\). The main difference lies in the expression of the function \(\varphi_p(z)\).

**Proof:** The proof of Theorem \[1\] is based on the same approach as for [10, Theorem 1]. One first creates a link between the Stieltjes transform \(m_{\nu_p}\) and \(D(C_1, C_2; f)\) using Cauchy’s integral formula:
\[
\frac{1}{p} \sum_{i=1}^p f(\lambda_i(C_1 C_2)) = \int_{\Gamma} f(t) d\nu_p(t)
\]
\[
= \frac{1}{2\pi i} \int_{\Gamma} \left[ f(t) \Psi_p'(z) - f(z) \Psi_p'(t) \right] dz
\]
\[
= \frac{1}{2\pi i} \int_{\Gamma} f(z) m_{\nu_p}(z) dz
\]
with \(\Gamma\) a contour surrounding the support \(\text{supp}(\nu_p)\) of \(\nu_p\). To relate the unknown \(m_{\nu_p}\) to the observable \(m_{\mu_p}\), we proceed as follows. By first conditioning on \(\hat{C}_1, \hat{C}_1 \hat{C}_2 \hat{C}_2^2\) seen as a sample covariance matrix for the samples \(C_1^2 C_2^2 \hat{z}_{(2)}\), for which [11] allows one to relate \(m_{\nu_p}\) to the Stieltjes transform of the eigenvalue distribution \(\zeta_p\) of \(C_2^2 \hat{C}_1 \hat{C}_2\). The latter is yet another sample covariance matrix for the samples \(C_2^2 \hat{C}_1 \hat{C}_2 \hat{C}_2^2\); exploiting [11] again creates the connection from \(m_{\zeta_p}\) to \(m_{\nu_p}\). This entails the two equations:
\[
m_{\mu_p}(z) = \varphi_p(z) m_{\zeta_p}(z) + o_p(1)
\]
\[
m_{\nu_p}(z) \sqrt{\Psi_p(z)} = m_{\zeta_p}(z) + o_p(1).
\]
where \(\Psi_p(z) = 1 - \frac{p}{n_2} - \frac{p}{n_z} m_{\zeta_p}(z)\). Successively plugging \[5\]–\[6\] into \[4\] by means of two successive appropriate changes of variables, we obtain Theorem \[1\].

Theorem \[1\] takes the form of a complex integral which, for generic choices of \(f\), needs be numerically evaluated. In the specific case of present interest where \(f(z) = \sqrt{z}\), this complex integral can be evaluated as follows.

**Theorem 2.** Let \(\lambda_1, \ldots, \lambda_p \geq 0\), with \(\lambda_i(\hat{C}_1 \hat{C}_2)\), and define \(\xi_j^{(1)} = \{ \xi_j \}_{j=1}^p\) the (increasing) eigenvalues of \(\Lambda - \frac{1}{n_1} \Lambda \Lambda^T\) and \(\Lambda - \frac{1}{n_2} \Lambda \Lambda^T\), respectively, where \(\lambda = (\lambda_1, \ldots, \lambda_p)^T\), \(\Lambda = \text{diag}(\lambda)\) and \(\sqrt{\cdot}\) is understood entry wise. Then, under Assumption \[7\]
\[
D(C_1, C_2; \sqrt{\gamma}) - \hat{D}(X_1, X_2; \sqrt{\gamma}) \overset{a.s.}{\longrightarrow} 0
\]
where, if \(n_1 \neq n_2\),
\[
\hat{D}(X_1, X_2; \sqrt{\gamma}) = 2 \sqrt{n_1} \sqrt{\frac{1}{n_2} \sum_{j=1}^p \frac{1}{\sqrt{x_j}}}
\]
\[
+ 2 \sqrt{n_2} \sum_{p=1}^p \int_{\xi_j^{(1)}} \frac{\varphi_p(x)}{\psi_p(x)} \psi_p(x) dx
\]
with \(\varphi_p, \psi_p\) defined in Theorem \[7\] and, if \(n_1 = n_2\),
\[
\hat{D}(X_1, X_2; \sqrt{\gamma}) = 2 \frac{n_1}{p} \sum_{j=1}^p \left( \sqrt{\lambda_j} - \sqrt{\xi_j} \right).
\]

