

Multiclass linear discriminant analysis with ultrahigh-dimensional features

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Abstract

Within the framework of Fisher's discriminant analysis, we propose a multiclass classification method which embeds variable screening for ultrahigh-dimensional predictors. Leveraging interfeature correlations, we show that the proposed linear classifier recovers informative features with probability tending to one and can asymptotically achieve a zero misclassification rate. We evaluate the finite sample performance of the method via extensive simulations and use this method to classify posttransplantation rejection types based on patients' gene expressions.

KEYWORDS

Fisher's multiclass discriminant analysis, jointly informative features, marginally informative features, multivariate screening, ultrahigh-dimensional classification

1 | INTRODUCTION

Ultrahigh-dimensional data, wherein the number of features p is in the exponential order of the sample size n , have now been routinely collected. For example, in the motivating kidney transplant study (Fletcher *et al.*, 2004), 62 posttransplant kidney tissue samples have been assayed on 12 625 genes. Distinguishing four types of tissues, namely, those from normal donors (C), well-functioning kidneys (TX), kidneys with acute rejection (AR), and kidneys with acute dysfunction but no rejection (NR), based on their molecular biomarkers is important in balancing the need for immunosuppression to prevent rejection and in minimizing drug-induced toxicities.

Linear discriminant analysis (LDA) is a widely used classification method with ready implementability and close relationships with many modern machine learning techniques (Dorfer *et al.*, 2015; Cai *et al.*, 2018; Gorban *et al.*, 2018). In high-dimensional settings, LDA using all features leads to poor results (Fan and Fan, 2008), and high-dimensional LDA is often preceded by variable selection procedures. Many variable selection methods are based on regularization approaches (Guo, 2010; Cai and Liu, 2011; Witten and Tibshirani, 2011; Fan *et al.*,

2012; Mai *et al.*, 2012; Xu *et al.*, 2014; Gaynanova *et al.*, 2016; Safo and Ahn, 2016), which require iterative estimation of high-dimensional parameters, including computation of a $p \times p$ precision matrix (Xu *et al.*, 2014). It is unclear whether these regularization methods can be directly applied to ultrahigh-dimensional classification. Furthermore, the conditions that guarantee selection consistency may fail to hold for ultrahigh-dimensional cases.

Computationally more efficient screening methods (Fan and Fan, 2008; Fan and Lv, 2008; Pan *et al.*, 2016; Yu *et al.*, 2016) and Bayesian methods (Johnson and Rossell, 2012; Johnson, 2013; Nikooienejad *et al.*, 2016; Rossell and Rubio, 2018) have also been developed for (ultra)high-dimensional variable selection. However, most of the screening methods in literature require the informative features to have strong marginal discriminant effects and ignore the interfeature correlations, therefore are not designed for weak signal selection. While in ultrahigh-dimensional settings, many marginally weak signals have strong predictive effects on the outcome classes. As shown in Figure 1, gene *IPO5* does not have a sufficient power to distinguish tissues with C and AR rejection types. However, jointly with gene

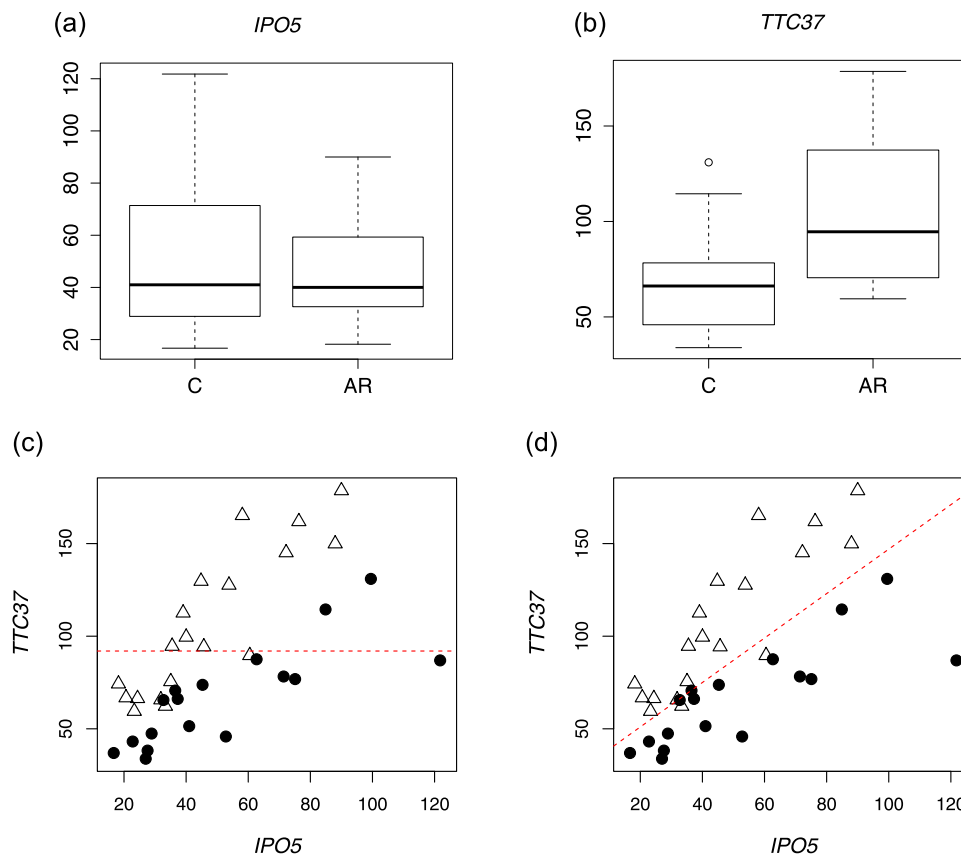


FIGURE 1 The roles of MI and JI features on classification. A, The marginal scores of gene *IPO5* between C and AR in the kidney transplant data are similar and thus *IPO5* is not MI. B, The marginal scores of a MI gene *TTC37*. C, Classification of C (circles) and AR (triangles) based on gene *TTC37*. D, Classification based on both *TTC37* and *IPO5* gives the better performance than *TTC37* only. JI, jointly informative; MI, marginally informative [Color figure can be viewed at wileyonlinelibrary.com]

TTC37, a marginally informative (MI) feature, the classification accuracy can be much improved. In this case, we call gene *IPO5* a marginally weak but jointly informative (JI) feature.

