An Adaptive Quasi-Newton Algorithm for Eigensubspace Estimation

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Abstract—In this paper, we derive and discuss a new adaptive quasi-Newton eigen-estimation algorithm and compare it with the RLS-type adaptive algorithms and the quasi-Newton algorithm proposed by Mathew et al. through experiments with stationary and nonstationary data.

Index Terms—Adaptive estimation, deflation, eigen-subspace, principal component.

I. INTRODUCTION

It is well known that the gradient-descent-based adaptive principal component analysis (PCA) algorithms are slow to converge. Bannour et al. [3] and Yang [4] derive RLS-type adaptive eigendecomposition algorithms from neural network and optimization frameworks, respectively. Strictly speaking, they are gradient descent algorithms with time-varying, self-tuning step sizes and are still slow to converge. In [1], Mathew et al. present a rapidly convergent quasi-Newton adaptive algorithm derived from a cost function based on the penalty function method of optimization. This algorithm sequentially estimates the eigenvector corresponding to the smallest eigenvalue up to the eigenvector corresponding to the largest eigenvalue. In many applications, however, we are only interested in estimating the principal eigenvectors. Reversing the estimation order of this algorithm (i.e., from the “largest” eigenvector to the “smallest” eigenvector sequentially) is not straightforward if we employ the Hessian approximation method proposed in [1] directly in the quasi-Newton algorithm.

In this paper, using a different method to approximate Hessian, we derive a novel quasi-Newton-based adaptive eigensubspace estimation algorithm that first estimates the principal eigenvector and then estimates the minor eigenvectors sequentially. Moreover, in our algorithm, the choice of the penalty coefficient does not need any a priori information of the data covariance matrix while in the Mathew et al. algorithm, this choice needs a priori information of data. In next section, we study the landscape of the cost function.

Section III proposes the new Quasi-Newton algorithm and gives the convergence analysis. The extensive experiments with stationary and nonstationary data are given in Section IV. Section V contains concluding remarks.

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Fig. 1. Convergence of the first four principal eigenvectors with the (a) QN1, (b) QN2, (c) PASTd, and (d) RLS algorithms.

In order to compute the inverse of Hessian efficiently, we define

\[ Q_k = -R_k + 2\mu \left( w_k^T w_k - 1 \right) I_N. \]  

Therefore, the inverse of the Hessian becomes

\[ H_k^{-1} = \left( 2Q_k + 8I w_k w_k^T \right)^{-1} \]

\[ = \frac{1}{2} \left[ Q_k^{-1} - Q_k^{-1} w_k w_k^T Q_k^{-1} \right]. \]

Assuming \( w_k \) is close to the solution \( w^* \) at time instant \( k \), the inverse of \( Q_k \) can be approximated as

\[ Q_k^{-1} \approx \frac{1}{2\mu} \left( w_k^T w_k - 1 \right) \left[ I_N + \frac{R_k}{2\mu} \left( w_k^T w_k - 1 \right) \right]. \]

After simplification, we get the following adaptive quasi-Newton algorithm:

\[ w_{k+1} = L_k w_k, \]

\[ L_k = \frac{w_k^T w_k}{\tau_k} \left[ 2\mu \left( w_k^T w_k - 1 \right) I_N + R_k \right] \]

\[ \tau_k = \mu \left( w_k^T w_k - 1 \right)^2 + 2\mu w_k^T w_k \left( w_k^T w_k - 1 \right) \]

\[ + w_k^T R_k w_k \]

\[ R_k = \begin{cases} R_k \quad &i = 1 \\ R_k^{-1} - w_k w_k^T R_k &2 \leq i \leq p \end{cases} \]
Fig. 2. Convergence performance of QN1 and QN2 algorithm to compute the first four eigenvectors under nonstationary data sequences with $\mu = 15$ and $\theta = 0.9955$. (a) QN1. (b) QN2.

Fig. 3. Convergence performance of the QN1 and QN2 algorithms to track the first eigenvector under nonstationary data sequences with $\mu = 15$ and $\theta = 0.9955$.

The approximate Hessian is obtained by dropping one term of the Hessian in [1]. We reached an approximate Hessian through geometric progression approximation. In such a geometric progression approximation, more terms can be used to keep the approximate Hessian closer to the exact Hessian.

We have the following result for the convergence of the above algorithm.

**Assumption A2):** Each $R_k$ is bounded with probability one (w.p. 1), symmetric, real, non-negative definite, and $\lim_{k \to \infty} E[R_k] = R$ w.p. 1.

**Theorem 3:** Let A1) and A2) hold. For any $\mu > 0$ and initial values $w_0$ sufficiently close to the desired eigenvectors, if process $\{w_k\}$ is generated by the quasi-Newton algorithm (9)–(13), then a) $\lim_{k \to \infty} w_k = \phi_i$; and b) $\lim_{k \to \infty} \tau_k = \lambda_i$ for $i = 1, 2, \ldots, p$, where $\lambda_i$ is the $i$th eigenvalue of $R$, and $\phi_i$ is the corresponding eigenvector of $R$.

