

Identification of Stellar Population in Galactic Spectra using the Hierarchical Decision Ensemble

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Abstract

In order to understand the process of formation and evolution of a galaxy, it is very important to identify the ages of stellar populations that make it up. The main contribution of this work is an efficient machine learning method, insensitive to emission lines and robust in the presence of noise, to determine the age of stellar populations in observational galactic spectra. It is a novel method for constructing ensembles where the decision in the prediction is done subdividing the data and organizing them in a tree structure. Experimental results show that our method yields a better accuracy than the traditional feed-forward Neural Network Ensemble.

Introduction

In astronomy, the very high spectral resolution achieved recently and the availability of very large databases such as the Sloan Digital Sky Survey (SDSS) (see (Abazajian 2003)) allows new kinds of research in spectroscopy.

To analyze the properties of celestial objects, astronomers use spectra. In a spectrum, all the information about the chemical composition of the object, is encoded. A spectrum is a plot of flux against wavelength and it is formed for three components: the continuum, absorption lines, and emission lines. The continuum is a smooth spectrum generated by solid objects and dense gases that emit radiation in a wide range of wavelength; the absorption and emission lines represent different atomic interactions in specific wavelengths of the light; they can be appreciated in Figure 1, as downward and upward lines respectively.

For astronomers, the amount of data that they need to analyze has become too large to be treated manually, therefore, a variety of machine learning methods has started to be used to perform automatic determination of parameters in spectra. Using those methods, astronomers can evaluate large amounts of information in an acceptable period of time (see (Coryn A. L. & Von-Hippel 1998), (Fuentes & Gulati 2001), (Snider 2001), (Weaver 2000)).

The most common approach that has been used consists of fitting a spectrum to a synthetic model. This approach often yields erroneous results when applied to observational

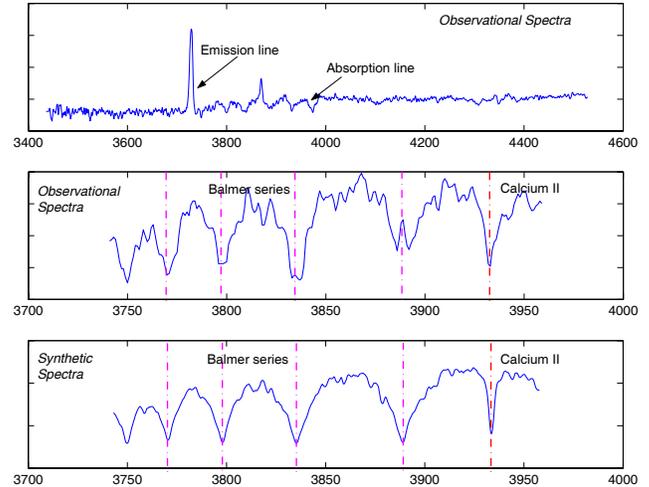


Figure 1: (1)Observational Spectrum, (2)Observational and (3)Synthetic spectra in the region of interest [3743Å – 3958Å] showing the Balmer series and the Calcium II lines

spectra, mainly because of the presence of emission lines (see figure 1), which are not taken into consideration in standard synthetic models. Those lines are considerably stronger than absorption lines and therefore they define the shape of the spectrum; also there is the problem of certain distortions such as noise, discretization effects and generalized dilution.

From a theoretical point of view, in order to understand galaxy formation and evolution, it is very important to identify the principal periods of star formation, but the task of determining the ages of stellar populations in galactic spectra is not trivial, since spectra are normally noisy and present other hard to model features.

In this work we propose a Machine Learning method that is insensitive to emission lines and robust in the presence of noise, to determine the age of stellar populations in synthetic and observational galactic spectra. It is a novel method for constructing ensembles where the decision in the prediction is done subdividing the data and organizing them in a tree structure, obtaining a good approximation when the instance reaches a leaf node or in the path behind.