Furthermore, many of the high-dimensional classification methods aforementioned are designed for binary classifications. Multiclass classification is more challenging than binary cases (Hastie *et al.*, 2009; Gaynanova *et al.*, 2016). Most multiclass classification methods rely on sequential binary classifications by way of one-versus-the rest (Bishop, 2006), direct pairwise comparison (Bishop, 2006), direct graph traversal (Platt *et al.*, 2000), error-correcting output coding (Allwein *et al.*, 2000), multiclass objective functions (Weston and Watkins, 1998), sequential approaches (Cai and Liu, 2011; Witten and Tibshirani, 2011; Mai *et al.*, 2012), or simultaneous canonical vector estimation (Gaynanova *et al.*, 2016; Mai *et al.*, 2019). However, the choice of reduction method from multiclass to binary is on a case-by-case basis and is not a trivial task (Allwein *et al.*, 2000). In particular, commonly used pairwise comparisons are involved with a large number of individual classifiers, which is likely to incur misclassification

error and numerical instability with small sample sizes (Wu *et al.*, 2004). On the other hand, LDA can perform multiclass classification without resorting to pairwise comparisons.

The covariance-enhanced discriminant analysis method proposed by Xu *et al.* (2014) requires estimating the $p \times p$ -dimensional precision matrix of the covariates and is not computationally feasible in ultrahigh-dimensional settings. The pairwise sure independent screening for multiclass LDA (pairwiseLDA) proposed by Pan *et al.* (2016) uses the independence rule and ignores the interfeature correlations. It cannot detect marginally weak signals. Furthermore, the marginal sliced inverse regression (SIR) for model-free feature selection and multiclass classification proposed by Yu *et al.* (2016) selects the linear combinations of features, and cannot select individual features. It therefore lacks interpretation for the selected features. The general sparse multiclass LDA (mLDA) proposed by Safo and Ahn (2016) projects the original feature space to a low-dimensional canonical subspace, and therefore cannot select individual features either. The work of Cai *et al.* (2018) is designed for analyzing dependent data with a large number of

samples, but not particularly for high-dimensional feature selection.

We propose an ultrahigh-dimensional multiclass classification method within the framework of Fisher's LDA. Our proposal, termed mLDA, embeds a computationally feasible screening procedure, specially designed for detection of weak signals by accounting for inter-feature correlations. We show that the proposed method can recover all the informative features, including both MI and JI features, with probability tending to one and can achieve an asymptotically negligible misclassification rate.

The rest of the paper is organized as follows. Section 2 introduces mLDA and Section 3 develops its theoretical properties. In Section 4, the performance of the proposed method is evaluated using simulation studies. We apply the proposed procedure to analyze the renal transplantation data in Section 5 and conclude the paper with a discussion in Section 6. Technical details are provided in the Web Appendices.

2 | ULTRAHIGH-DIMENSIONAL MULTICLASS CLASSIFICATION

2.1 | Notation

Denote by \mathbf{A}' the transpose of a $p \times p$ matrix \mathbf{A} and by A_{jk} the (j, k) th entry of \mathbf{A} , where $1 \leq j, k \leq p$. Let $|\mathcal{S}|$ be the cardinality of a set \mathcal{S} and \mathcal{S}^c be the complement of \mathcal{S} . We denote the trace of \mathbf{A} as $\text{tr}(\mathbf{A})$, the minimum and maximum eigenvalues of \mathbf{A} as $\lambda_{\min}(\mathbf{A})$ and $\lambda_{\max}(\mathbf{A})$, and the operator norm and the Frobenius norm as $\|\mathbf{A}\| = \lambda_{\max}^{1/2}(\mathbf{A}'\mathbf{A})$ and $\|\mathbf{A}\|_F = \text{tr}(\mathbf{A}'\mathbf{A})^{1/2}$, respectively. Let $\mathbf{X} = (X_1, \dots, X_p)'$ be a p -dimensional vector of features. We refer to X_j as feature j for short, $1 \leq j \leq p$. Denote by Σ the covariance matrix of \mathbf{X} and by $\Omega = \Sigma^{-1}$ the precision matrix. Let $\mathcal{G}(\mathcal{V}, \mathcal{E}; \Omega)$ be the graph induced by Ω , where $\mathcal{V} = \{1, \dots, p\}$ is the vertex set and \mathcal{E} is the edge set. An edge refers to a pair of two vertices, j and j' , which satisfies $\Omega_{jj'} \neq 0$. For a subset $\mathcal{V}_l \subset \mathcal{V}$, denote by Ω_l the principal submatrix of Ω with its row and column indices restricted to \mathcal{V}_l . Denote by \mathcal{E}_l the corresponding edge set. A subgraph $\mathcal{G}(\mathcal{V}_l, \mathcal{E}_l, \Omega_l)$ is a connected component in Ω if any two vertices in \mathcal{V}_l are connected, and for $j \in \mathcal{V}_l^c$, then $\Omega_{jj'} = 0$ for any $j' \in \mathcal{V}_l$. We write $\mathcal{G}(\Omega) = \mathcal{G}(\mathcal{V}, \mathcal{E}; \Omega)$ for short when there is no confusion.

2.2 | MI features and JI features

Consider a K -class classification problem, where $K \geq 2$. Denote by Y the class membership and assume that the covariate vector \mathbf{X} satisfies

$$\mathbf{X} \{Y = k\} \sim N(\boldsymbol{\mu}_k, \Sigma), \quad k = 1, \dots, K,$$

where $\boldsymbol{\mu}_k = (\mu_{k1}, \dots, \mu_{kp})'$ is a p -dimensional mean vector of class k and Σ is the common covariance matrix for all K classes. Let $(Y_1, \mathbf{X}_1), \dots, (Y_n, \mathbf{X}_n)$ be n independent observations of (Y, \mathbf{X}) , where $\mathbf{X}_i = (X_{i1}, \dots, X_{ip})'$, $i = 1, \dots, n$. Denote by n_k the number of observations in class k such that $\sum_{k=1}^K n_k = n$. For a pair of classes k and k' , Fisher's rule, which can also be considered as a Bayes rule with equal prior probabilities, assigns an observation to class k over class k' if $(\mathbf{X} - \boldsymbol{\mu}_k/2)' \Sigma^{-1} \boldsymbol{\mu}_k > (\mathbf{X} - \boldsymbol{\mu}_{k'}/2)' \Sigma^{-1} \boldsymbol{\mu}_{k'}$. This naturally leads to the following classification rule:

$$\hat{Y} = \arg \max_{1 \leq k \leq K} (\mathbf{X} - \boldsymbol{\mu}_k/2)' \Sigma^{-1} \boldsymbol{\mu}_k. \quad (1)$$

When $p < n$, $\boldsymbol{\mu}_k$ and Σ can be estimated by $\hat{\boldsymbol{\mu}}_k = \sum_{i: Y_i=k} \mathbf{X}_i/n_k$ and $\hat{\Sigma} = \sum_{k=1}^K \sum_{i: Y_i=k} (\mathbf{X}_i - \hat{\boldsymbol{\mu}}_k)(\mathbf{X}_i - \hat{\boldsymbol{\mu}}_k)'/(n - K)$. However, when $p > n$, (1) is ill-posed as $\hat{\Sigma}$ is singular. Hence, a variable selection procedure is usually required to precede classification under some sparsity assumption. It can be shown that a sufficient and necessary condition for feature j , where $1 \leq j \leq p$, to be informative is

$$\mathcal{S}_0 = \left\{ 1 \leq j \leq p: \sum_{j'=1}^p \Omega_{jj'} (\mu_{kj'} - \mu_{k'j'}) \neq 0 \right. \\ \left. \text{for some } 1 \leq k < k' \leq K \right\}.$$