**Proof:** a) can be completed by first proving $\lim_{k \to \infty} w_k = \rho \rho_i$ and then proving $\rho_i = \pm 1$; b) immediately follows a).

## IV. SIMULATIONS

In this section, we compare our adaptive PCA algorithm (QN1) with the following:

1) the Mathew et al. quasi-Newton algorithm (QN2);
2) the Yang RLS algorithm (PASTd);
3) the Bannour et al. RLS algorithm (RLS).

We demonstrate these comparisons with two different simulation methods. In both methods, high-dimensional data sequences are generated.

### A. Method 1

In this simulation method, we use the first covariance matrix from [6] multiplied by 2. We generate samples of 10-dimensional Gaussian data with mean zero and covariance $R$. The eigenvalues of the this covariance matrix $R$ in decreasing order are

$$
11.7096 \quad 5.5644 \quad 3.4175 \quad 2.0589 \quad 0.7873 \\
0.5878 \quad 0.1743 \quad 0.1423 \quad 0.1213 \quad 0.1007.
$$

Clearly, the first four eigenvalues are significant, and corresponding eigenvectors are important. We use **direction cosine** (DC) to measure the convergence and accuracy of the adaptive algorithms. It is defined as

$$
\text{DC}_k^i = \frac{\langle w_k^i, \phi_i \rangle}{\|w_k^i\| \|\phi_i\|}, \quad i = 1, 2, \ldots, p \quad (p \leq N)
$$

where $w_k^i$ is the estimated $i$th eigenvector of $R_k$ corresponding to the $i$th eigenvalue at the $k$th update, and $\phi_i$ is the actual eigenvector of $R$ computed by a conventional numerical analysis method. Clearly, $\lim_{k \to \infty} \text{DC}_k^i = 1$ if $\lim_{k \to \infty} w_k^i = \phi_i$.

We start our algorithm with $w_0^i = 0.8 \cdot \mathbf{i}_i$ for $i = 1, 2, \ldots, p$ $(p \leq N)$, where $\mathbf{i}_i$ denotes the $i$th column of $N \times N$ identity matrix $I_N$. For other adaptive algorithms, the initial values are QN2: $w_0^i = 0.8 \cdot \mathbf{i}_i$; $R_0^{-1} = 100 \cdot I_N$; PASTd: $w_0^i = 0.8 \cdot \mathbf{i}_i$; $P_0^i = 100 \cdot \mathbf{i}_i$.
Suppose we do not know any *a priori* information on data sequence, we simply choose $\mu = 1$ to compute the first four principal eigenvectors for all of the adaptive algorithms. Fig. 1 shows the convergence of these adaptive algorithms.

The simulation results demonstrate that QN1 and QN2 algorithms offer faster convergence rates than the PASTd and RLS algorithms. We also see that the QN2 algorithm does not estimate the first three principal eigenvectors when $\mu = 1$. In the QN2 algorithm, $\mu$ has to be larger than $\frac{\lambda_{\text{max}}}{2}$ or $\mu > \frac{\text{trace}(R)}{2}$. It turns out that the QN2 algorithm requires a $\mu > \frac{\lambda_{\text{max}}}{2} \approx 5.8908$ to estimate all eigenvectors. Clearly, in the QN2 algorithm, an ad hoc selection of $\mu$ is risky for applications where we do not have *a priori* information on the data covariance matrix $R$ or where there are jumps in eigenvalues of $R$ in the time-varying data sequences. However, when a large suitable penalty coefficient $\mu$ is chosen, the QN2 algorithm exhibits very similar convergence performance as the QN1 algorithm for stationary signals.

From experimental results, we also find that the QN1 and QN2 algorithms have good tracking properties, even for minor eigenvectors, whereas PASTd and RLS do not have this performance. The results are omitted here.