While still assuming an integral form (when \(n_1 \neq n_2\)), this formulation no longer requires the arbitrary choice of a contour \(\Gamma\) and significantly reduces the computational time to estimate \(D(C_1, C_2; \sqrt{\gamma})\). For \(n_1 = n_2\), a case of utmost practical interest, the expression is completely explicit and computationally only requires to evaluate the eigenvalues \(\xi_j\) of \(\Lambda - \frac{1}{n_1} \Lambda \Lambda^T\). The latter being a (negative definite) rank-1 perturbation of \(\Lambda\), by Weyl’s interlacing lemma [12], the \(\xi_j\)’s are interlaced with the \(\lambda_j\)’s as
\[
\xi_1 \leq \lambda_1 \leq \xi_2 \leq \ldots \leq \xi_p \leq \lambda_p.
\]
As the \(\lambda_j\)’s are of order \(O(1)\) with respect to \(p\), \(|\lambda_j - \xi_j| \leq |\lambda_j - \lambda_j| = O(p^{-1})\), therefore explaining why the expression of \(\hat{D}(X_1, X_2; \sqrt{\gamma})\) is of order \(O(1)\).
Proof: The $\xi_i$ and $\eta_i$, as defined in the theorem statement, are the respective zeros of the rational functions $1 - \frac{p}{n_1} - \frac{p}{n_2}z\tilde{m}_p(z)$ and $1 - \frac{p}{n_2} - \frac{p}{n_2}z\tilde{m}_p(z)$ (see [10, Appendix B]). Thus, $\varphi_p$ and $\psi_p$ can be expressed under the rational form:

$$\varphi_p(z) = \sum_{j=1}^{p} \frac{z - \lambda_j}{\prod_{i=1}^{P} (z - \lambda_i)} \psi_p(z) = \prod_{j=1}^{p} \frac{z - \xi_j}{\prod_{i=1}^{P} (z - \xi_i)} \psi_p(z),$$

Evaluating the estimate from Theorem 1 for $f(z) = \sqrt{z}$ then requires to evaluate a complex integral involving rational functions and square roots of rational functions. Since the complex square root is multivalued, a careful control of "branch-cuts" is required. To perform this calculus, we deform the integration contour $\Gamma$ of Theorem 1 into $\Gamma$ as per Figure 1. In the case $n_1 \neq n_2$, the closed null-integral contour $\Gamma$ (blue in Figure 1) is the sum of the sought-for integral over $\Gamma$ and of four external contours:

1) Integrals over $\epsilon$-radius circles around $\xi_i$: those are null in the limit $\epsilon \rightarrow 0$, as confirmed by a change of variable $z = \xi_i + e^{i\epsilon}$ which allows one to bound the integrand;
2) Integrals over the real axis (in the $\epsilon \rightarrow 0$ limit):

$$A_2 = \frac{n_2}{\pi p} \sum_{j=1}^{p} \int_{\xi_j + \epsilon}^{\xi_j - \epsilon} \sqrt{\varphi_p \psi_p}(x) \left[ \frac{2 \psi_p'(z)}{\psi_p(z)} \right] \frac{d}{dx} \left( \varphi_p(x) \psi_p(x) \right) dx$$

$$A_2 = \frac{2n_2}{\pi p} \sum_{j=1}^{p} \int_{\xi_j}^{\xi_j + \epsilon} \sqrt{\varphi_p \psi_p}(x) \left[ \frac{d}{dx} \left( \varphi_p(x) \psi_p(x) \right) \right] dx$$

$$A_2 = \frac{2n_2}{\pi p} \sum_{j=1}^{p} \int_{\xi_j}^{\xi_j + \epsilon} \sqrt{\varphi_p \psi_p}(x) \psi_p'(x) dx$$

$$A_2 = \frac{2n_2}{\pi p} \sum_{j=1}^{p} \int_{\xi_j}^{\xi_j + \epsilon} \frac{1}{\sqrt{\varphi_p \psi_p}(x)} \left( \eta_j \right)$$

3) Integrals over the $\epsilon$-radius circles around $\eta_j$, with $\epsilon \rightarrow 0$

$$A_2 = \frac{n_2}{\pi p} \sum_{j=1}^{p} \int_{\xi_j}^{\xi_j + \epsilon} \sqrt{\varphi_p \psi_p}(x) \left[ \frac{1}{\psi_p(x)} \right] \left( \eta_j \right)$$

which thus compensates the last ($\epsilon$-diverging term) in $A_2$.

