

# Title

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# 1 Introduction

Bla bla bla.

# 2 Theory

Bla bla bla.

# 3 Results

Our final result of the fit is shown in figure ??.

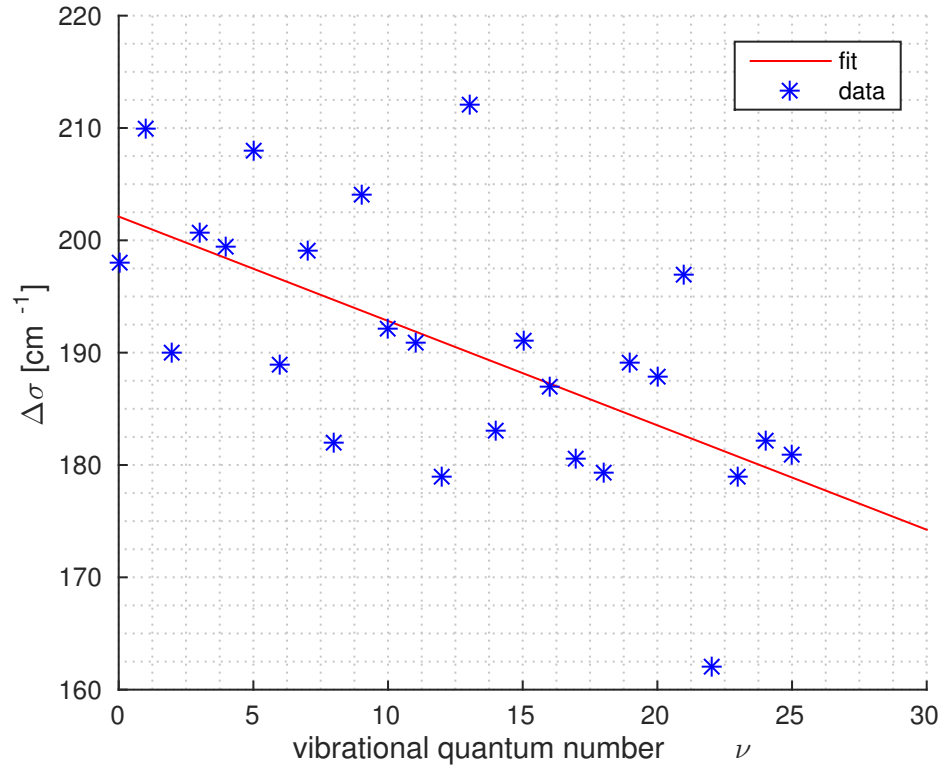


Figure 1: Fitting the data to the to the linear equation  $\Delta\sigma = \omega_e - 2\omega_e x_e(1 + \nu)$  with inserted peaks.  $R^2 = 0.3828$

$\omega_e = 214.50cm^{-1}$  and  $\omega_e x_e = 0.614cm^{-1}$  was retrieved from NiST[?] is the reference value or theoretical value. The result we got from the last fit is  $\omega_e 202.5803^{-1}$  which is a 5.557% deviation from the theoretical value,  $\omega_e x_e = 0.46463cm^{-1}$  which is a 24.3277% deviation and  $D_e = 22081.5347cm^{-1}$  that is a 17.8751% deviation.

You can see the peaks that are used in figure ?? at table ??.

Vibrational quantum number $\nu$	$\Delta\sigma$ [ $10^4cm^{-1}$ ]
4	1.8828
11	1.7455
18	1.6131
20	1.5763
25	1.4855

Table 1: Inserted peaks that are used in figure ??

## 4 Discussion

## References

- [1] ChemeDDL.org <http://www.chemeddl.org/alfresco/service/api/node/content/workspace/SpacesStore/8a367d14-7f6d-4fef-a7df-eb19579da11b/IodineSpectrum.pdf?guest=true>
- [2] NiST <http://webbook.nist.gov/cgi/cbook.cgi?ID=C7553562&Units=SI&Mask=1000>