Temperature analysis

May 3, 2025

1 Temperature analysis

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1.1 Introduction

This notebook provides a comprehensive analysis of temperature measurements derived from spectral data. The workflow includes loading and preprocessing the data, identifying key spectral features, and calculating vibrational and rotational temperatures using both analytical methods and simulation-based fitting. The results are then visualized and compared to assess the accuracy and reliability of the methods.

1.2 Imports

[1]:
```
# %matplotlib ipympl # for interactive plots

import imageio.v2 as imageio
import matplotlib.pyplot as plt
import scipy.stats as stats
import numpy as np
import pandas as pd
import os
```

[2]:
```
import sys
sys.path.append("../src")
```

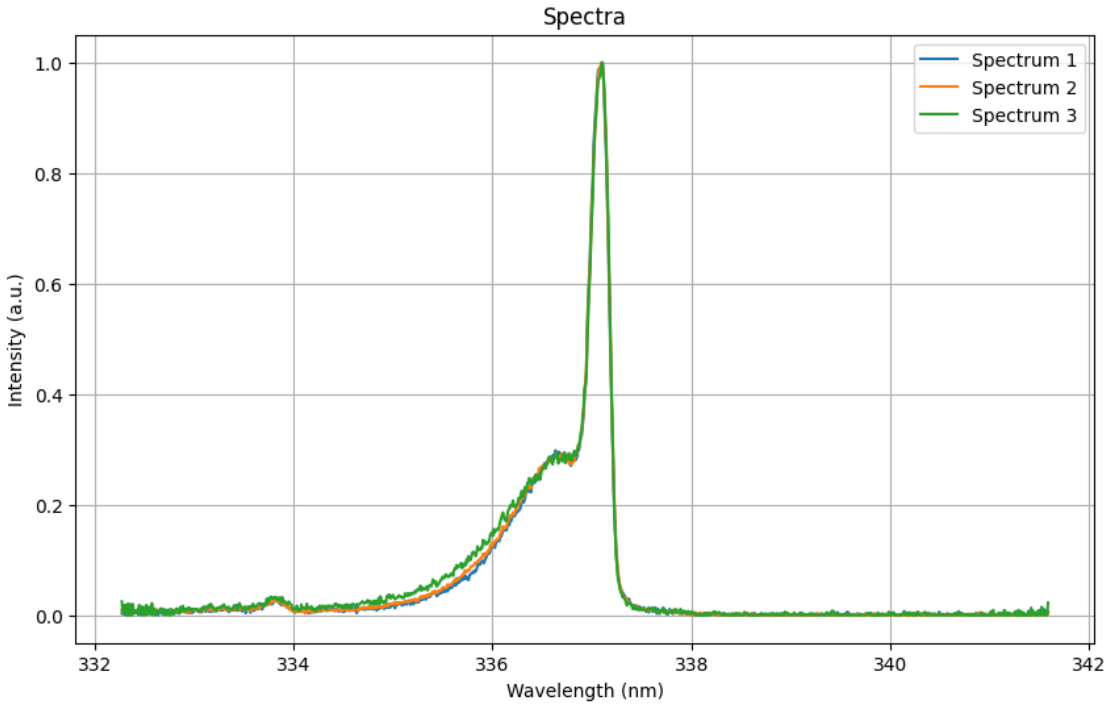
```
import xspectra.simulation as sim
import xspectra.utils as utils
import xspectra.visualization as vis
```

1.3 Load data

[3]:
```
folder_data = "data/temperature_analysis"
datas = np.array([np.loadtxt(os.path.join(folder_data, d), delimiter="\t") for␣
 ↪d in os.listdir(folder_data) if d.endswith(".txt")])
datas.shape
```