In order to demonstrate the tracking abilities of the two quasi-Newton algorithms (QN1 and QN2) under a nonstationary environment, we provide a time-varying scenario in which the covariance matrix of data sequences change suddenly. The first
500 data samples of the data sequence have the covariance \( R \), as described above. After 500 samples, we abruptly change to covariance \( R' \), which is obtained from the second matrix in [6] multiplied by 30. Accordingly, the eigenvalues change from
\[
\begin{align*}
11.7996 & \quad 5.5644 & \quad 3.4175 & \quad 2.0589 & \quad 0.7873 \\
0.5878 & \quad 0.1743 & \quad 0.1423 & \quad 0.1213 & \quad 0.1007
\end{align*}
\]
to
\[
\begin{align*}
178.8639 & \quad 101.2184 & \quad 86.0674 & \quad 65.3080 & \quad 22.4001 \\
18.3543 & \quad 11.3537 & \quad 7.6971 & \quad 1.4791 & \quad 2.2881.
\end{align*}
\]
As discussed in Section III, we use forgetting factor \( \theta \) to reduce the effect of old data and enhance the tracking ability in the nonstationary case. Therefore, we use iterative equation (5) to update \( R_k \) in the QN1 algorithm. Accordingly, we use
\[
R_k^{-1} = \frac{k}{(k-1)\theta} \left[ R_{k-1}^{-1} - \frac{R_{k-1}^{-1} x_k x_k^T R_{k-1}^{-1}}{(k-1)\theta + x_k^T R_{k-1}^{-1} x_k} \right]
\]
to update \( R_k^{-1} \) in the QN2 algorithm. Fig. 2 shows the simulation results for both quasi-Newton adaptive algorithms when \( \mu = 15 \) and \( \theta = 0.9955 \). We observe that the QN2 algorithm does not track the first principal eigenvector, whereas the QN1 algorithm tracks all of eigenvectors quite well. Fig. 3 shows the results of QN1 and QN2 algorithms to track the first principal eigenvector when \( \mu = 1000 \) and \( \theta = 0.9955 \). It demonstrates that the QN1 algorithm converges faster and tracks the changes in data much better than the QN2 algorithm. It is clear that QN1 is superior to QN2 under a nonstationary environment.

B. Method 2

In this simulation, we follow the experimental setup in [1], where the data \( x_k \) is generated as
\[
x_k = A \sin(0.4\pi k) + A \sin(0.48\pi k + \vartheta) + v_k \quad k = 1, 2, \ldots
\]
The initial phase \( \vartheta \) is uniformly distributed in \([\pi, \pi]\), and \( v_k \) is white noise with zero mean and unit variance. Scalar \( A \) determines the desired signal-to-noise ratio (SNR) as \( 10 \log(A^2/2) \). Here, we choose a window of length \( N \) and the dimension of the covariance as \( 6 \). We estimate the first \( p = 4 \) principal eigenvectors.

The initial values for each algorithm are the same as those used in Method 1. We run 100 Monte Carlo simulations for each algorithm. We only compare the QN1 algorithm with the RLS-type algorithms since the QN2 algorithm has very similar tracking performance as QN1 under stationary environment for a large value of penalty coefficient. Nonstationary experiments show that the QN1 algorithm has superior tracking performance than the QN2 algorithm. Further consideration should be given to the computational complexity of the algorithms since the QN1 and QN2 algorithms have the computational complexity of \( O(pN^2) \), whereas the PASTd and RLS algorithms offer \( O(pN) \) complexity.

APPENDIX

This Appendix proves Theorem 2. From the objective function, we define the following energy functions:
\[
E_i(w^1, \ldots, w^i) = -w^{iT} R w^i + \sum_{j=1}^{i-1} w^j \cdot w^j \cdot w^j \cdot R w^j + \mu \left( w^{iT} w^i - 1 \right)^2
\]
for \( i = 1, \ldots, p \) \((p \leq N)\). Let us assume that \( \rho(\vartheta) = 0 \) for some \( i \leq p \), i.e.,
\[
W = [ \rho(\vartheta(1)), \ldots, \rho(i-1)\vartheta(i-1), 0, \rho(i+1)\vartheta(i+1), \ldots, \rho(p)\vartheta(p) ].
\]
Next, we perturb the \( i \)th column of \( W \) by \( \delta \vartheta(\vartheta) \). We observe that
\[
E_i (\rho(1)\vartheta(1), \ldots, \rho(i-1)\vartheta(i-1), \delta \vartheta(\vartheta)), \ldots, \rho(i-1)\vartheta(i-1), 0) = -\delta \vartheta(\vartheta) \lambda(i) + \mu(\delta \vartheta(\vartheta)^2 - 1)^2 - \mu.
\]
Clearly, the energy function decrease for some \( \delta \vartheta(\vartheta) < \sqrt{2/(\lambda(i)/\mu)} \).

