

Predicting physical properties of Lyman alpha forest with deep learning

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Lyman alpha (Lya) forest records how light of a quasar interacts with neutral hydrogen atoms in intergalactic medium (IGM). These systems provide much information to cosmological studies. However, the conventional line detection methods are restricted to spectra with high signal-to-noise ratio (SNR) and are time-consuming to apply in analysis. In this work, we apply convolution neural networks (CNN) to scan through quasar spectra for Ly α absorption line detection and simultaneously predict physical properties such as column densities (logNHI), redshifts (z), and Doppler widths (b) of the identified Ly α systems. For systems with logNHI ≤ 18.0 , our CNN, trained with simulated mock spectra, provides a state-of-art prediction in physical properties. We apply this model to build a catalogue of Ly α absorption systems for 300 quasar spectra from High Resolution Echelle Spectrometer (HIRES).

(logNHI, z, b) at the central pixel in a segmentation to train the CNN (as shown below).



using PYIGM¹ with a column density distribution function (CDDF) physical properties (logNHI, b) from simulated spectra are used in (2012). the training process. Additional SNR is added and this stabilises the CNN predictions when applying to spectra with different SNR (figure shown as below).

1: <u>https://pyigm.readthedocs.io/en/latest/#</u>



Abstract

Methodology — Our CNN model is trained with several segments of Results with mock spectra — The CNN model identifies Ly α abso- Results — We match the catalogue of Rudie+12 with our quasar spectra. We scan through a spectrum and use the labels rption lines with logNHI < 18.0 from a mock quasar spectrum with an CNN model with a probability threshold of 0.5 using a additional SNR of 10.0. An accuracy of over 0.85 for the identification is redshift difference of 0.0001 (~6.52 pixels). The CNN shows reached when applying a probability threshold of 0.5 to the predictions. a comparable result to the one using simulated mock spectra The prediction of logNHI and b has a root-mean-square error (RMSE) of 0.243 cm^{-2} and 5.35 km/s, respectively.



Results — Our CNN model trained with simulated mock spectra picked up corresponding Ly α systems listed in Rudie+12 (examples shown as below).



with a RMSE of 0.251 cm⁻² and 8.89 km/s (5.37 if excluding systems with b>45 km/s) for logNHI and b, respectively.



Observational spectra — Spectroscopic data from the High Resolution The above figure combined with the histogram of logNHI Mock spectra (for training) — Mock spectra for training is created Echelle Spectrometer (HIRES) has $R \sim 45,000$ (FWHM=7.0 km/s). We and b, the CNN model shows great consistency with the apply our CNN model to them and compare our identification and measurement in Rudie+12. The mis-prediction (>1 σ) in b from Prochaska et al. (2014) and FWHM=7.0 km/s. Labels of predictions with a catalogue of Lyα absorption systems from Rudie et al. occurs when a system has a broad feature (i.e. large b), our machine tends to under-estimate b for this kind of system.





<u>Summary & Future Plan</u> — The CNN model trained with simulated mock spectra shows a state-of-art predicting ability to observational data. This takes ~5 mins to generate a catalogue of Ly α forest for one quasar spectrum. We apply this technique to build a catalogue of 300 HIRES spectra. The paper of this work is in preparation. This technique can be easily adapted to other spectroscopic surveys. More importantly, with the generated Ly α forest catalogue, several studies of IGM can be carried out such as the thermal state of IGM. Follow-up works are in progress.