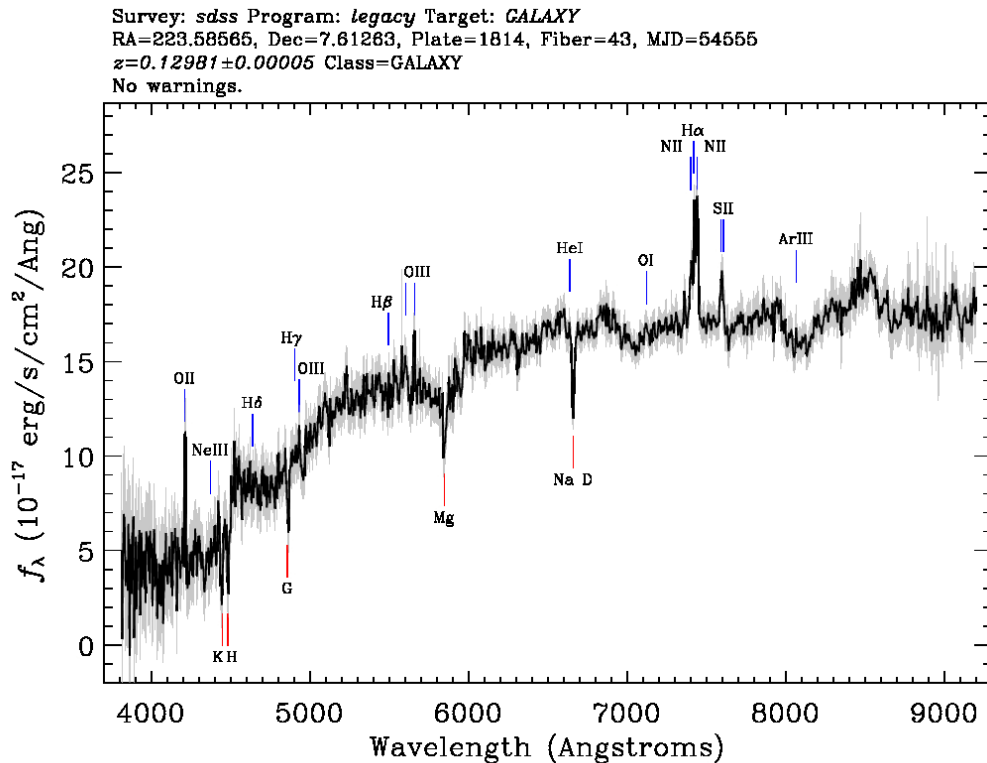


# Tutorial in plotting SDSS spectrum in Python (emission lines + absorbtion lines)

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## 1 Example from SDSS Webpage



```
[1]: import numpy as np
import matplotlib.pyplot as plt
import matplotlib.image as mpimg
from matplotlib.ticker import AutoMinorLocator
from astropy.io import fits
```

## 2 Prepare the Spectrum Data

```
[2]: # input the target name

gal_name = 'SDSS J145420.55+073645.4'

# Read the FITS data

spec_file = '/home/phylmf/Code/idl_ppxf_run/spec-1814-54555-0043.fits'
spec_hdu = fits.open(spec_file)

# load the SDSS spectrum

gal_spec = spec_hdu[1].data

# get the inferred spectroscopic redshift by SDSS team
# there is conflicts in z value between the FITS file and SDSS website
# for consistence with our IDL version tutorial, we apply the former one value

gal_z = float(spec_hdu[2].data['Z'])

gal_z_err = float(spec_hdu[2].data['Z_ERR'])

wave = 10 ** gal_spec['loglam'] / (1 + gal_z)

flux = gal_spec['flux']

ivar = gal_spec['ivar']

noise = ivar ** (- 0.5)

# red the galaxy image

gal_img = mpimg.imread("/home/phylmf/Code/idl_ppxf_run/py_ppxf/
↳spec-1814-54555-0043.jpeg")

# read the basic information:

gal_ra = spec_hdu[0].header['RA']
gal_dec = spec_hdu[0].header['DEC']

gal_plateid = spec_hdu[0].header['PLATEID']
gal_mjd = spec_hdu[0].header['MJD']
gal_fiberid = spec_hdu[0].header['FIBERID']

gal_redshift = '$z_{spec}$=' + str("{:.4f}".format(gal_z)) + '±' + str("{:.4f}".
↳format(gal_z_err))
```