[3]:
```
(3, 1024, 2)
```

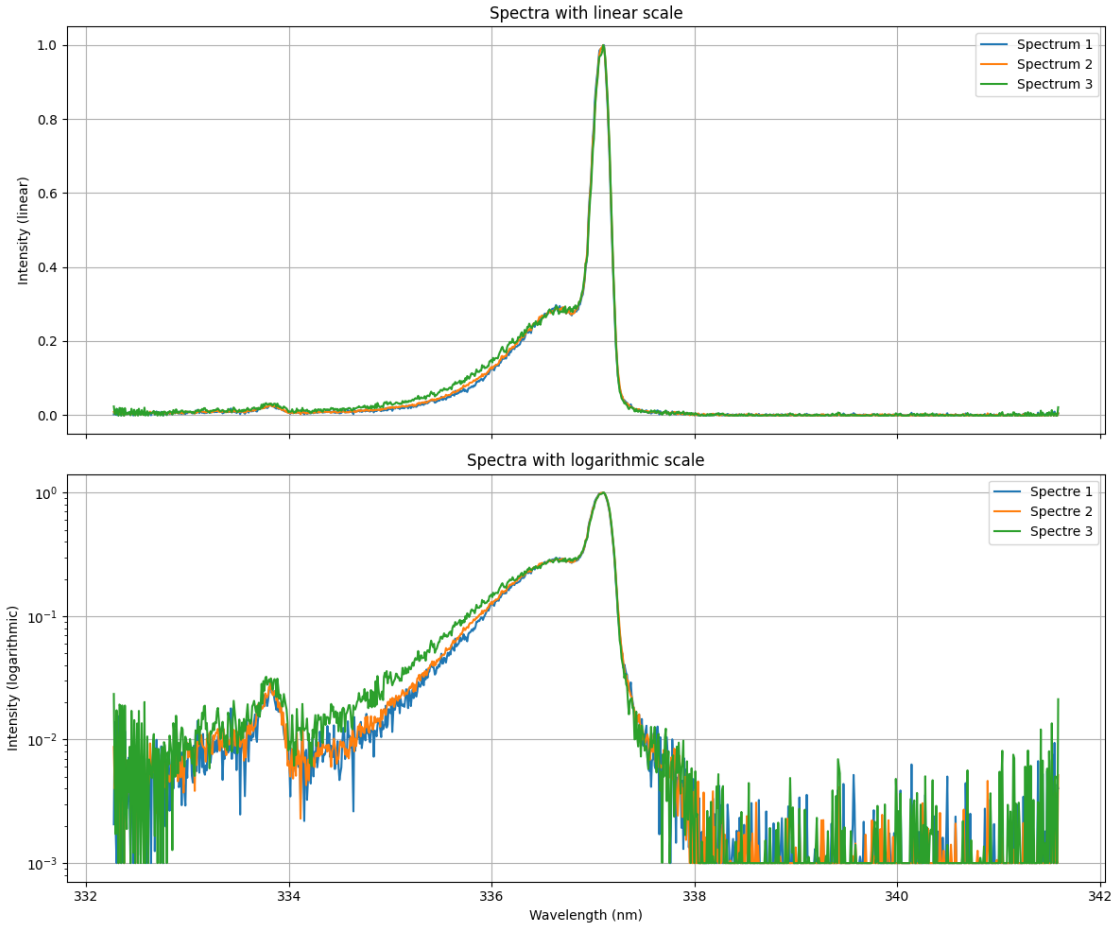
[4]:
```
plt.figure(figsize=(10, 6))

for i, data in enumerate(datas):
    plt.plot(data[:, 0], data[:, 1], label=f'Spectrum {i + 1}')

plt.xlabel('Wavelength (nm)')
plt.ylabel('Intensity (a.u.)')
plt.title('Spectra')
plt.legend()
plt.grid()
plt.show()
```



1.4 Background removing

[5]:
```
for i, data in enumerate(datas):
    datas[i] = utils.delete_background(data)
```

[6]:
```
fig, axs = plt.subplots(2, 1, figsize=(12, 10), sharex=True)

# Graphique avec échelle linéaire
for i, data in enumerate(datas):
    axs[0].plot(data[:, 0], data[:, 1], label=f'Spectrum {i + 1}')
axs[0].set_ylabel('Intensity (linear)')
axs[0].set_title('Spectra with linear scale')
axs[0].legend()
axs[0].grid()

# Graphique avec échelle logarithmique
epsilon = 1e-3
for i, data in enumerate(datas):
    axs[1].plot(data[:, 0], [x if x > epsilon else epsilon for x in data[:,␣
 ↪1]], label=f'Spectre {i + 1}')
axs[1].set_yscale('log')
axs[1].set_xlabel('Wavelength (nm)')
axs[1].set_ylabel('Intensity (logarithmic)')
axs[1].set_title('Spectra with logarithmic scale')
axs[1].legend()
axs[1].grid()

plt.tight_layout()
plt.show()
```



1.5 Calculation of the vibrational temperature

We identify the dominant peak at $337.5\,nm$ as well as secondary peaks. We take $334.44\,nm$ as the second reference.

Looking at the theoretical spectrum, we obtain that: - the dominant peak at $\lambda_1 = 337.5\,nm$ corresponds to the transition $C^3\Pi(\nu' = 0) \rightarrow B^3\Pi_g(\nu'' = 0)$ - the dominant peak at $\lambda_2 = 334.44\,nm$ corresponds to the transition $C^3\Pi(\nu' = 1) \rightarrow B^3\Pi_g(\nu'' = 1)$

Following the theoretical formulas developed in the simulation jupyter notebook, we have the following emission ratio:

$$r = \frac{\epsilon_1}{\epsilon_2} = \frac{n_1\nu_1}{n_2\nu_2} = \frac{g_{e1}(2J_1+1)e^{-\frac{T_{e1}}{kT_{el}} - \frac{G(\nu_1)}{kT_{vib}} - \frac{F(J_1)}{kT_{rot}}}}{g_{e2}(2J_2+1)e^{-\frac{T_{e2}}{kT_{el}} - \frac{G(\nu_2)}{kT_{vib}} - \frac{F(J_2)}{kT_{rot}}}} \frac{\nu_1}{\nu_2}$$

By eliminating the electronic degeneracies, which are equal, as well as the effect of rotations, we arrive at:

$$r_{12} = \frac{\nu_1}{\nu_2} \exp\left(\frac{T_{e2} - T_{e1}}{kT_{el}} + \frac{G(\nu_2) - G(\nu_1)}{kT_{vib}}\right)$$

Then, since we start from the same electrical energy level for both ($C^3\Pi$), we have $T_{e1} = T_{e2}$, hence:

$$r_{12} = \frac{\nu_1}{\nu_2} \exp\left(\frac{G(\nu_2) - G(\nu_1)}{kT_{vib}}\right)$$

Otherwise, we can use a third peak: $\lambda_3 = 331.735\,nm$ corresponding to the transition $C^3\Pi(\nu' = 2) \rightarrow B^3\Pi_g(\nu'' = 2)$ to find the two unknowns.

