

SUPPLEMENTAL MATERIAL:

R computation details

RGMM were implemented in R by adding a Ridge regularization term νI to the covariance matrix Σ_c , where the parameter ν has been tuned using a 5-fold cross-validation. HDDA method [2] has been implemented thanks to the R package *HDclassif* [1]. Implemented using the R package *e1071* [7], SVM require the setting of the kernel scale parameter and the soft margin cost, tuned by a 5-fold cross-validation. Random Forest method has two parameters to be tuned: the number of trees in the forest and the Bagging number of variables. We decided to set the number of trees to 500, and the Bagging parameter was tuned by a 5-fold cross-validation [5]. The Random Forest method was implemented thanks to the R package *randomForest* [6]. As there is no standard R package to implement FMLM, we made our own functional adaptation of the multinomial logit model in decomposing the data on a cubic spline basis (degree 3) [10], then applying the standard multivariate multinomial logit model on the obtained coefficients (we used the R package *splines* [8] to compute the cubic spline basis and the R package *nnet* [11] to build the multinomial logit model). The spline decomposition requires the choice of the number of pieces of polynomial functions (or the number of knots), tuned by a 5-fold cross-validation. NPFD requires the choice the pseudometric δ , the k-nearest neighbours parameter k and the kernel function K (the original R implementation codes of the NPFDFA method are freely available online at <http://www.math.univ-toulouse.fr/~ferraty/>. We made an adaptation the function *funopadi.knn.lcv* in order to make global the automatic choice of the k-nearest neighbours parameter k by a 5-fold cross-validation. However, Ferraty and Vieu [4] showed that the choice of the kernel function K has a negligible impact on the NPFD estimator's convergence rate compared to the the pseudometric δ and the k-nearest neighbours parameter k ; thus, we arbitrarily fixed it as the standard quadratic kernel. Among several others, we chose in this study δ as the MPLSR (Multiple Partial Least Squares Regression) pseudometric, because it seems to be particularly adapted for hyperspectral data classification [12]. **FSVM is based on an approach proposed by Rossi and Villa [9], computing the standard SVM on the coefficients of functional data decomposition by a cubic splines basis. FRF applies a Random Forest approach for functional data classification, decomposing the data on a cubic splines basis, then applying the standard Random Forest on the obtained coefficients. For both last methods, parameters have been tuned by a 5-fold cross-validation in the same way than their multivariate version.**

References

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