The approach we propose consists of fitting just a few

lines of the spectrum that contain the information related to the age of the stellar populations. It has the additional advantage that it does not use regions with strong emission lines. Our experimental results show the efficacy of the proposed method applied to a difficult pattern recognition problem.

The organization of the remainder of this paper is as follows: Section 2 describes the data used in the experiments. Section 3 describes the process used to characterize the spectra. Section 4 presents the ensemble method used to predict the age of stellar populations. Section 5 presents experimental results using synthetic and observational spectra, and Section 6 presents conclusions and future work.

Data

The spectra are defined by 2 vectors; the first one is the wavelength and the second is the amount of flux in that point, the difference between a specific wavelength and the next is the sampling rate; the measurement unit used in astronomy for denoting the wavelength is the Angstrom (abbreviated \AA) and represents 1×10^{-10} meter.

The observational data consists of 1157 galactic spectra obtained from the Sloan Digital Sky Survey (see (Abazajian 2003)) with a sampling rate of 1\AA and a range from 3435\AA to 8315\AA .

A galactic spectrum is composed of stars, nebula and interstellar dust. In general terms an observational spectrum can be considered as the combination of a few main stellar populations. Those populations are formed by stars with similar cinematic properties, ages and chemical composition. Thus, to analyze the observational spectra, we generated simulated galactic spectra as the sum of three synthetic models of stellar populations with different ages. In order to make the simulated data as similar to the observational galactic spectra as possible, we added noise and gaussian smoothing.

The synthetic models (\vec{m}_x) are composed by 14 high resolution spectra (see (Cerviño M. 2000)) at 0.3\AA sampling and a range from 3000\AA to 6900\AA . The experiments were calibrated using three populations with different ages and proportions:

- *age1*: A young population with ages between $10^{7.0}$ and $10^{7.6}$ years, representing a starburst component, five spectra.
- *age2*: An intermediate population with ages between $10^{8.0}$ and $10^{8.6}$ years, four spectra.
- *age3*: An old population with ages between $10^{9.0}$ and $10^{9.6}$ years, representing the bulge component, five spectra.

In this work we focused on five spectral lines that encode most of the information related to the age of the stellar populations; those lines are the Balmer series and Calcium II (see figure 1) that are characteristic of the young and the old stellar populations respectively. This subset of the spectra consist of 717 points with wavelength between 3743\AA and 3958\AA .

Generation of Simulated Spectra. A simulated spectrum \vec{S}_{sim} is a combination of synthetic models that rep-

resents a galaxy with three principal stellar populations of different ages. Define M_{models} the matrix containing the 14 synthetic models; as

$$M_{models} \in \mathbb{R}^{14 \times 717} = \begin{pmatrix} \vec{m}_1 \\ \vdots \\ \vec{m}_{14} \end{pmatrix} \quad (1)$$

$\vec{p}_1 \in \mathbb{R}^{1 \times 5}$, $\vec{p}_2 \in \mathbb{R}^{1 \times 4}$ and $\vec{p}_3 \in \mathbb{R}^{1 \times 5}$ are vectors representing the proportion of the stellar populations: *age1*, *age2* and *age3* respectively, each with just one non-zero element that represents one specific age in the simulated galactic spectra, where:

$$\sum_{i=1}^3 \sum \vec{p}_i = 1 \quad (2)$$

\vec{ages} is the vector that results of concatenating \vec{p}_1 , \vec{p}_2 and \vec{p}_3 . It contains the proportions of ages, with just three non-zero elements:

$$\vec{ages} \in \mathbb{R}^{1 \times 14} = (\vec{p}_1, \vec{p}_2, \vec{p}_3) \quad (3)$$

Then *Ssim* is constructed as follows:

$$\vec{S}_{sim} = \text{smooth}(\vec{ages} * M_{models}) + \vec{S}_{noise} \quad (4)$$

where $\vec{S}_{noise} \in \mathbb{R}^{1 \times 717}$ is a random vector of the same size as \vec{S}_{sim} that represents gaussian noise. The function *smooth* is used for smoothing the spectra simulating the limited resolution of the observations; it uses gaussian window filtering with a standard deviation of 0.70 and a window size ranging from 5 to 11.