4) Residues in the $\lambda_j$ poles

$$A_4 = \frac{2n_2}{p} \lim_{z \rightarrow \lambda_j} \sum_{j=1}^{p} \sqrt{(\varphi_p \psi_p)(z)} = \frac{2n_2}{p} \sum_{j=1}^{p} \sqrt{\lambda_j}.$$ 

Putting these terms together entails the result of the theorem for the case where $n_1 \neq n_2$. For $n_1 = n_2$, it suffices to take the limit of the expression as $\xi_j \rightarrow \eta_j$. This yields:

$$\hat{D}(X_1, X_2; \sqrt{\cdot}) = \frac{2n_1}{p} \sum_{j=1}^{p} \sqrt{x_j}$$

$$\hat{D}(X_1, X_2; \sqrt{\cdot}) = \frac{2n_1}{p} \sum_{j=1}^{p} \sqrt{x_j}$$

where $\Gamma_{\xi_j}$ is an $\epsilon$-radius circular contour around $\xi_j$. The second equality is obtained by deforming the real integral in the complex plane (see [13] for complex analysis details). The result unfolds by letting $z = \xi_i + e^{i\theta}$.

Consequently, we obtain the following $n, p$-consistent estimate for the Wasserstein distance $D_W(C_1, C_2)$ of (1).

**Corollary 1** (Consistent Estimate of $D_W(C_1, C_2)$). Under Assumption 1

$$\frac{1}{p} D_W(C_1, C_2) = \frac{1}{p} \text{tr}(\hat{C}_1 + \hat{C}_2) - 2\hat{D}(X_1, X_2; \sqrt{\cdot}) \xrightarrow{a.s.} 0$$

for $\hat{D}(X_1, X_2; \sqrt{\cdot})$ given by Theorem 2.

**Remark 1** (Estimation of $\|C_1 - C_2\|_F^2$). The Frobenius distance between two covariance matrices also falls under the scope of the present article for the function $f(z) = z$.

$$D_F(C_1, C_2) = \|C_1 - C_2\|_F^2 = \text{tr}(C_1^2 + C_2^2) - 2\text{tr}(C_1C_2)$$

Then under Assumption 1 and along with the fact that $\frac{1}{p} \text{tr}C_2^2$ can be estimated consistently from $\frac{1}{p} \text{tr}\hat{C}_2^2 = \frac{1}{n_1} \left( \frac{1}{p} \text{tr}\hat{C}_1^2 \right)^2$,

$$\frac{1}{p} D_F(C_1, C_2) = \frac{1}{p} \text{tr}\hat{C}_2^2 - \frac{1}{n_1} \left( \frac{1}{p} \text{tr}\hat{C}_1 \right)^2$$

$$\frac{1}{p} D_F(C_1, C_2) = \frac{1}{p} \text{tr}\hat{C}_2^2 - \frac{1}{n_1} \left( \frac{1}{p} \text{tr}\hat{C}_1 \right)^2$$

In this case, $\hat{D}(X_1, X_2; \cdot)$ assumes the simple expression

$$\hat{D}(X_1, X_2; \cdot) = \frac{1}{p} \sum_{j=1}^{p} \lambda_j = \frac{1}{p} \text{tr}\hat{C}_1 \hat{C}_2$$

which follows from $\frac{1}{p} \text{tr}\hat{C}_1 \hat{C}_2 = \frac{1}{p} \text{tr}C_1C_2 \xrightarrow{a.s.} 0$ (by elementary probability arguments) or equivalently from a residue calculus based on Theorem 2 for $f(z) = z$. 

IV. SIMULATIONS AND APPLICATIONS

In this section, we first corroborate our theoretical findings by comparing the results of our proposed estimator on synthetic Gaussian data. We then provide an application of our results to improved covariance matrix estimation based on few samples.

A. Confirmation of our results on synthetic data

We here compare the classical plug-in estimate of the Wasserstein distance (that is (1) with $\bar{C}_a$, $a = 1, 2$) with our proposed estimate in Corollary [4]. Table I lists the results obtained for Toeplitz matrices $C_1, C_2$ estimated based on various values of $p, n_1, n_2$. While our proposed estimator is designed under a large $p, n_1, n_2$ assumption (as per Assumption [4]), it achieves competitive performances even for small values of $p$, corroborating our findings in [8] for other classes of covariance matrix distances.