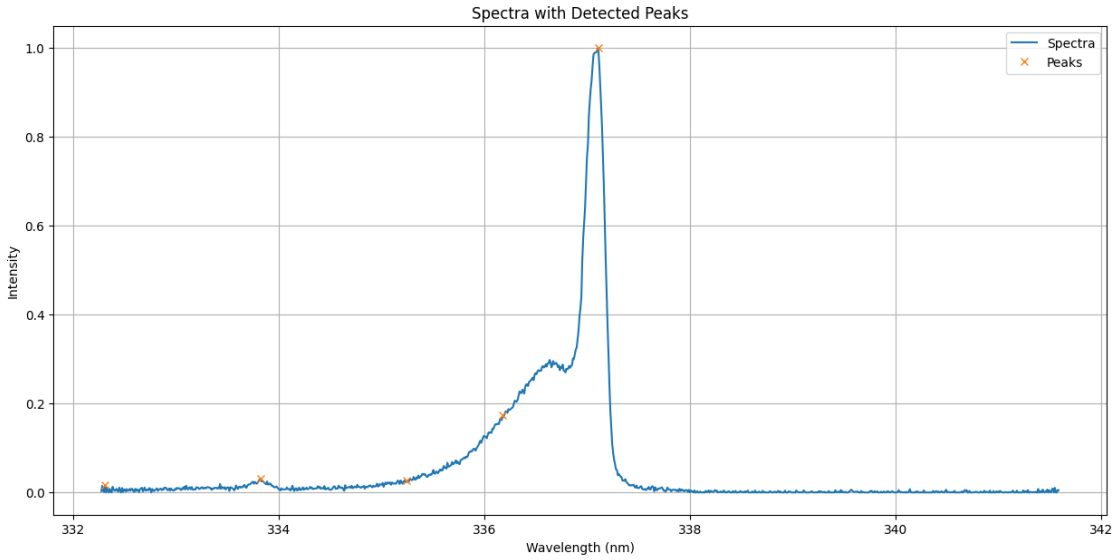
We don't have a clear enough spectrum to look at the rotational levels.

$$T_{vib} = \frac{G(\nu_2) - G(\nu_1)}{k \ln\left(r_{12} \cdot \frac{\lambda_1}{\lambda_2}\right)}$$

Using the wavelength values $\nu_1 = 337.5\,nm$ and $\nu_2 = 334.44\,nm$, as well as the ratio r_{12} calculated from the intensities of the corresponding peaks in the measured spectrum, we can determine T_{vib}.

Theoretically, $G(\nu_2 = 1) - G(\nu_1 = 0) = 5.973 \times 10^{-20} - 2.019 \times 10^{-20}\,J = 3.954 \times 10^{-20} J$

[7]:
```
from scipy.signal import find_peaks

peaks, _ = find_peaks(datas[0,:,1], height=0.01, distance=100)

# Afficher les résultats
plt.figure(figsize=(15, 7))
plt.plot(datas[0,:,0], datas[0,:,1], label="Spectra")
plt.plot(datas[0,peaks,0], datas[0,peaks,1], "x", label="Peaks")
plt.xlabel("Wavelength (nm)")
plt.ylabel("Intensity")
plt.title("Spectra with Detected Peaks")
plt.legend()
plt.grid()
plt.show()

# Afficher les longueurs d'onde des pics détectés
print(f"Index of the peaks : {peaks}")
print("Detected peaks at wavelengths:", datas[0,peaks,0])
```




```
Index of the peaks : [  4 169 324 427 529]
Detected peaks at wavelengths: [332.31041 333.8222  335.23876 336.17808
337.10659]
```

We can get automatically get the two index of the main transitions.

[8]:
```
i_primary, i_secondary = utils.find_index_primary_peaks(datas[0,:,0], datas[0,:
  ↪,1], height=0.01, distance=100)
print(f"Index of the primary peaks   : {i_primary:3d}   - Wavelengths :␣
  ↪{datas[0,i_primary,0]:3.2f}")
print(f"Index of the secondary peaks : {i_secondary:3d}   - Wavelengths :␣
  ↪{datas[0,i_secondary,0]:3.2f}")
```

```
Index of the primary peaks   : 529   - Wavelengths : 337.11
Index of the secondary peaks : 169   - Wavelengths : 333.82
```

We note the indices of the lines that interest us: - $i = 529$ for the line $\nu' = 0 \leftrightarrow \nu'' = 0$ - $i = 169$ for the line $\nu' = 1 \leftrightarrow \nu'' = 1$

[9]:
```
T_vib = [utils.calculate_temperature_single(data[:,1], i1=i_primary,␣
  ↪i2=i_secondary) for data in datas]

for i, t in enumerate(T_vib):
    print(f"Temperature for spectrum {i + 1}: {t:.0f} K")
```

```
Temperature for spectrum 1: 808 K
Temperature for spectrum 2: 825 K
Temperature for spectrum 3: 826 K
```

Here are the vibrational temperatures of the three spectra:

- Temperature for spectrum 1: 808 K

- Temperature for spectrum 2: 825 K
- Temperature for spectrum 3: 826 K

1.6 Calculation of the rotational temperature - By analysing the R branch

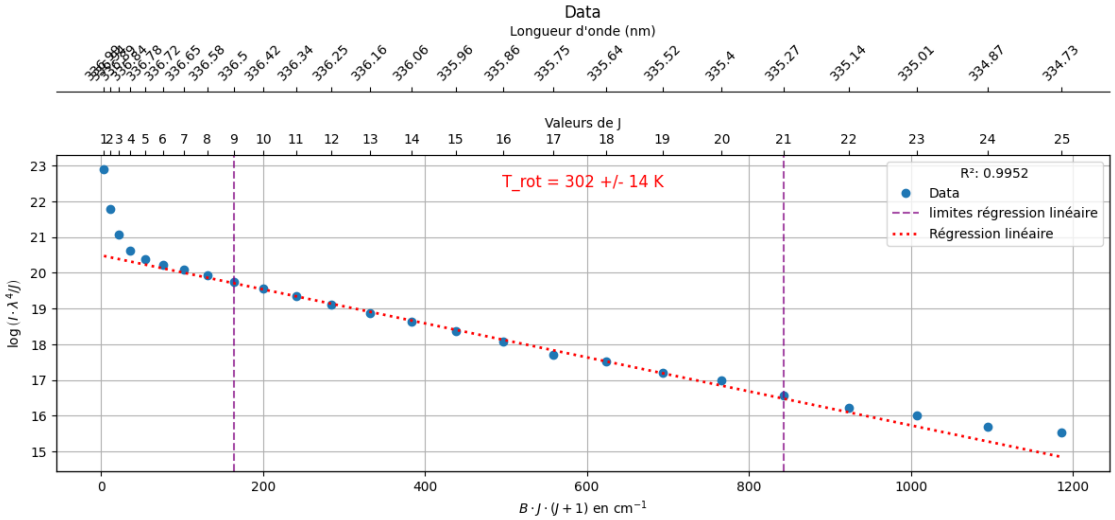
[10]:
```
T_rot_R_branch = [utils.compute_Trot_with_branch(data[:,0], data[:,1],␣
 ↪J_range=(8, 20), certainty=0.95) for data in datas]

for i, t in enumerate(T_rot_R_branch):
    print(f"Temperature for spectrum {i + 1}: {t[0]:.0f} +/- {t[1]:.0f} K")
```

```
Temperature for spectrum 1: 302 +/- 14 K
Temperature for spectrum 2: 312 +/- 8 K
Temperature for spectrum 3: 363 +/- 16 K
```