Each simulated galactic spectrum is defined by \vec{ages} . The pattern recognition task consists of predicting the non-zero elements that best fit the line profiles, producing a resultant spectrum that is as similar to the observed one as possible.

Data properties. Spectra are not chaotic signals, on the contrary, a spectrum has a defined behavior depending on its spectral parameters. In a spectrum there are several lines that always appear in certain wavelengths; those lines can be weaker or stronger depending on the structure of the object. In this case, the width of the Calcium line and the lines in the Balmer series encode the relevant information about the age of its populations.

As we can see in Table 1, the resolution in the synthetic and the observational spectra are different, and there are emission lines in the observational spectra that are not considered in the synthetic ones. These differences are very important, and they complicate the task of fitting the whole spectrum, that is the reason we decide to fit just few lines where there are not strong emission lines.

To characterize the properties of the five lines, we extract the width of the lines in 12 regions; it is a key step in the preprocessing, because we eliminate most of the noise effects and maintain the relevant information about each line. The widths are the features that we use to characterize each spectrum and make the age prediction. In this stage we can eliminate most of the noise in the spectrum because we do

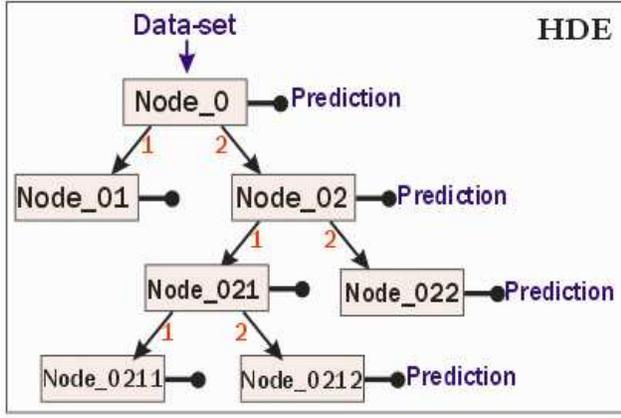


Figure 2: Tree structure of the Hierarchical Decision Ensemble

not focus on each flux point, which could be contaminated, but on general features such as the spectrum shape in a specific area. We extract 12 widths per line in 5 lines, obtaining a total of 60 features.

Hierarchical Decision Ensemble (HDE)

The hierarchical ensemble (figure 2) is a tree-based algorithm (see (Quinlan 1986)); in each step we train a classifier using a training set and predict the parameters for a validation set. The spectra corresponding to the predicted parameters are reconstructed and compared with the validation set. The predictions of those spectra that have an error that is smaller than a predefined threshold are accepted, and those spectra that are not accepted are added to the training set along with the reconstructed spectra; this new training set is divided into two subsets using a clustering algorithm, then each subgroup is used to train new classifiers in the next level of the tree (see figure 3), this is done until all the errors are lower than the threshold or until a predefined number of levels is reached.

In the prediction stage, each instance in the test set, is evaluated for the first node using the trained classifier. The prediction is employed to reconstruct a simulated spectrum. If the RMS error in the comparison is lower than a predefined threshold, the prediction is accepted. If the error is higher than the threshold, the trained divider in the node is used to determine the next path of the instance. This process is repeated until the instance obtains an accepted prediction

Table 1: Summary of properties of the data test

PROPERTY	SYNTHETIC	OBSERVATIONAL
Number of spectra	3600	1155
Resolution	0.3Å	1Å
Spectral range	[3435-8315]Å	[3000-6900]Å
Emission lines	No	Yes
Noise	10%	Un known
Number of features	60	60