Under the faithfulness condition that any MI features must belong to \mathcal{S}_0 , the informative features consist of the following two mutually exclusive sets:

$$\mathcal{S}_1 = \{1 \leq m \leq p: \mu_{km} - \mu_{k'm} \neq 0 \\ \text{for some } 1 \leq k < k' \leq K\}$$

and

$$\mathcal{S}_2 = \left\{ j \in \mathcal{S}_1^c: \sum_{m \in \mathcal{S}_1} \Omega_{jm} (\mu_{km} - \mu_{k'm}) \neq 0 \right. \\ \left. \text{for some } 1 \leq k < k' \leq K \right\}, \quad (2)$$

where \mathcal{S}_1 and \mathcal{S}_2 contain the MI and JI features, respectively.

Though identifying marginally weak features is challenging in general, the JI features can be found by

searching the connected components in Ω which contain at least one feature in \mathcal{S}_1 . Furthermore, Theorem 1 in Section 3 shows that the connected components in Ω can be accurately recovered by thresholding the corresponding sample covariance matrix.

2.3 | Algorithm of mLDA

Given the training dataset $\{Y_i, \mathbf{X}_{ij}\}_{i=1}^n$, denote by $\bar{X}_j^{(k)} = n_k^{-1} \sum_{i:Y_i=k} X_{ij}$ the sample mean of feature j within class $k \in \{1, \dots, K\}$. Denote by $\tilde{\Sigma}$ the thresholded sample covariance matrix. That is, $\tilde{\Sigma}_{jj'} = \hat{\Sigma}_{jj'} \mathbf{1}(|\hat{\Sigma}_{jj'}| \geq \alpha)$, $1 \leq j, j' \leq p$, where $\hat{\Sigma}_{jj'}$ is the (j, j') th entry of $\hat{\Sigma}$, $\mathbf{1}(\cdot)$ is the indicator function, and α is a threshold. For a pair of classes (k, k') , $1 \leq k < k' \leq K$, select the set $\hat{\mathcal{S}}_1(k, k')$ containing indices m which satisfy

$$|\bar{X}_m^{(k)} - \bar{X}_m^{(k')}| > \tau,$$

where $\tau > 0$ is a thresholding parameter controlling the size of $\hat{\mathcal{S}}_1(k, k')$. Denote by $\hat{\mathcal{S}}_1 = \bigcup_{1 \leq k < k' \leq K} \hat{\mathcal{S}}_1(k, k')$ the set containing all of the MI features.

With $\hat{\mathcal{S}}_1$, we use the recursive labeling algorithm (Shapiro and Stockman, 2002) to identify the connected components in $\mathcal{G}(\tilde{\Sigma})$, the graph introduced by $\tilde{\Sigma}$, that contain any features in $\hat{\mathcal{S}}_1$. Suppose there are $B \leq |\hat{\mathcal{S}}_1|$ such connected components, say, $\hat{\mathcal{C}}_l$, $l = 1, \dots, B$, each containing at least one MI feature. Let $\mathcal{U} = \bigcup_{l=1}^B \hat{\mathcal{C}}_l$ with $u = |\mathcal{U}|$. Notice that $\hat{\mathcal{S}}_1 \subseteq \mathcal{U}$.

Let $\tilde{\Sigma}_l$ be the principal submatrix of $\tilde{\Sigma}$ with the row and column indices restricted to $\hat{\mathcal{C}}_l$, and compute $\hat{\Omega}_l = (\tilde{\Sigma}_l)^{-1}$. Let $\hat{\Omega}^u = \text{diag}(\hat{\Omega}_1, \dots, \hat{\Omega}_B)$ be a block diagonal matrix of dimension $u \times u$. Under the sparsity assumption, $\hat{\mathcal{C}}_l$ are of small sizes and u is much smaller than p . To detect the JI features, we only need to consider \mathcal{U} as the candidate set for them. Thus, \mathcal{S}_2 can be estimated by

$$\begin{aligned} \hat{\mathcal{S}}_2 &= \{j \in \mathcal{U} \cap \hat{\mathcal{S}}_1^c : \sum_{j' \in \hat{\mathcal{S}}_1} \hat{\Omega}_{jj'}^u (\bar{X}_{j'}^{(k)} - \bar{X}_{j'}^{(k')})| \\ &\geq \nu_n \text{ for some } 1 \leq k < k' \leq K\}, \end{aligned} \tag{3}$$

where $\nu_n > 0$ is a thresholding parameter controlling the size of selected JI features.

We denote by $\hat{\mathcal{S}}_0 = \hat{\mathcal{S}}_1 \cup \hat{\mathcal{S}}_2$ the set containing all the informative features. Finally, for a new observation with covariate vector \mathbf{X}_{new} , we determine the class membership by

$$\arg \max_{1 \leq k \leq K} d_k,$$

where d_k is the Fisher discriminant statistics for class k , defined by

$$d_k = (\mathbf{X}_{\text{new}}^s - \hat{\boldsymbol{\mu}}_k^s / 2)' \hat{\Omega}^s \hat{\boldsymbol{\mu}}_k^s. \tag{4}$$

Here $\hat{\boldsymbol{\mu}}_k$ and $\hat{\Omega}^u$ are estimated from the training data and $\mathbf{X}_{\text{new}}^s$, $\hat{\boldsymbol{\mu}}_k^s$, and $\hat{\Omega}^s$, respectively, are subvectors or submatrices of \mathbf{X}_{new} , $\hat{\boldsymbol{\mu}}_k$, and $\hat{\Omega}^u$ with the elements indexed by $\hat{\mathcal{S}}_0$.

For ease of understanding, Figure 2 depicts the flowchart of the mLDA algorithm. The proposed mLDA algorithm utilizes the dependence between the MI and JI features, and when the informative features are sparse, the proposed screening procedure to identify $\hat{\mathcal{S}}_0$ is computationally feasible, as we will demonstrate in Section 4. Moreover, as shown in Section 3, setting the tuning parameters as $\tau = O((r \log p)^s)$, $\alpha = O(n^{\xi-1/2})$ and $\nu_n = O((r(\log p) \exp(n^\xi))^{s'})$, for some $0 < r < 1$, $0 < s < 1/2$, $0 < \xi < 1$ and $0 < s' \leq 1/2$ guarantees that mLDA has selection consistency and a zero asymptotic misclassification rate.