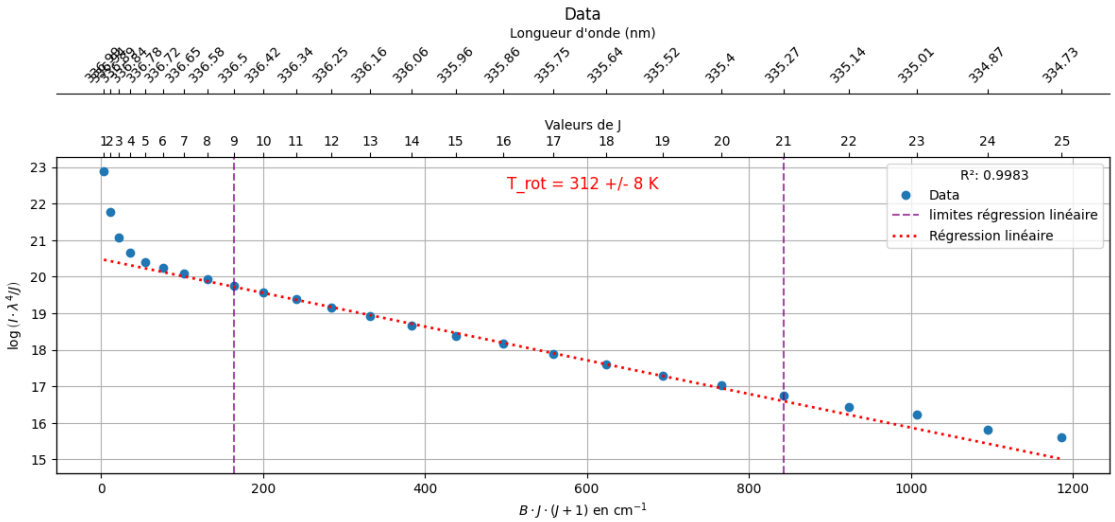
1.6.1 Verification

We van verify the regression by plotting it

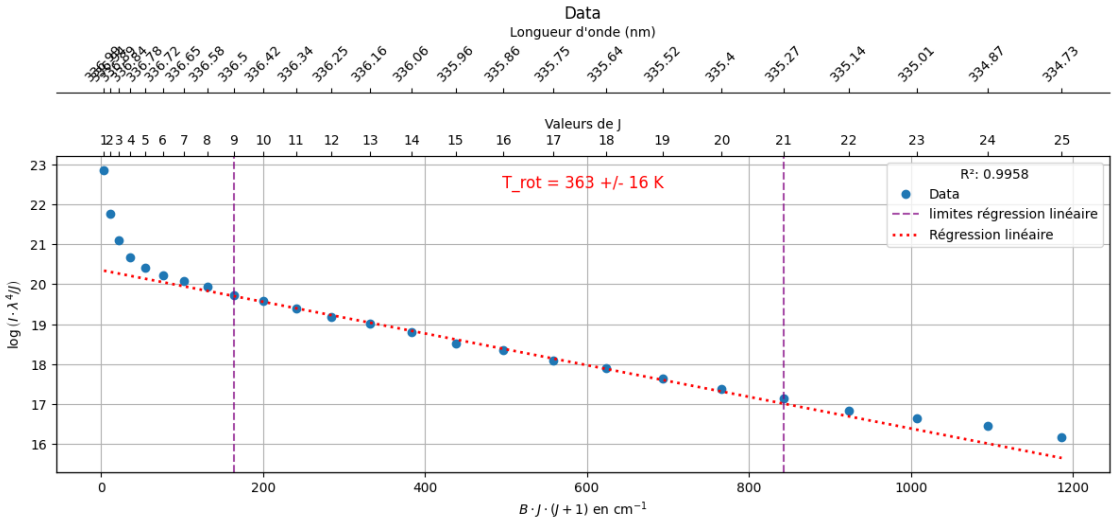
[11]:
```
# vis.show_result_calculation_Trot(datas[0,:,0], datas[0,:,1], J_range=(8, 20),␣
 ↪certainty=0.95, output_file="./res/Trot_R_branch.png")
vis.show_result_calculation_Trot(datas[0,:,0], datas[0,:,1], J_range=(8, 20),␣
 ↪certainty=0.95)
```



[12]:
```
vis.show_result_calculation_Trot(datas[1,:,0], datas[1,:,1], J_range=(8, 20),␣
 ↪certainty=0.95)
```



[13]:
```
vis.show_result_calculation_Trot(datas[2,:,0], datas[2,:,1], J_range=(8, 20),␣
↪certainty=0.95)
```



1.7 Calculation of the rotational temperature - By fitting with the simulation spectrum

[14]:
```
llims = (335.3, 338)
masks = (llims[0] <= datas[:, :, 0] ) & (datas[:, :, 0] <= llims[1])
```

[15]:
```
def find_wavelenth_deviation(wavelength, spectrum):
    MAXIMUM_SIMULATION_SPECTUM = 337.18 # nm
```

```
    return MAXIMUM_SIMULATION_SPECTUM-wavelength[spectrum.argmax()]
```

[16]:
```
decalage_vs_simulation = find_wavelenth_deviation(datas[0, :, 0], datas[0, :,␣
↪1])
decalage_vs_simulation
```

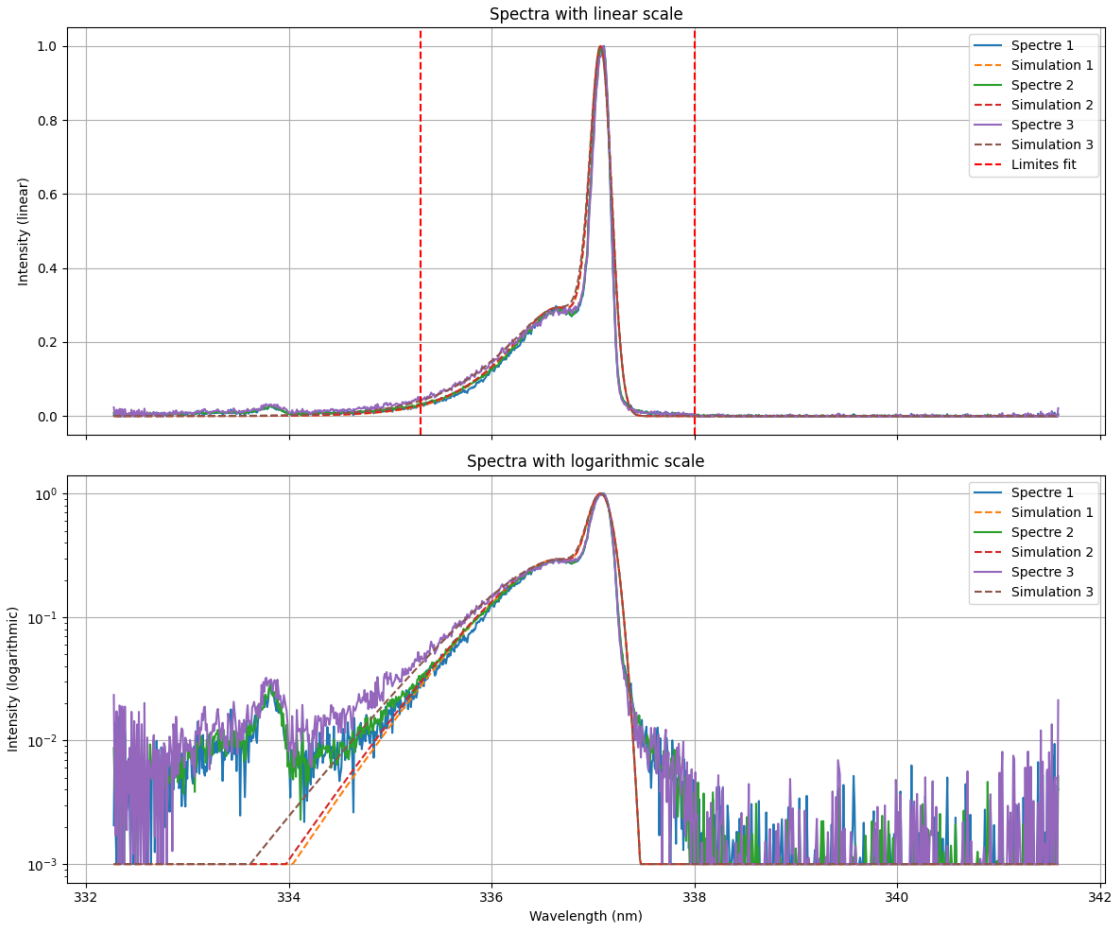
[16]:
```
np.float64(0.07341000000002396)
```