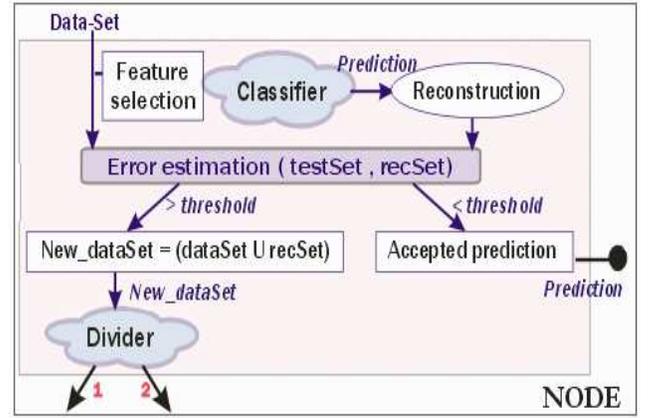


Figure 3: Internal node in HDE

or until it reaches a leaf node. The prediction and the error produced by each node are stored to make a final decision at the end of the process.

Feature Selection. We calculate the inter-class and intra-class variations for each age. These measurements are used to analyze the data behavior and we prefer those features that have low intra-class variation and high inter-class variation.

Let $X_j \in \mathbb{R}^{n_j \times m}$ be the matrix of instances relevant to the class j where $\bar{\mu}_j$ is the matrix of means in class j and $\bar{\mu}$ is the matrix of total means, n_j is the total number of instances in class j and c is the total number of classes (see (Fisher 1936)).

Define $E \in \mathbb{R}^{m \times m}$ as the inter-class variation and $I \in \mathbb{R}^{m \times m}$ as the intra-class variation:

$$E = \sum_{j=1}^c n_j \cdot (\bar{\mu}_j - \bar{\mu})(\bar{\mu}_j - \bar{\mu})^T \quad (5)$$

$$I = \sum_{j=1}^c \sum_{i=1}^{n_j} (X_{j,i} - \bar{\mu}_j)(X_{j,i} - \bar{\mu}_j)^T \quad (6)$$

Classifier. This module is independent of the algorithm, in this paper we have used standard feedforward neural networks and locally weighted linear regression, but we could use any supervised machine learning algorithm. Neural networks were chosen because of their speed at prediction time and generalization ability and LWR was introduced to add diversity to the tree. The classifier is trained using half of the total data, leaving a subset for validating its generalization ability in the specific level of the tree.

Error Estimation and Data Set Regeneration. The error estimation process uses the reconstructed galactic spectra $f(pred)$ and calculates its difference from the spectra in the test set. The resulting error is compared with a threshold, after that, we modify the data set adding those reconstructed spectra that have an RMS error higher than the threshold and removing those spectra in the test set that have an RMS error lower than the threshold. Both thresholds were set experimentally to accept about 15% of the predictions and to add about 40% of new reconstructed spectra.

Divider. In the divider module we make a clustering process; this stage is independent of the algorithm, too. In this work we use k-means (Alsabti & Singh 1998), but it could be done using any clustering algorithm. Each cluster is then passed to the next level of the tree to train a new classifier. For the experiments we use $k = 2$ for simplicity and to generate small trees, but this parameter can be modified depending on the problem.

Experimental Results

In this work we made four experiments, the task was fitting the five lines of the spectra and predicting the parameters corresponding to the age and proportion of each stellar population. Using the predicted parameters we reconstruct the suggested spectrum and compare it with the original; considering that a prediction is better than another when the reconstructed spectrum produced by this prediction is more similar to the original than the reconstructed spectrum produced by the other prediction. Bellow we will explain briefly the experiments.

ANN-E Feedforward-Backpropagation Neural Network Ensemble. It consists of 5 classifiers trained independently, using five different data sets of 3600 spectra, and combining their predictions by average.

LWR-E Locally Weighted Regression Ensemble. It is made in a similar way as the ANN-E.

ANN-LWR Heterogeneous Ensemble. It is formed by 3 classifiers trained with neural networks and 2 classifiers trained with locally weighted regression.