3 | THEORETICAL PROPERTIES

Under regularity conditions (A1) to (A11) listed in appendix, mLDA possesses theoretical properties, such as the sure screening property and asymptotic vanishing postscreening misclassification rate.

For feature j , denote by $\mathcal{C}_{[j]}$ and $\hat{\mathcal{C}}_{[j]}$ the vertex sets of the connected component containing j in the graphs induced by Ω and $\tilde{\Sigma}$, respectively.

Theorem 1. *For any feature j , $1 \leq j \leq p$, suppose that $\mathcal{C}_{[j]} = O(\exp(n^\xi))$ for the ξ given in condition (A3), then, together with conditions (A5) and (A7), we have*

$$P(\mathcal{C}_{[j]} = \hat{\mathcal{C}}_{[j]}) \rightarrow 1 \text{ as } n \rightarrow \infty.$$

Therefore, the principal submatrices of Ω corresponding to the relevant connected components can be estimated by inverting the corresponding submatrices of $\tilde{\Sigma}$. For a properly chosen thresholding parameter α , Bickel and Levina (2008) and Fan *et al.* (2011) showed that the estimated precision matrix using $\tilde{\Sigma}$ is consistent.

Theorem 2 (Sure screening property). *Under conditions (A1) to (A9) and (A11),*

$$P(\mathcal{S}_0 \subseteq \hat{\mathcal{S}}_0) \rightarrow 1 \text{ as } n \rightarrow \infty.$$

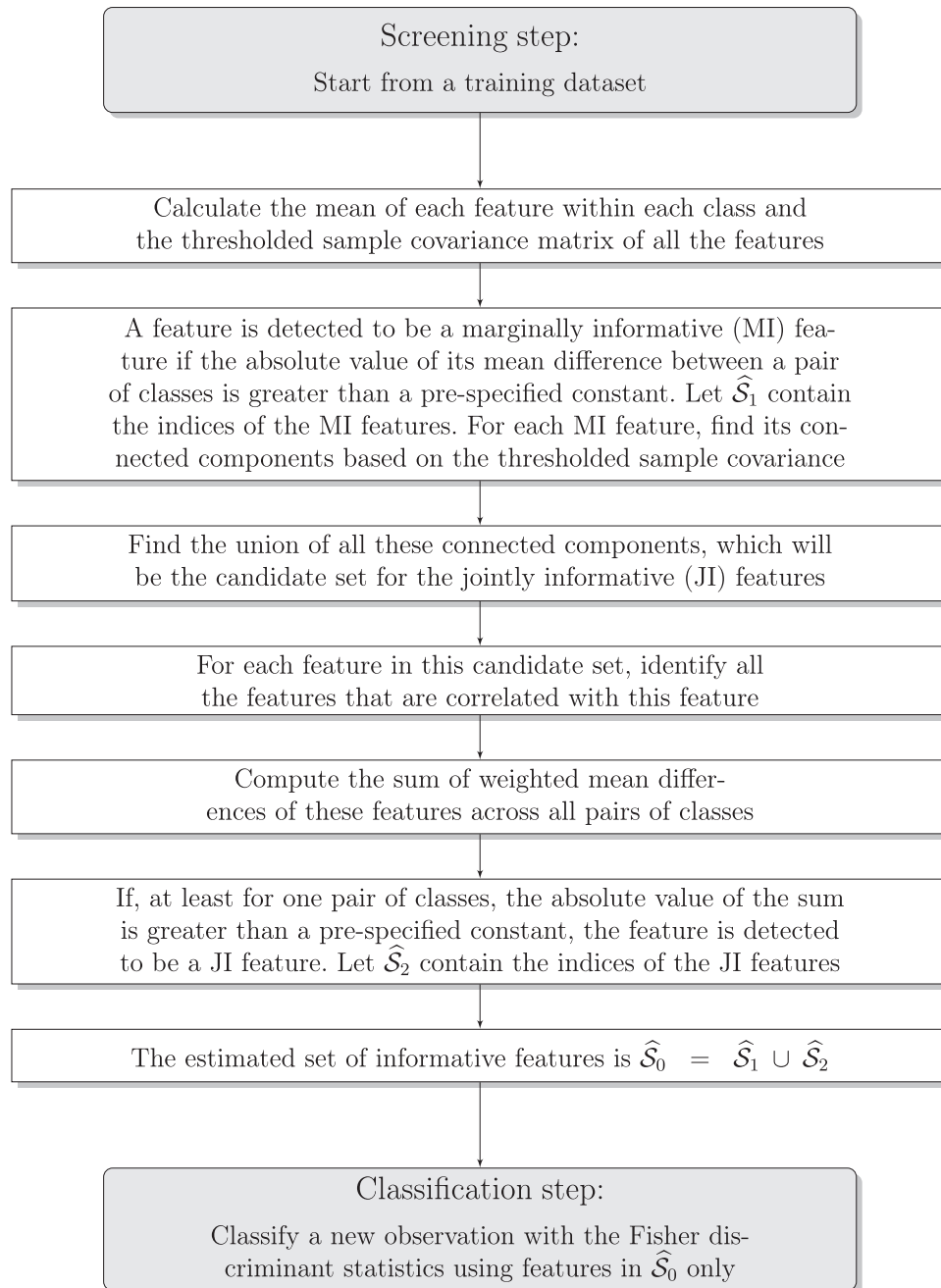


FIGURE 2 Flowchart of the mLDA procedure. mLDA: multiclass linear discriminant analysis

Theorem 3 (False-positive control property). *Under conditions (A1) to (A8), (A10) and (A11), for any $\zeta = o(n \log p)$, we have*

$$P(|\widehat{\mathcal{S}}_0 \cap \mathcal{S}_0^c| \leq \zeta^{-1} |\mathcal{S}_0^c|) \rightarrow 1 \text{ as } n \rightarrow \infty.$$

Remark. Theorems 2 and 3 hold for any distributions satisfying (A1) given in appendix. Condition (A1) characterizes a rich family of distributions, including distributions with polynomial tails such as the t distribution.