[17]:
```
decalage_vs_simulation = 0.1 # Or you can fix it manually (you can even make it␣
↪trainable for the fit - check bellow)
```

[18]:
```
# Calcul du spectre de simulation
T_rot_assumption = [t[0] for t in T_rot_R_branch]
simulation_spectrum = [sim.get_spectrum(datas[0, :, 0]+decalage_vs_simulation,␣
↪T_el=1_000, T_vib=T_vib[0], T_rot=T_rot_assumption[i], sigma_exp=0.1) for i␣
↪in range(3)]
```

[19]:
```
fig, axs = plt.subplots(2, 1, figsize=(12, 10), sharex=True)

# Graphique avec échelle linéaire
for i, data in enumerate(datas):
    axs[0].plot(data[:, 0], data[:, 1], label=f'Spectre {i + 1}')
    axs[0].plot(datas[0, :, 0], simulation_spectrum[i], label=f'Simulation␣
↪{i+1}', linestyle='--')

axs[0].axvline(x=llims[0], color='r', linestyle='--', label='Limites fit')
axs[0].axvline(x=llims[1], color='r', linestyle='--')

axs[0].set_ylabel('Intensity (linear)')
axs[0].set_title('Spectra with linear scale')
axs[0].legend()
axs[0].grid()

# Graphique avec échelle logarithmique
epsilon = 1e-3
for i, data in enumerate(datas):
    axs[1].plot(data[:, 0], [x if x > epsilon else epsilon for x in data[:,␣
↪1]], label=f'Spectre {i + 1}')
    axs[1].plot(datas[0, :, 0], [x if x > epsilon else epsilon for x in␣
↪simulation_spectrum[i]], label=f'Simulation {i+1}', linestyle='--')
axs[1].set_yscale('log')
axs[1].set_xlabel('Wavelength (nm)')
axs[1].set_ylabel('Intensity (logarithmic)')
axs[1].set_title('Spectra with logarithmic scale')
axs[1].legend()
axs[1].grid()
```

```
plt.tight_layout()
plt.savefig("./res/spectrum_simulation.png")
plt.show()
```



[20]:
```
T_rot_sim = np.zeros(len(datas))
elargissement_sim = np.zeros(len(datas))
decalage_sim = np.zeros(len(datas))

for i, data in enumerate(datas):
    s, t, elargissement, decalage = utils.get_best_fit(data[masks[0],0],␣
 ↪data[masks[0],1],
                                T_vib=T_vib[0],
                                elargissement=0.1,
                                w_decalage=0,
                               T_range=(100, 1_200),
                               elargissement_range=(0.05,0.12),
                               w_decalage_range=(-2, 2),
                               verbose=True,
```

```
                                nb_steps=5)
    print()

    T_rot_sim[i] = t
    elargissement_sim[i] = elargissement
    decalage_sim[i] = decalage
```