<table>
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<th>$p$</th>
<th>$D_W(C_1, C_2)$ Classical</th>
<th>Proposed</th>
</tr>
</thead>
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<tr>
<td>2</td>
<td>0.0110 0.0127 0.0120</td>
<td></td>
</tr>
<tr>
<td>4</td>
<td>0.0175 0.0198 0.0183</td>
<td></td>
</tr>
<tr>
<td>8</td>
<td>0.0208 0.0232 0.0206</td>
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</tr>
<tr>
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<td></td>
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<tr>
<td>32</td>
<td>0.0233 0.0339 0.0234</td>
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<tr>
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<tr>
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<td>0.0239 0.0667 0.0244</td>
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<tr>
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</tr>
<tr>
<td>512</td>
<td>0.0241 0.1951 0.0245</td>
<td></td>
</tr>
</tbody>
</table>

(error < 5%) (error > 50%) (error > 100%) (error > 300%)

Table I

ESTIMATORS OF THE WASSERSTEIN DISTANCE BETWEEN $C_1$ AND $C_2$

with $|C_1|_{ij} = 2^{i+j}, |C_2|_{ij} = 4^{i+j}, \chi^2/2 \sim N(0, C_a); n_1 = 1024$

and $n_2 = 2048$ for different $p$. Averaged over 100 trials.

B. Application to covariance matrix estimation

As a concrete application, Theorem [4] may be used to improve the actual estimation of covariance matrices under a small number $n \sim p$ of sample data, as similarly performed in [14] for other covariance matrix distances.

The idea is as follows: we first particularize Theorem [4] and Theorem [2] to the case where one of the covariance matrices, say $C_1$, is known, by taking $c = 0$ (i.e., $n_1 \to \infty$ for all fixed $p$). This gives access to estimates for $D_W(M, C_2; \sqrt{\gamma})$ for all deterministic positive definite matrices $M$. We then minimize this estimated distance over $M$ in order to estimate $C_2$ by means of a gradient descent ascent.

For $C_1$ known, we redefine $\mu_p = \frac{1}{p} \sum_{i=1}^{p} \delta_{\lambda_i(C_1, C_2)}$ and obtain, as a corollary of Theorem [1], Theorem 3.

Theorem 3. Let $\Gamma \subset \{ z \in \mathbb{C}, \text{real}[z] > 0 \}$ a contour surrounding $\cup_{p=1}^{\infty} \supp(\mu_p)$. Then,

$$D(C_1, C_2; f) = -\frac{1}{2\pi i c_2} \int_{\Gamma} F(-\mu_p(z)) \frac{dz}{\bar{z}} \xrightarrow{a.s.} 0$$

with $m_{\mu_p}(z) = \frac{p}{n_z} m_{\psi_p}(z) + \frac{n_z}{n_z} z$ and $F^*(z) = f\left(\frac{1}{z}\right)$.

Proof: For $C_1$ known ($c_1 \to 0$), $\varphi_p(z) = z$, and the estimator of Theorem [1] yields:

$$\tilde{D}(X_1, X_2; f) = \frac{1}{2\pi} \int_{\Gamma} \int_{\Gamma} f\left(\frac{z}{\psi_p(z)}\right) \psi_p^2(z) \frac{dz}{c_2}.$$
(precisely, $v_p = \frac{1}{2}(\delta_1 + \delta_3 + \delta_4 + \delta_5)$) and various estimators of $C$: the sample covariance matrix (SCM), the state-of-the-art “non-linear shrinkage” estimators QuEST1 [16] (based on a Frobenius distance minimization) and QuEST2 [17] (based on a Stein loss minimization), and the result of the gradient descent approach proposed in this section. For fair comparison, the iterative QuEST1, QuEST2 and our proposed method are all initialized at $M_0$ the linear shrinkage estimator from [18]. Note that our proposed choice of $C$ is particularly suited to mimic an “optimal transport” problem of displacing the eigenvalues of $M_0$ to the discrete four positions of the eigenvalues of $C$.

In addition to the computational simplicity of our gradient-descent approach with respect to the QuEST estimators (see the numerical method details in [19]), the figure demonstrates significant gains brought by our proposed approach for large values of $p/n$, where the SCM particularly fails.

V. CONCLUDING REMARKS

Interestingly, while the Fisher distance or Kullbach-Liebler divergence, which depend on logarithms of $p/n$ values of $\nu_p$, are not satisfying if $p < n$, $n_2$ from Assumption [1] are fundamental to our proofs. Precisely, the variable changes exploited in the proof of Theorem [1] to reach a contour $\Gamma_c$ correctly surrounding $supp(\nu_p)$ from a contour $\Gamma$ surrounding $supp(\mu_p)$ are not satisfying if $c_1 > 1$ or $c_2 > 1$. These surprising difficulties need clarification.

Another point of interest lies in the comparative advantage of exploiting a particular covariance matrix distance in specific scenarios. For instance, it may seem that ill-conditioned matrices should be more tolerated by Wasserstein distance estimators than by Fisher distance estimators. Yet, this aspect is not obvious in our proofs and also deserves more insights.

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