Given the training samples $\mathcal{D} = \{Y_i, \mathbf{X}_i\}_{i=1}^n$, we can assess the conditional misclassification rate for a class k of mLDA by

$$R_{\text{mLDA}}(k; \mathcal{D}) = P(\arg \max_{1 \leq l \leq K} (\mathbf{X}_{\text{new}}^S - \widehat{\boldsymbol{\mu}}_l^S / 2)' \widehat{\boldsymbol{\Omega}}^S \widehat{\boldsymbol{\mu}}_l^S \neq k | Y_{\text{new}} = k; \mathcal{D}),$$

where $\widehat{\boldsymbol{\mu}}_l^S$, and $\widehat{\boldsymbol{\Omega}}^S$ are estimated from \mathcal{D} . As pointed out in Shao *et al.* (2011), by the dominated convergence

theorem, it suffices to focus on the conditional misclassification rate, instead of the unconditional misclassification rate. We define the overall conditional misclassification rate of mLDA as

$$R_{\text{mLDA}}(\mathcal{D}) = K^{-1} \sum_{k=1}^K R_{\text{mLDA}}(k; \mathcal{D}).$$

Theorem 4 (Asymptotic vanishing postscreening misclassification rate). *For any pair of classes $1 \leq k < k' \leq K$, let*

$$\Delta_p^2(k, k') = \min_{1 \leq k < k' \leq K} (\boldsymbol{\mu}_k^0 - \boldsymbol{\mu}_{k'}^0)' \boldsymbol{\Omega}^0 (\boldsymbol{\mu}_k^0 - \boldsymbol{\mu}_{k'}^0),$$

where the superscript “0” denotes subvectors or submatrices with indices restricted to \mathcal{S}_0 . Let $\Delta_p^2 = \min_{1 \leq k < k' \leq K} \Delta_p^2(k, k')$. Under conditions (A2) to (A11), when classifying \mathbf{X}_{new} based on features selected from the screening step, for sufficiently large n , we have

$$R_{\text{mLDA}}(\mathcal{D}) \leq K\Phi(-(1 + O_p(a_n))^{1/2}(1 + O_p(\rho_n))^{1/2}\Delta_p/2) + o_p(1),$$

where Φ is the standard normal distribution function, ρ_n is given in (A11) and

$$a_n \equiv \min_{1 \leq k < k' \leq K} \max \left\{ \frac{|\mathcal{S}_0|^{1/2}}{n^{1/2}\Delta_p(k, k')}, \frac{|\mathcal{S}_0|}{n\Delta_p^2(k, k')}, \frac{1}{\Delta_p^2(k, k')} \right\}.$$

Furthermore, if $\Delta_p^2 \min\{n/|\mathcal{S}_0|, 1\} \rightarrow \infty$, then $R_{\text{mLDA}}(\mathcal{D}) \rightarrow 0$.

Notice that the condition $\Delta_p^2 \min\{n/|\mathcal{S}_0|, 1\} \rightarrow \infty$ is weaker than $n\Delta_p^2/p \rightarrow \infty$, which is required for achieving an asymptotic zero misclassification rate under the independent rule (Fan and Fan, 2008).

4 | SIMULATION STUDIES

We compared the finite sample performance of mLDA with that of other ultrahigh-dimensional classification methods, including the penalizedLDA (Witten and Tibshirani, 2011), the regularized risk minimization package (bmrn) (Teo *et al.*, 2010), the multigroup sparse discriminant analysis (MGSDA) (Gaynanova *et al.*, 2016), pairwiseLDA (Pan *et al.*, 2016), SIR (Yu *et al.*, 2016), the feature annealed independence rule (MS) (Fan and Fan, 2008), and the sure independence screening (SIS) (Fan and Lv, 2008). As an oracle benchmark, we also applied Fisher’s rule with informative features known a priori. We

first investigated the cases where the variance-covariance matrices of the features were equal across different classes, and, hence, the classes were linearly separable. We specifically considered the following two models.

Model 1 (multivariate normal distribution): Set $K = 3$ with class sizes $n_1 = n_2 = n_3 = 100$ and $p = 10\,000$. Variables 1 to 30 were generated from a multivariate normal distribution with the means specified as in Table 1. These 30 features were divided into six independent blocks: X_1 – X_5 , X_6 – X_{10} , X_{11} – X_{15} , X_{16} – X_{20} , X_{21} – X_{25} , and X_{26} – X_{30} . Features within the block were governed by the same covariance structures such as compound symmetry (CS), first-order autocorrelation (AR1), banded, star, and “unstructured” (Un). The explicit form of the last three covariance structures was given in (5). When the covariance structure required a correlation coefficient parameter, we used $\rho = 0.35$. In this case, variables 5 to 10, 15 to 20, 25 to 30 were MI features, whereas X_1 – X_4 , X_{11} – X_{14} were considered JI features for class pair (1, 2), X_{21} – X_{24} were considered JI features for class pair (1, 3), and X_1 – X_4 , X_{11} – X_{14} , and X_{21} – X_{24} were considered JI features for class pair (2, 3). The remaining noninformative 9970 features were independently generated from $N(0, 1)$ and were independent of the first 30 variables.

Model 2 (multivariate t distribution): It was the same as model 1 except that variables 1 to 30 were generated from the multivariate t distribution with four degrees of freedom and the remaining noninformative 9970 features were independently generated from the univariate t_4 distribution and were independent of the first 30 features:

$$\begin{aligned} \text{Banded: } & \begin{pmatrix} 1 & \rho & 0 & 0 & 0 \\ \rho & 1 & \rho & 0 & 0 \\ 0 & \rho & 1 & \rho & 0 \\ 0 & 0 & \rho & 1 & \rho \\ 0 & 0 & 0 & \rho & 1 \end{pmatrix}, & \text{Star: } & \begin{pmatrix} 1 & \rho & \rho & \rho & \rho \\ \rho & 1 & 0 & 0 & 0 \\ \rho & 0 & 1 & 0 & 0 \\ \rho & 0 & 0 & 1 & 0 \\ \rho & 0 & 0 & 0 & 1 \end{pmatrix}, \\ \text{Un: } & \begin{pmatrix} 1 & \rho & 0 & 0 & \rho \\ \rho & 1 & \rho & \rho & 0 \\ 0 & \rho & 1 & \rho & 0 \\ 0 & \rho & \rho & 1 & 0 \\ \rho & 0 & 0 & 0 & 1 \end{pmatrix}. \end{aligned} \tag{5}$$

TABLE 1 Means of the informative features

Features	Class 1	Class 2	Class 3
X_1 – X_4 , X_{11} – X_{14} , X_{21} – X_{24}	0	0	0
X_5 , X_{15}	0	2.5	0
X_6 – X_{10} , X_{16} – X_{20}	1.5	–1.5	–1.5
X_{25}	0	0	2.5
X_{26} – X_{30}	–1.5	–1.5	1.5

Even though condition (A7) provides the orders of the tuning parameters so that mLDA will render the desired theoretical properties, these orders do not provide specific ranges of the tuning parameters with given p and n in practice. In our numerical studies, we used fivefold cross-validation to choose the optimal tuning parameters, α , τ and ν_n .

To assess the performance in feature selection, we used false positives (FP), false negatives (FN), and the minimum number of features needed to include all informative features (MMS). To assess the classification performance, we used the number of misclassified cases (ER). The simulation results were reported in Table 2. It appears that mLDA had the lowest FP, FN, ER, and MMS under various covariance structures. When features did not follow multivariate Gaussian distributions, ERs tended to be larger across all of the methods, compared to the multivariate Gaussian. However, mLDA consistently outperformed the other methods. More simulation results with different correlations were reported in Table S6 in the Web Appendices.