### 3 Prepare Spectra Features Lines

- You can refer to the SDSS webpage about SSSP line index information via the link given below:
- ([https://data.sdss.org/datamodel/files/SSPP\\_REDUX/RERUN/PLATE4/output/param/ssppOut\\_lineinfo](https://data.sdss.org/datamodel/files/SSPP_REDUX/RERUN/PLATE4/output/param/ssppOut_lineinfo))
- SDSS DR6 line table:
- (<https://classic.sdss.org/dr6/algorithms/linestable.php>)
- UV/Opt. Line table with physical explanation(recommended!):
- (<http://astronomy.nmsu.edu/drewski/tableofemissionlines.html>)
- Anyway, you can always refer to Xueguang Zhang's book:
  - [O II]  $\lambda 3727, 3729\text{\AA}$  双线, 但是在很多的星系光谱中, 由于双线间的波长间隔过于狭窄, 只能分辨出一个中心波长在  $3728\text{\AA}$  的单峰的发射线
  - H $\theta$   $\lambda 3798.976\text{\AA}$ , 常表现为吸收线的特征, 如果是发射线, 一般都较弱。
  - H $\eta$   $\lambda 3836.47\text{\AA}$ , 常表现为吸收线的特征, 如果是发射线, 一般都较弱。
  - Ca II H & K  $\lambda 3934.777, 3969.588\text{\AA}$ , 显著的吸收线特征。
  - 4000 $\text{\AA}$  break, 明显的吸收特征, 出现在年老的星系光谱中。
  - H $\delta$   $\lambda 4102.89\text{\AA}$ , 常表现为吸收线的特征, 如果是发射线, 一般都较弱。
  - G band  $\lambda 4305.61$ , 星系光谱中常见的吸收特征。
  - H $\gamma$   $\lambda 4341.68\text{\AA}$ , 常见的吸收线特征, 有时也是能分辨的较强的发射线特征。
  - [O III]  $\lambda 4364.436\text{\AA}$ , 常见的强的 [O III] 发射线。
  - He II  $\lambda 4687\text{\AA}$ , 常见的较强的 He II 发射线。
  - H $\beta$   $\lambda 4862.68\text{\AA}$ , 常见较强的发射线
  - [O III]  $\lambda 4960.295, 5008.24\text{\AA}$ , 几乎是光学波段最强的窄发射线, 而且是双线。
  - Mg I b  $\lambda 5176.7\text{\AA}$ , 常见的 Mg I 吸收线。
  - Na I D  $\lambda 5895.6\text{\AA}$ , 常见的 Na I D 吸收线。
  - [O I]  $\lambda 6302.046, 6365.536\text{\AA}$ , 光学波段常见的 [O I] 发射线, 而且是双线。
  - [N II]  $\lambda 6549.86, 6585.27\text{\AA}$ , 光学波段常见的 [N II] 发射线, 而且是双线。
  - H $\alpha$   $\lambda 6564.61\text{\AA}$ , 光学波段最强 Balmer 发射线。
  - [S II]  $\lambda 6718.29, 6732.67\text{\AA}$ , 光学波段常见的 [S II] 发射线, 而且是双线。
  - Ca II  $\lambda 8498, 8542, 8662\text{\AA}$ , 靠近红外波段的 Ca II 吸收线, 而且是三线特征。

```
[3]: # emission lines

el_name = [r'$O_{\mathrm{II}}$',
            r'$Ne_{\mathrm{III}}$',
            r'$H_{\mathrm{\delta}}$',
            r'$H_{\mathrm{\gamma}}$',
            r'$O_{\mathrm{III},4363}$',
            r'$H_{\mathrm{\beta}}$',
            r'$O_{\mathrm{III},4959,5007}$',
            r'$He_{\mathrm{I}}$',
```

```

        r'$H_{\alpha}$',
        r'$N_{\mathrm{II},6548,6583}}$',
        r'$S_{\mathrm{II},6716,6731}}$',
        r'$Ar_{\mathrm{III}}$']

el_color = 'blue'

el_wave = {}
el_wave[el_name[0]] = [3728]
el_wave[el_name[1]] = [3868.760]
el_wave[el_name[2]] = [4101.742]
el_wave[el_name[3]] = [4340.471]
el_wave[el_name[4]] = [4363.210]
el_wave[el_name[5]] = [4861.333]
el_wave[el_name[6]] = [4958.911, 5006.843]
el_wave[el_name[7]] = [5875.624]
el_wave[el_name[8]] = [6562.819]
el_wave[el_name[9]] = [6548.050, 6583.460]
el_wave[el_name[10]] = [6716.440, 6730.810]
el_wave[el_name[11]] = [7135.790]

# absorption line

al_name = [r'K', r'H', r'G', r'Mg', r'Na']

al_color = 'red'

al_wave = {}
al_wave[al_name[0]] = [3934.777]
al_wave[al_name[1]] = [3969.588]
al_wave[al_name[2]] = [4305.61]
al_wave[al_name[3]] = [5176.7]
al_wave[al_name[4]] = [5895.6]

```

## 4 Plot the Spectrum and Save as Figure File

```

[4]: # set up plot (symbols in LaTeX Math Mode)

fig, ax = plt.subplots(figsize = (10,6), dpi = 300)

text_kwargs = dict(horizontalalignment='left',
                    verticalalignment='top',
                    linespacing = 5)

ax.set_xlabel('$\lambda_{\mathrm{rest}}$ (\AA)', fontsize = 15)
ax.set_ylabel('$F_{\lambda} (10^{-17} \mathrm{erg/s/cm}^2/\mathrm{\AA})$', fontsize = 15)