We next prove that if \( \vartheta(\vartheta) \neq \vartheta(i) \) for some \( i = 1, \ldots, p \) then the critical points \( W \) are unstable equilibrium points of (15). We obtain the Hessian of \( E_i \) with respect to \( w^i \) as
\[
\nabla_{w^i}^2 E_i(w^1, \ldots, w^i) = -2R + \sum_{j=1}^{i-1} \sum_{j=1}^{i-1} R w^j w^j \cdot R w^j + 4\mu \left( w^{iT} w^i - 1 \right) I + 8\mu w^i w^i \cdot R^T.
\]
We prove the result by induction. We note that \( \nabla_{w^i}^2 E_i(\rho(1)\vartheta) = -2R + 4\mu(\rho^2 - 1)I + 8\mu \rho \vartheta(\vartheta)^2 \vartheta(\vartheta)^2 \). The eigenvectors of \( \nabla_{w^i}^2 E_i(\rho(1)\vartheta) \) are \( \vartheta(1), \ldots, \vartheta(N) \). The eigenvalues are \( 8\mu \rho^2 \) for \( \vartheta(1) \) and \( 2\lambda(i) - \lambda(i) \) for \( \vartheta(i) \), \((r > 1)\). Clearly, \( \nabla_{w^i}^2 E_i(\rho(1)\vartheta) \) is positive definite. On the other hand, we observe that
\[
\nabla_{w^i}^2 E_i(\rho_r \vartheta_r) = -2R + 4\mu(\rho_r^2 - 1)I + 8\mu \rho_r^2 \vartheta_r(\vartheta_r)^2 \quad (r > 1)
\]
We see that the eigenvectors of \( \nabla_{w^i}^2 E_i(\rho_r \vartheta_r) \) are \( \vartheta(1), \ldots, \vartheta(N) \). However, the eigenvalues are \( -2(\lambda(i) - \lambda(i)) \) for \( \vartheta(1) \) and \( 8\mu \rho_r^2 \) for \( \vartheta(i) \), \((r > 1)\). Clearly, \( \nabla_{w^i}^2 E_i(\rho_r \vartheta_r) \) for \( r > 1 \) is an indefinite matrix. Thus, \( \rho(1)\vartheta(1) \) is a stable equilibrium point for the energy function \( E_i \), whereas \( \rho(1)\vartheta(1) \) for \( r > 1 \) are unstable equilibrium points of \( E_i \).
We next assume that \((\rho_1 \phi_1, \ldots, \rho_{i-1} \phi_{i-1}, \rho_i \phi_i)\) are the stable equilibrium points for the energy function \(E_{i-1}\). Then, for \(\phi(i) = \phi_i\), we have from (15)

\[
\nabla^2_{\omega_i} E_i(\rho_1 \phi_1, \ldots, \rho_{i-1} \phi_{i-1}, \rho_i \phi_i) = -2R + 4\mu \left(\rho_i^2 - 1\right) I + 8\mu \rho_i^2 \phi_i \phi_i^T + 2 \sum_{j=1}^{i-1} \rho_j^2 \phi_j \phi_j^T.
\]

The eigenvectors of \(\nabla^2_{\omega_i} E_i(\rho_1 \phi_1, \ldots, \rho_{i-1} \phi_{i-1}, \rho_i \phi_i)\) are \(\phi_1, \ldots, \phi_N\). The eigenvalues are \(2\lambda_i + \lambda_j^2 / \mu\) for \(\phi_i\), and \(2(\lambda_i - \lambda_j)\) for \(\phi_j\) (\(r > i\)).

Thus, \(\nabla^2_{\omega_i} E_i(\rho_1 \phi_1, \ldots, \rho_{i-1} \phi_{i-1}, \rho_i \phi_i)\) is positive definite.

On the other hand, if \(\phi(i) = \phi_i(r > i)\) then

\[
\nabla^2_{\omega_i} E_i(\rho_1 \phi_1, \ldots, \rho_{i-1} \phi_{i-1}, \rho_i \phi_i) = -2R + 4\mu (\rho_i^2 - 1) I + 8\mu \rho_i^2 \phi_i \phi_i^T + 2 \sum_{j=1}^{i-1} \rho_j^2 \phi_j \phi_j^T,
\]

The eigenvalues of \(\nabla^2_{\omega_i} E_i(\rho_1 \phi_1, \ldots, \rho_{i-1} \phi_{i-1}, \rho_i \phi_i)\) are \(-2(\lambda_i - \lambda_j)\) for \(\phi_i\), and \(-8\mu \rho_i^2\) for \(\phi_j\) (\(r > i\)). Thus, \(\nabla^2_{\omega_i} E_i(\rho_1 \phi_1, \ldots, \rho_{i-1} \phi_{i-1}, \rho_i \phi_i)\) is an indefinite matrix. We therefore see that \((\rho_1 \phi_1, \ldots, \rho_i \phi_i)\) are the stable equilibrium points of \(E_i\).

We conclude that \(W = [\rho_1 \phi_1 \rho_2 \phi_2 \cdots \rho_p \phi_p]\), where \(\rho_i = \pm \sqrt{1 + \lambda_i^2 / 2\mu}\) are stable equilibrium points of the energy functions (15), whereas \(W = [\rho_1 \phi_1(1) \rho_2 \phi_1(2) \cdots \rho_p \phi_1(p)]\), where \(\rho_i = 0\) for \(i > p\) or \(\phi_1(i) \neq \phi_i\) are unstable equilibrium points.

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