HDE Hierarchical Decision Ensemble. Combining Feedforward-Backpropagation Neural Networks and Locally Weighted Regression nodes; restricting the tree growth to 5 levels to obtain small trees keeping the balance between accuracy and efficiency.

We compared the performance of the HDE with the three ensembles made in a traditional way; finding that the generalization ability of the HDE was better than the other experiments realized here for the presented problem. The summary of results is shown in Table 2 and Table 3.

Table 2: RMS-Error obtained in synthetic galactic spectra using traditional ensembles and the HDE

ALGORITHM	PREDICTION TIME	SYNTHETIC GALACTIC SPECTRA
		<i>Profile fitting</i>
ANN-E	0.8 sec.	0.0451
LWR-E	1.5 sec.	0.0435
ANN-LWR	1.1 sec.	0.0447
HDE	1.7 sec.	0.0431

For the synthetic data we generated 10800 simulated galactic spectra and tested the methods in two stages, in the first one, we used ten fold cross validation, where we divide the data into 10 subsets of equal size, train the classifier 10 times, each time leaving out one of the subsets and using it for testing the algorithm. At the end of the process

Table 3: RMS-Error obtained in observational spectra using traditional ensembles and the HDE

ALGORITHM	OBSERVATIONAL GALACTIC SPECTRA	
	<i>Profile fitting</i>	<i>Reconstructed spectra</i>
ANN-E	0.348	0.411
LWR-E	0.323	0.425
ANN-LWR	0.357	0.467
HDE	0.291	0.321

we calculate the Root Mean Square Error for each instance and average all the errors to obtain a confidence measurement. In a secondary stage we used a test set consisting of 3600 new examples which have not been seen before by the algorithm, with parameters generated randomly. Both experiments were similar in precision. Here we present the average of results obtained in the two stages. To predict the parameters in the observational galactic spectra we used the ensembles generated in the previous step.

The prediction time required for the traditional ensembles is about 1.1 seconds per spectra and the time consumed by the HDE is about 1.7 seconds per spectra. Although the time taken by HDE is higher than the time taken by the normal ensembles, the errors in the profiles and the reconstructed spectra using the HDE prediction are better than the best obtained using the traditional ensembles prediction, as we can see in Figure 4. As expected, the RMS-error is higher for the observational spectra than for the synthetic spectra, because the profiles of the synthetic lines are better defined and there are no emission lines inside the absorption lines as in the observational spectra.

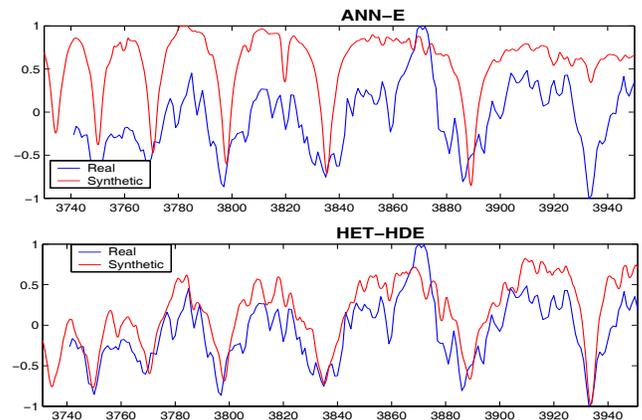


Figure 4: Fitting in an observational spectrum

Conclusions and Future Work

We have introduced a novel method for constructing ensembles using the *divide and conquer* principle, which is specially effective when there are several distinct subgroups that share a behavior.

To determine the age of stellar populations in synthetic and observational galactic spectra, the accuracy of the HDE

is better than the one obtained using traditional ensemble methods.

We presented an efficient machine learning method for identification of ages in galactic spectra, which is robust in the presence of noise, and in addition is insensitive to strong emission lines.

For future work we will experiment with different algorithms for the classifier and the divider modules and we will examine the behavior of the HDE using different data sets.

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