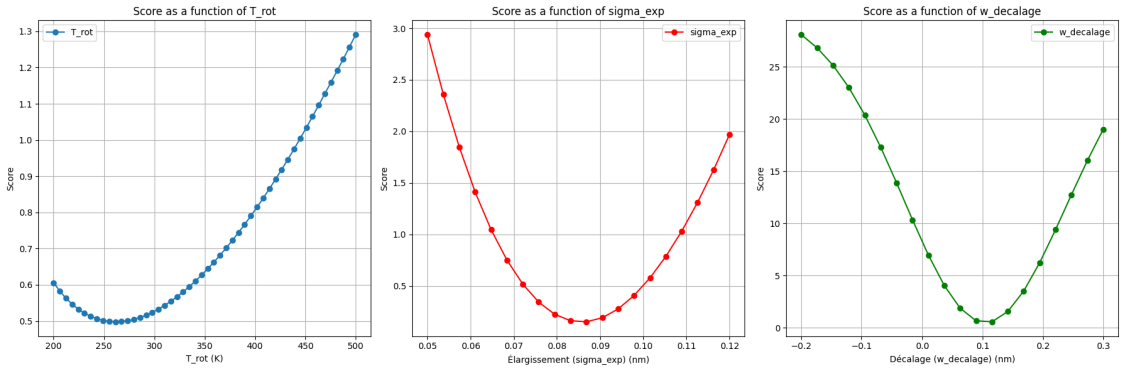
```
Iteration   1 | Score:    6.989 | Elargissement:   0.08 nm | T_rot:    570 K |
Décalage:   0.09 nm
Iteration   2 | Score:    8.080 | Elargissement:   0.09 nm | T_rot:    282 K |
Décalage:   0.10 nm
Iteration   3 | Score:    8.263 | Elargissement:   0.09 nm | T_rot:    257 K |
Décalage:   0.10 nm
Iteration   4 | Score:    8.278 | Elargissement:   0.09 nm | T_rot:    255 K |
Décalage:   0.10 nm
Iteration   5 | Score:    8.280 | Elargissement:   0.09 nm | T_rot:    255 K |
Décalage:   0.10 nm

Iteration   1 | Score:    6.614 | Elargissement:   0.08 nm | T_rot:    580 K |
Décalage:   0.09 nm
Iteration   2 | Score:    7.682 | Elargissement:   0.09 nm | T_rot:    297 K |
Décalage:   0.10 nm
Iteration   3 | Score:    7.870 | Elargissement:   0.09 nm | T_rot:    272 K |
Décalage:   0.10 nm
Iteration   4 | Score:    7.890 | Elargissement:   0.09 nm | T_rot:    270 K |
Décalage:   0.10 nm
Iteration   5 | Score:    7.891 | Elargissement:   0.09 nm | T_rot:    269 K |
Décalage:   0.10 nm

Iteration   1 | Score:    6.455 | Elargissement:   0.08 nm | T_rot:    608 K |
Décalage:   0.09 nm
Iteration   2 | Score:    7.464 | Elargissement:   0.08 nm | T_rot:    334 K |
Décalage:   0.10 nm
Iteration   3 | Score:    7.665 | Elargissement:   0.09 nm | T_rot:    309 K |
Décalage:   0.10 nm
Iteration   4 | Score:    7.694 | Elargissement:   0.09 nm | T_rot:    305 K |
Décalage:   0.10 nm
Iteration   5 | Score:    7.698 | Elargissement:   0.09 nm | T_rot:    305 K |
Décalage:   0.10 nm
```

1.8 Influence of different parameters on the fit

[21]:
```
# Calculer les scores pour chaque valeur de T_rot
T_rot_range = np.linspace(200, 500, 50)
sigma_exp_range = np.linspace(0.05, 0.12, 20)
w_decalage_range = np.linspace(-0.2, 0.3, 20)
```

```
filtered_wavelengths_target = datas[0, masks[0], 0]
filtered_spectrum_target = datas[0, masks[0], 1]

scores_T_rot = np.zeros_like(T_rot_range)
scores_sigma_exp = np.zeros_like(sigma_exp_range)
scores_w_decalage = np.zeros_like(w_decalage_range)

for idx, T_rot in enumerate(T_rot_range):
    scores_T_rot[idx] = utils.compute_score_fit(
        filtered_spectrum_target,
        sim.get_spectrum(filtered_wavelengths_target + decalage_vs_simulation,␣
 ↪T_el=T_vib[0], T_rot=T_rot, sigma_exp=0.1)
    )

for idx, sigma_exp in enumerate(sigma_exp_range):
    scores_sigma_exp[idx] = utils.compute_score_fit(
        filtered_spectrum_target,
        sim.get_spectrum(filtered_wavelengths_target + decalage_vs_simulation,␣
 ↪T_el=T_vib[0], T_rot=T_rot_sim[0], sigma_exp=sigma_exp)
    )

for idx, w_decalage in enumerate(w_decalage_range):
    scores_w_decalage[idx] = utils.compute_score_fit(
        filtered_spectrum_target,
        sim.get_spectrum(filtered_wavelengths_target + w_decalage,␣
 ↪T_el=T_vib[0], T_rot=T_rot_sim[0], sigma_exp=0.1)
    )

# Tracer le graphe
fig, axs = plt.subplots(1, 3, figsize=(18, 6))

axs[0].plot(T_rot_range, scores_T_rot, marker='o', label="T_rot")
axs[0].set_xlabel("T_rot (K)")
axs[0].set_ylabel("Score")
axs[0].set_title("Score as a function of T_rot")
axs[0].grid()
axs[0].legend()

axs[1].plot(sigma_exp_range, scores_sigma_exp, marker='o', label="sigma_exp",␣
 ↪color='red')
axs[1].set_xlabel("Élargissement (sigma_exp) (nm)")
axs[1].set_ylabel("Score")
axs[1].set_title("Score as a function of sigma_exp")
axs[1].grid()
axs[1].legend()
```

```
axs[2].plot(w_decalage_range, scores_w_decalage, marker='o',␣
 ↪label="w_decalage", color='green')
axs[2].set_xlabel("Décalage (w_decalage) (nm)")
axs[2].set_ylabel("Score")
axs[2].set_title("Score as a function of w_decalage")
axs[2].grid()
axs[2].legend()

plt.tight_layout()
plt.savefig("./res/scores_function.png", dpi=300, bbox_inches='tight')
plt.show()
```