We next investigated the performance of mLDA when the variance-covariance matrices differed across classes so that the classes were not linearly separable. We compared the classification performance of mLDA with various nonlinear classification methods, including the mixed discriminant analysis (Hastie and Tibshirani, 1995), the quadratic discriminant analysis (Ripley, 1996), the regularized discriminant analysis (Hastie et al., 1995), the shrunken-centroids regularized discriminant analysis (Guo et al., 2005), neural network (Ripley, 1996), kernel support vector machine (Hsu and Lin, 2002), k -nearest neighbors (Torgo, 2010), and naive Bayes (Ng and Jordan, 2001).

Model 3 (heterogeneous covariance): We set $K = 3$, $n_1 = n_2 = n_3 = 100$, and $p = 10\,000$. The variables in each class were simulated from multivariate normal distributions with the same mean structure as in model 1 and a class-specific covariance matrix.

To increase the heterogeneity and the level of nonlinearity of the class boundaries, we let different classes have different correlation coefficients in the covariance matrices. The “unstructured” covariance matrix in (5) was used for each of the six blocks within the first 30 features. Class-specific correlation coefficients were set to be ρ_1 , ρ_2 , and ρ_3 for the first, second, and third classes, respectively. The remaining noninformative 9970 features were independently generated from $N(0, 1)$ and were independent of the first 30 features.

The results reported in Tables S4 and S5 in the Web Appendices showed that, in most cases considered, mLDA still outperformed the other linear classification methods, in terms of selecting JI features and classifica-

tion accuracy. The mLDA procedure also outperformed the nonlinear classification methods in classification accuracy when the classes were nearly linear separable. As expected, the classification performance of mLDA deteriorated as the classes became more linearly inseparable.

5 | CLASSIFICATION OF POSTTRANSPLANT REJECTION TYPES

We applied the proposed mLDA to classify postkidney transplant rejection types based on patients' gene expressions. The kidney transplant study (Fletcher et al., 2004) had a total of 62 kidney tissue samples taken from 17 normal donor kidneys (C), 19 well-functioning kidneys more than 1-year posttransplant (TX), 13 biopsy-confirmed acute rejection (AR), and 13 acute dysfunction with no rejection (NR). Each sample was microarrayed (by HG-U95Av2 GeneChips; Affymetix) with 12 625 genes from kidney biopsies and peripheral blood lymphocytes at transplant.

For comparisons, we also considered regularization methods including the regularized optimal affine discriminant (ROAD) by Fan et al. (2012), the linear programming discriminant (LPD) by Cai and Liu (2011), the covariance-enhanced discriminant analysis (CED) by Xu et al. (2014), and screening methods including MS (Fan and Fan, 2008), SIS and the iterative SIS (ISIS) by Fan and Lv (2008). Since ROAD, LPD, and CED cannot handle ultrahigh-dimensional data, we performed variable selection using mLDA before applying the corresponding regularization method.

The classification performance was assessed by the leave-one-out procedure. The thresholding parameters τ , α , and ν_n in (3) were chosen by fivefold cross-validation. ROAD, LPD, and the R package SIS, which implements the SIS and ISIS methods, cannot handle categorical outcomes with $K > 2$. When implementing them, we first carried out pairwise comparisons between rejection types, and then used the majority vote to decide the final membership. The binary classification approaches inadvertently produced ties, which made the final class membership assignment difficult. When a tie occurred, we randomly assigned a class membership among the tied rejection types. Contrarily, mLDA, which performed multiclass classification without resorting to pairwise comparisons, did not encounter the tie issue. It turned out that the numbers of misclassified tissues given by mLDA, ROAD, LPD, CED, MS, SIS, and ISIS were 6, 9, 12, 8, 15, 16, and 13, respectively.

TABLE 2 Comparisons with the competing methods

	Model 1					Model 2				
	CS	ARI	Band	Star	Un	CS	ARI	Band	Star	Un
FP										
mLDA	9.4 (0.9)	9.5 (0.9)	16.2 (1.0)	16.3 (2.1)	17.4 (2.3)	7.5 (0.9)	8.5 (1.0)	15.8 (2.1)	19.6 (3.5)	28.0 (4.1)
MS	19.8 (1.2)	24.0 (1.2)	59.4 (2.4)	41.6 (1.3)	59.4 (2.1)	18.2 (1.2)	26.4 (1.1)	59.1 (2.3)	39.3 (2.1)	40.7 (2.0)
pairwiseLDA	16.7 (2.3)	18.9 (2.6)	24.1 (3.0)	20.5 (3.0)	25.7 (3.1)	17.1 (2.6)	20.6 (2.7)	25.5 (3.1)	24.1 (2.9)	27.3 (3.2)
SIR	34.2 (5.7)	35.6 (6.8)	37.1 (5.9)	38.3 (6.0)	37.8 (7.2)	33.9 (6.3)	36.2 (7.7)	35.1 (6.4)	37.7 (6.6)	39.3 (7.8)
bmm	24.0 (4.6)	26.6 (5.1)	23.9 (4.4)	25.4 (5.2)	27.1 (5.6)	24.3 (6.3)	25.7 (5.9)	25.9 (6.0)	26.1 (6.1)	28.8 (6.6)
MGSDA	45.1 (7.2)	43.2 (6.7)	49.0 (6.9)	47.1 (7.0)	47.3 (7.3)	48.8 (8.7)	49.3 (7.9)	50.1 (7.4)	47.4 (7.6)	51.2 (8.5)
SIS	0.8 (0.0)	1.8 (0.5)	2.2 (0.7)	0.3 (0.01)	1.1 (0.01)	1.4 (0.1)	2.1 (0.3)	2.2 (0.4)	1.0 (0.2)	1.0 (0.1)
FN										
mLDA	0.3 (0.9)	0.6 (1.2)	0.5 (1.5)	1.0 (1.4)	1.1 (1.5)	1.7 (1.7)	2.3 (1.9)	2.5 (2.1)	4.7 (2.1)	10.3 (1.1)
MS	12.3 (0.9)	12.0 (0.9)	12.4 (1.2)	13.1 (1.0)	11.9 (1.4)	11.5 (0.9)	12.4 (0.8)	11.9 (0.8)	11.8 (0.9)	11.8 (1.4)
pairwiseLDA	12.2 (0.9)	12.7 (1.1)	12.1 (1.0)	13.0 (1.0)	12.9 (1.1)	12.3 (1.0)	13.1 (1.0)	12.4 (1.0)	13.3 (1.1)	13.4 (1.1)
SIR	12.0 (0.0)	12.0 (0.0)	12.0 (0.0)	12.0 (0.0)	12.0 (0.0)	12.0 (0.0)	12.0 (0.0)	12.0 (0.0)	12.0 (0.0)	12.0 (0.0)
bmm	12.0 (0.0)	12.0 (0.0)	12.0 (0.0)	12.0 (0.0)	12.0 (0.0)	12.1 (0.2)	12.0 (0.0)	12.0 (0.0)	12.1 (0.2)	12.1 (0.2)
MGSDA	13.0 (2.4)	12.7 (1.8)	13.4 (2.5)	13.5 (2.2)	13.7 (2.5)	14.1 (3.2)	13.6 (3.0)	13.9 (3.3)	14.2 (3.1)	13.8 (3.2)
SIS	20.7 (1.5)	20.9 (1.3)	20.7 (1.3)	20.4 (1.2)	21.0 (1.3)	20.9 (1.4)	21.3 (1.3)	21.6 (1.5)	20.5 (1.5)	21.0 (1.5)
MMS										
mLDA	38.7 (2.1)	36.2 (3.4)	42.1 (3.0)	39.8 (5.8)	44.0 (3.2)	48.9 (8.5)	49.1 (9.1)	44.5 (4.5)	36.3 (5.6)	48.8 (5.9)
MS	9856 (286)	9811 (243)	9773 (275)	9832 (263)	9804 (286)	9733 (293)	9864 (266)	9770 (231)	9699 (269)	9634 (247)
bmm	9673 (223)	9526 (284)	9725 (257)	9766 (279)	9422 (247)	9567 (265)	9327 (251)	9764 (270)	9327 (284)	9334 (255)
SIS	9463 (232)	9634 (243)	9721 (257)	9644 (234)	9579 (261)	9842 (270)	9756 (225)	9688 (291)	9720 (256)	9591 (248)
ER										
mLDA	6.1 (2.1)	6.5 (2.3)	6.3 (2.1)	2.7 (1.8)	5.0 (2.2)	10.4 (3.0)	10.5 (3.0)	7.6 (2.4)	4.1 (2.1)	9.7 (2.9)

(Continues)

TABLE 2 (Continued)

	Model 1					Model 2				
	CS	ARI	Band	Star	Un	CS	ARI	Band	Star	Un
MS	10.2 (2.6)	10.3 (2.7)	9.4 (2.5)	3.0 (2.7)	7.8 (2.6)	11.8 (3.2)	11.8 (3.0)	11.0 (3.1)	5.4 (2.1)	10.7 (3.1)
pairwiseLDA	8.8 (3.2)	9.2 (3.6)	8.4 (3.3)	8.0 (3.1)	8.3 (3.4)	12.4 (3.7)	11.5 (3.9)	11.7 (3.5)	10.3 (3.4)	11.4 (3.5)
SIR	43.2 (14.6)	42.1 (13.2)	39.7 (13.5)	36.5 (13.8)	44.0 (13.9)	42.9 (14.3)	41.8 (14.5)	40.5 (14.3)	41.4 (13.6)	46.7 (14.7)
penalizedLDA	47.9 (12.1)	43.3 (14.0)	45.8 (13.2)	47.7 (12.8)	49.3 (11.9)	48.0 (14.7)	49.2 (15.2)	50.4 (15.0)	52.1 (14.1)	51.5 (13.5)
bmm	28.1 (6.2)	24.6 (6.7)	29.3 (6.3)	28.4 (6.5)	33.3 (6.8)	31.6 (7.1)	28.7 (6.9)	30.9 (7.0)	32.4 (7.3)	36.8 (7.1)
MGSDA	9.3 (3.7)	8.7 (3.6)	9.9 (3.7)	10.0 (4.0)	7.5 (4.1)	11.3 (4.3)	12.0 (4.2)	10.4 (4.5)	9.7 (4.5)	12.5 (4.7)
SIS	17.9 (3.5)	17.3 (2.8)	7.4 (3.4)	5.0 (3.1)	17.7 (3.9)	19.7 (3.4)	19.6 (3.8)	9.3 (3.1)	5.5 (3.5)	20.2 (4.8)
Oracle	5.2 (2.0)	6.0 (1.9)	5.8 (1.9)	2.4 (1.6)	4.6 (2.1)	9.7 (3.1)	8.9 (3.3)	6.5 (3.0)	3.2 (2.2)	8.8 (3.0)

Abbreviation: ARI, first-order autocorrelation covariance matrix; Band, banded structured covariance matrix; CS, compound symmetry covariance matrix; ER, the number of misclassified cases; FN, average number of false negatives; FP, average number of false positives; LDA, linear discriminant analysis; MGSDA, multigroup sparse discriminant analysis; MMS, the minimum number of features needed to include all informative features; SIR, sliced inverse regression; SIS, sure independence screening; Star, star-shape structured covariance matrix; Un, “unstructured” covariance matrix.

In model 1, features were generated from multivariate normal distribution and in model 2, features were generated from multivariate t distribution. Numbers in parentheses are interquartile ranges for MMS and standard deviations for FP, FN, and ER. MMS is reported only for methods that output the full ranks of all features. The competing methods include penalizedLDA (Witten and Tibshirani, 2011), the regularized risk minimization package (bmm) (Teo et al., 2010), the MGSDA (Gaynanova et al., 2016), pairwiseLDA (Pan et al., 2016), marginal SIR (Yu et al., 2016), the feature annealed independence rule (MS) (Fan and Fan, 2008), and the SIS (Fan and Lv, 2008).

For each gene, we computed the frequency of its being selected during the leave-one-out procedure. The top 10 genes with the highest selection frequency were given in Table 3. Among them, the JI genes *CEACAM8*, *RNASE3*, *TCN1*, *BPI*, and *CRISP3* were all highly correlated with the MI gene *TCF12*, while the JI gene *IGHV3-23* was highly correlated with the MI gene *HLA-G*. Our results have biological explanations. For example, the identified gene *TCN1* encodes a member of the vitamin B12 binding protein family and vitamin B12 inefficiency can cause kidney injury (Gowder, 2014). Gene *BPI* fold-containing family A member 2/parotid secretory protein is associated with acute kidney injury (Kota et al., 2017). Gene *TCF12* is expressed in the forming collecting ducts in the developing kidney as well as in the liver (Lazzaro et al., 1992), while *HLA-G* expression in biliary epithelial cells is associated with allograft acceptance in liver-kidney transplantations (Xiao et al., 2013). Our study has also identified some novel genes, such as genes *CEACAM8* (a carcinoembryonic antigen related to cell adhesion), *RNASE3*

(associated with allergic rhinitis), and *CRISP3* (strongly upregulated in prostate carcinomas), which are all JI features and have not been reported in transplant literature.

TABLE 3 Top 10 genes selected from the posttransplant rejection study

Genes	Selection frequency	MI or JI
<i>CEACAM8</i>	0.90	JI
<i>TCN1</i>	0.85	JI
<i>HLA-G</i>	0.84	MI
<i>BPI</i>	0.82	JI
<i>GUSBP11</i>	0.82	MI
<i>IGHV3-23</i>	0.79	JI
<i>TAF6L</i>	0.74	MI
<i>RNASE3</i>	0.73	JI
<i>CRISP3</i>	0.66	JI
<i>TCF12</i>	0.58	MI

Abbreviation: JI, jointly informative; MI, marginally informative.

6 | DISCUSSION

The proposed mLDA can be easily extended to accommodate non-Gaussian covariates. Indeed, the results in Section 4 hinted that mLDA works well for the heavy-tailed t distribution. Our proposed mLDA is based on the assumption of the common covariance matrix across classes. We have examined the performance of our proposal when the covariance matrices vary across classes. Our empirical results show that mLDA performs reasonably well even under the misspecified models, as long as the common covariance assumption is not severely violated. The performance of mLDA might deteriorate when the covariance matrices do differ much across classes. For these cases, the nonlinear classification methods, such as quadratic discriminant analysis, may be more appropriate.

Algorithms can be developed more efficiently. For instance, when the whole feature space can be divided into uncorrelated subspaces, such as blocks of chromosomes in a genome or different functional regions in a brain, parallel computing on multiple partitioned feature spaces can be implemented. We are also cognizant that there might be some marginally weak signals that may not be detected by our method, such as the marginally weak signals with joint pooled effects (Li and Leal, 2008). In this case, the proposed method may be used with other weak signal detection techniques to boost weak signal selection and classification performance. We will pursue this.

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SUPPORTING INFORMATION

Additional supporting information may be found online in the Supporting Information section at the end of the article. The Web Appendix, referenced in Section 4, is available with this paper at the Biometrics website on Wiley Online Library.

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APPENDIX: TECHNICAL CONDITIONS

To derive the sure screening property and asymptotic vanishing postscreening misclassification rate for mLDA, we assume the following conditions.

- (A1)** For any positive integer $l \geq 1$, $E|X_j|^l \leq l!C^l$ for some constant $C > 0$, $1 \leq j \leq p$.
- (A2)** (Faithfulness condition) For any j satisfying $\mu_{kj} - \mu_{k'j} \neq 0$, $\sum_{j'=1}^p \Omega_{jj'}(\mu_{kj'} - \mu_{k'j'}) \neq 0$ for some $1 \leq k < k' \leq K$.
- (A3)** $p = O(\exp(n^\zeta))$ for some $0 < \zeta < 1$ and $C_{\max} = O(\exp(n^\xi))$ for some $0 < \xi < \zeta$, where C_{\max} is the maximum size of the connected components in Ω containing at least one MI signal. For any MI feature $j \in \mathcal{S}_1$, if $\mu_{kj} - \mu_{k'j} \neq 0$ for some pair (k, k') , $1 \leq k < k' \leq K$, we assume

that $|\mu_{kj} - \mu_{k'j}| > \sqrt{r \log p}$, where $0 < r < 1$ controls the overall strength of the MI features.

(A4) For any subset $\mathcal{V}_l \subset \{1, \dots, p\}$ with $|\mathcal{V}_l| = O(n)$,

$$\max_{1 \leq k \leq K} \sup_j E |X_{ij} 1(Y_i = k, j \in \mathcal{V}_l)|^{2t} < C_t < \infty$$

for some constants $t > 0$ and $C_t > 0$.

(A5) $\min_{(j,j') \in \mathcal{E}} |\Sigma_{jj'}| \geq C n^{(\xi-1)/2}$ and $\max_{(j,j') \notin \mathcal{E}} |\Sigma_{jj'}| = o(n^{(\xi-1)/2})$ for the ξ in (A3).

(A6) There exist positive constants κ_1 and κ_2 such that

$$0 < \kappa_1 < \lambda_{\min}(\Sigma) \leq \lambda_{\max}(\Sigma) < \kappa_2 < \infty.$$

(A7) $\tau = O((r \log p)^s)$ for some $0 < s < 1/2$ and r given in (A3), $\alpha = O(n^{(\xi-1)/2})$ and $\nu_n = O((r(\log p) \exp(n^\xi))^{s'})$ for some $0 < s' \leq 1/2$ and ξ given in (A3).

(A8) There exist positive constants c_1 and c_2 such that

$$0 < c_1 \leq \min_{1 \leq k \leq K} \frac{n_k}{n} \leq \max_{1 \leq k \leq K} \frac{n_k}{n} \leq c_2 < \infty$$

for all n .

(A9) For any pair of classes, the number of informative features that differentiate the pair is of $O(n^\varrho)$, for some uniform constant $\varrho > 0$ not depending on the pairs.

(A10) There are $p^{-\beta}$, $0 < \beta < 1$, fraction of features that are MI and $p^{-\gamma}$, $0 < \gamma < 1$, fraction of features that are JI.

(A11) $\rho_n = n^{-1} C_t C_{\max}^{4/t} \rightarrow 0$ as $n \rightarrow \infty$, where t and C_t are given in (A4) and C_{\max} is defined in (A3).

(A1) is required to prove that the connected components of Ω can be consistently recovered by $\hat{\Sigma}$. Gaussian distributed random variables satisfy (A1). (A2) ensures that all MI features belong to \mathcal{S}_0 . (A3) implies that for each connected component of a MI feature, its size cannot exceed the order of $\exp(n^\xi)$ for some $\xi \in (0, 1)$. This condition is required for consistently estimating precision matrices by thresholding covariance matrices and is also assumed by Bickel and Levina (2008). (A4) ensures that the connected components in Ω can be adequately estimated from the connected components in $\hat{\Sigma}$; see Shao *et al.* (2011) and Bickel and Levina (2008) for details. (A5) guarantees that the true zero entries in Ω can be reliably separated from the nonzero entries in the covariance matrix, and therefore the connected components containing at least one MI feature in the precision matrix can be detected accurately (Zhang *et al.*, 2014). (A6) is commonly assumed on design matrices in high-dimensional settings (Bickel and Levina, 2008; Fan *et al.*, 2011; Shao *et al.*, 2011). (A7) gives the order of tuning parameters τ , α , and ν_n to achieve the desired theoretical properties (Bickel and Levina, 2008; Jin, 2009; Fan *et al.*, 2011; Shao *et al.*, 2011). (A8) implies that the K classes are of comparable sample sizes and the sample size of each class goes to infinity when n goes to infinity (Fan and Fan, 2008; Shao *et al.*, 2011). (A9) requires that the numbers of informative features differentiating different pairs of classes have a uniform lower bound of the order $O(n^\varrho)$. (A10) indicates the sparsity of the MI and JI features and is required for proving Theorem 3. As β and γ get closer to 1, MI and JI features become sparser, respectively (Jin, 2009). (A11) is used to gauge the covariance and precision estimation errors (Bickel and Levina, 2008; Fan *et al.*, 2011; Shao *et al.*, 2011).