```

```

ax.set_xlim(3300., 8000.)
ax.set_ylim(-2., 33)
ax.minorticks_on()

# generate minor axis

x_minor_locator = AutoMinorLocator(n = 10)
y_minor_locator = AutoMinorLocator(n = 5)

# apply minor axis

ax.xaxis.set_minor_locator(x_minor_locator)
ax.yaxis.set_minor_locator(y_minor_locator)

ax.tick_params(axis = 'both', which = 'minor', length = 5, width = 1, direction=
    ↪ 'in', color = 'black')
ax.tick_params(axis = 'both', which = 'major', length = 10, width = 1,
    ↪ direction = 'in', color = 'black')
ax.tick_params(axis = 'both', labelsizе = 15)

# plot spectrum

ax.plot(wave, flux, color = 'black', linewidth = 0.6)
ax.fill_between(wave, flux - noise, flux + noise, color = 'gray', alpha = 0.5)

# plot emission lineq label

band_width = 5 # set band width for masking and find peak value
line_offset = 2

for i in range(len(el_name)):
    for j in range(len(el_wave[el_name[i]])):
        # for specifying the NII and H_alpha line
        if i in [8, 9]:
            line_mask = (wave > el_wave[el_name[i]][j] - 100) & (wave <
    ↪ el_wave[el_name[i]][j] + 100)
            line_wave = wave[line_mask]
            line_flux = flux[line_mask]
            line_peak_mask = (line_flux == np.max(line_flux))
            line_pos = el_wave[el_name[i]][j]
            line_peak = np.max(line_flux)

            if i == 8:
                line_peak = line_peak + 1.5
            elif i in [6, 10]:
                line_mask = (wave > el_wave[el_name[i]][0] - 5) & (wave <
    ↪ el_wave[el_name[i]][1] + 5)

```

```

        line_wave = wave[line_mask]
        line_flux = flux[line_mask]
        line_peak_mask = (line_flux == np.max(line_flux))
        line_pos = el_wave[el_name[i]][j]
        line_peak = np.max(line_flux)

    else:
        line_mask = (wave > el_wave[el_name[i]][j] - band_width) & (wave <
↪ el_wave[el_name[i]][j] + band_width)
        line_wave = wave[line_mask]
        line_flux = flux[line_mask]
        line_peak_mask = (line_flux == np.max(line_flux))
        line_pos = el_wave[el_name[i]][j]
        line_peak = np.max(line_flux)

    ax.text(line_pos, line_peak + line_offset, '----', color = el_color, ha=
↪ 'center', va = 'center', fontsize = 12, rotation=90)

    ax.text(line_pos, line_peak + line_offset + 1.5, el_name[i], color =
↪ 'black', ha = 'center', va = 'center', fontsize = 8)

# plot absorption lineq label

band_width = 2 # set band width for masking and find peak value
line_offset = 2

for i in range(len(al_name)):
    for j in range(len(al_wave[al_name[i]])):
        line_mask = (wave > al_wave[al_name[i]][j] - band_width) & (wave <
↪ al_wave[al_name[i]][j] + band_width)
        line_wave = wave[line_mask]
        line_flux = flux[line_mask]
        line_peak_mask = (line_flux == np.max(line_flux))
        line_pos = al_wave[al_name[i]][j]
        line_peak = np.max(line_flux)
        ax.text(line_pos, line_peak - line_offset, '----', color = al_color, ha=
↪ 'center', va = 'center', fontsize = 12, rotation=90)

        ax.text(line_pos, line_peak - line_offset - 1.5, al_name[i], color =
↪ 'black', ha = 'center', va = 'center', fontsize = 8)

# add galaxy information

ax.text(0.6, 0.29, 'MJD=' + str(gal_mjd), transform = ax.transAxes,

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        fontsize = 10, color = 'black', **text_kwargs, zorder = 10)

ax.text(0.6, 0.25, 'Plate=' + str(gal_plateid), transform = ax.transAxes,
        fontsize = 10, color = 'black', **text_kwargs, zorder = 10)

ax.text(0.6, 0.21, 'Fiber=' + str(gal_fiberid), transform = ax.transAxes,
        fontsize = 10, color = 'black', **text_kwargs, zorder = 10)

ax.text(0.6, 0.17, 'Ra=' + str(gal_ra), transform = ax.transAxes,
        fontsize = 10, color = 'black', **text_kwargs, zorder = 10)

ax.text(0.6, 0.13, 'Dec=' + str(gal_dec), transform = ax.transAxes,
        fontsize = 10, color = 'black', **text_kwargs, zorder = 10)

ax.text(0.6, 0.09, gal_redshift, transform = ax.transAxes,
        fontsize = 10, color = 'black', **text_kwargs, zorder = 10)

# Plot galaxy image

ax_inset = plt.axes([.72, 0.145, .2, .2])
ax_inset.imshow(gal_img)
ax_inset.axis('off')
ax_inset.set_title(gal_name, fontsize = 6)

# Save figure

fig_name = 'spectrum_' + 'SDSS J145420.55+073645.4.'

fig.savefig(fig_name + 'png', bbox_inches = 'tight', format = 'png', dpi = 300)
fig.savefig(fig_name + 'pdf', bbox_inches = 'tight', format = 'pdf', dpi = 300)

```