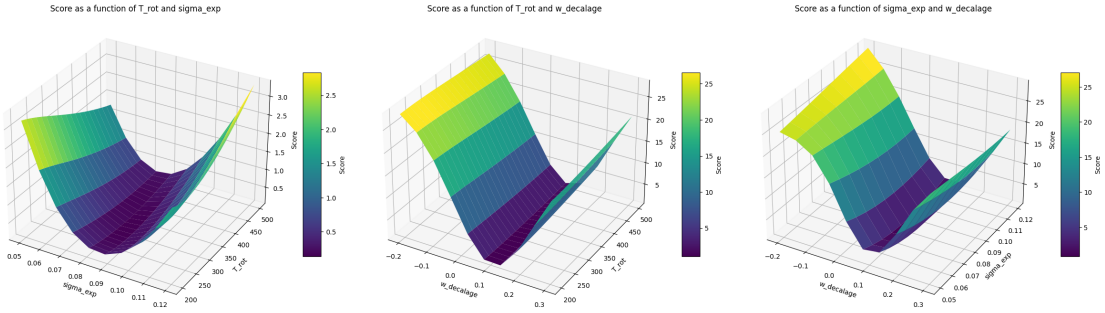
The functions are all convexe and that justifies the trichotomy algorithm used to compute the minimum.

We can even draw 3D graph for each combination.

[22]:
```
from itertools import combinations

# Définir les plages de paramètres
T_rot_range = np.linspace(200, 500, 20)
sigma_exp_range = np.linspace(0.05, 0.12, 10)
w_decalage_range = np.linspace(-0.2, 0.3, 10)

# Paramètres à combiner
params = {
    "T_rot": T_rot_range,
    "sigma_exp": sigma_exp_range,
    "w_decalage": w_decalage_range
}

param_names = list(params.keys())
```

[23]:
```
# Tracer les graphiques 3D alignés horizontalement
fig, axs = plt.subplots(1, 3, figsize=(24, 8), subplot_kw={'projection': '3d'})

# Générer toutes les combinaisons possibles de deux paramètres
for ax, (param1_name, param2_name) in zip(axs, combinations(param_names, 2)):
    param1_values = params[param1_name]
    param2_values = params[param2_name]

    # Initialiser une matrice pour stocker les scores
    scores = np.zeros((len(param1_values), len(param2_values)))

    # Calculer les scores pour chaque combinaison des deux paramètres
    for i, param1 in enumerate(param1_values):
        for j, param2 in enumerate(param2_values):
            # Définir les valeurs des paramètres en fonction de la combinaison␣
 ↪actuelle
            kwargs = {param1_name: param1, param2_name: param2, "T_el":␣
 ↪T_vib[0]}

            # Définir les valeurs par défaut pour les paramètres non inclus␣
 ↪dans la combinaison
            if param1_name != "T_rot" and param2_name != "T_rot":
                kwargs["T_rot"] = T_rot_sim[0]
            if param1_name != "sigma_exp" and param2_name != "sigma_exp":
                kwargs["sigma_exp"] = 0.1
            if param1_name != "w_decalage" and param2_name != "w_decalage":
                kwargs["w_decalage"] = decalage_vs_simulation

            # Calculer le score
            wavelengths = filtered_wavelengths_target + kwargs["w_decalage"]
            spectrum = sim.get_spectrum(wavelengths, T_el=kwargs["T_el"],␣
 ↪T_rot=kwargs["T_rot"], sigma_exp=kwargs["sigma_exp"])
            scores[i, j] = utils.compute_score_fit(filtered_spectrum_target,␣
 ↪spectrum)

    # Créer les grilles pour les paramètres
    param1_grid, param2_grid = np.meshgrid(param1_values, param2_values)

    # Tracer la surface
    surf = ax.plot_surface(param2_grid, param1_grid, scores.T, cmap='viridis',␣
 ↪edgecolor='none')

    # Ajouter une barre de couleur
    fig.colorbar(surf, ax=ax, shrink=0.5, aspect=10, label='Score')

    # Ajouter des labels
    ax.set_xlabel(param2_name)
```

```
    ax.set_ylabel(param1_name)
    ax.set_zlabel('Score')
    ax.set_title(f'Score as a function of {param1_name} and {param2_name}')

plt.tight_layout()
plt.show()
```



2 Conclusion

Spectre	Température vibrationnelle (K)	Température rotationnelle (K) - méthode analytique	Température rotationnelle (K) - fit simulation
Spectre 1	808	302 ±14	262
Spectre 2	825	312 ± 8	273
Spectre 3	826	363 ± 16	303

[24]:
```
data = {
    'Spectre': ['Spectre 1', 'Spectre 2', 'Spectre 3'],
    'T_vib (K)': T_vib,
    'T_rot (K) - R branch': [f"{t[0]:.0f} +/- {t[1]:.0f}" for t in␣
 ↪T_rot_R_branch],
    'T_rot (K) - fit simulation': [f"{t:.0f}" for t in T_rot_sim],
    'Broadening (nm)': [f"{e:.3f}" for e in elargissement_sim],
    'WDeviation (nm)': [f"{d:.3f}" for d in decalage_sim],
}

df = pd.DataFrame(data)
# print(df.to_markdown(index=False, tablefmt="grid"))
df
```

[24]:

	Spectre	T_vib (K)	T_rot (K) - R branch	T_rot (K) - fit simulation	Broadening (nm)	WDeviation (nm)
0	Spectre 1	808.232528	302 +/- 14	255	0.086	0.104
1	Spectre 2	825.049796	312 +/- 8	269	0.086	0.102
2	Spectre 3	825.847336	363 +/- 16	305	0.085	0.099

[ ]: