

Fine Structure Corrections in Optical Emission Spectra of Sodium

Bhaskar Mookerji Charles Herder

8.13 Experimental Physics I
MIT Department of Physics

11 December 2007

Goal:

- Characterize corrections to Bohr model and Rydberg equation.
- Validate predictions of quantum mechanics for hydrogen-like atoms.

CAUTION



**ELECTRONS
AT PLAY**

1 Introduction

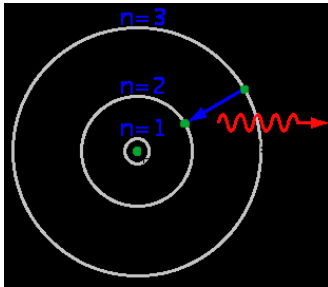
- QM of Hydrogenic Atoms
- Fine Structure Perturbation
- Sodium Doublet States

2 Experimental

- Czerny-Turner Monochromator (but I...)
- Peak resolution and their parameters

3 Results and Error Analysis

- 1800gvs/mm Mercury Calibration
- Error Accounting
- Measured Sodium Doublets
- Screening in s and d Angular Momentum States



$$H = \frac{\mathbf{p}^2}{2m} - \frac{Ze^2}{4\pi\epsilon_0} \frac{1}{r}. \quad (1)$$

$$\begin{aligned} E_n &= -hcR_\infty \frac{Z^2}{n^2} \\ &= -\frac{m_e e^4}{2(4\pi\epsilon_0)^2 \hbar^2} \frac{Z^2}{n^2} \quad (2) \end{aligned}$$

- Energy and electron orbitals are quantized.
- Electron transitions emit photons, causing discrete optical spectra.

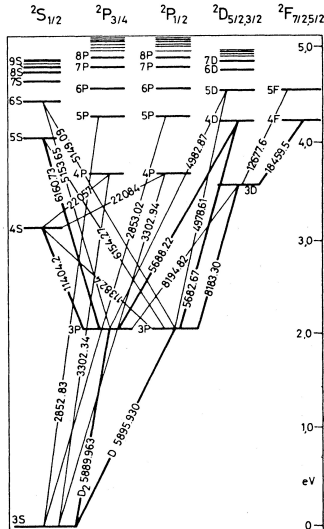
But wait!

- Electrons have angular momentum $L = 0, 1, \dots, n - 1$, $S = 1/2$
- Orbiting proton in electron frame sees a magnetic field \mathbf{B}

$$H_1 = -\boldsymbol{\mu} \cdot \mathbf{B} \rightarrow \left(\frac{e^2}{8\pi\epsilon_0} \right) \frac{1}{m^2 c^2 r^3} \mathbf{S} \cdot \mathbf{L} \quad (3)$$

- Now H , L^2 , S^2 , and $J = L + S$ commute!
- From first-order perturbation, angular momentum terms are no longer degenerate

$$E_{nj} = -\frac{E_n}{n^2} \left[1 + \frac{\alpha^2}{n^2} \left(\frac{n}{j + 1/2} - \frac{3}{4} \right) \right] \quad (4)$$



For states specified by $|nlm\rangle$,
transition probability related to

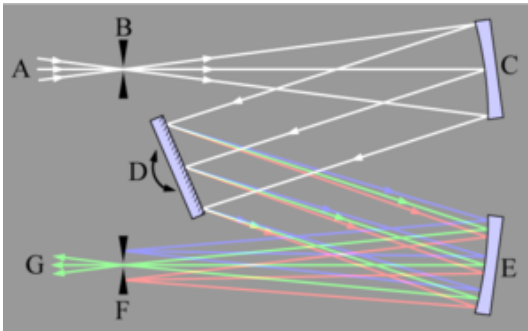
$$\langle n'l'm' | \mathbf{r} | nlm \rangle.$$

Only non-zero when,

$$\begin{aligned} \Delta l &= \pm 1 \\ \Delta m &= 0, \pm 1 \end{aligned} \quad (5)$$

‘Fine structure’: Energy
difference between P states

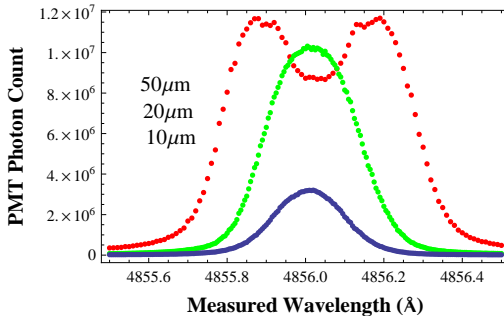
- Diffraction grating: 1800gvs/mm and 3600gvs/mm



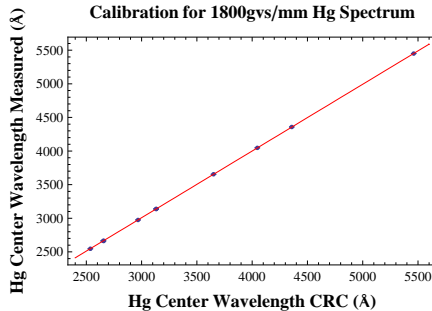
- (A) Emission source
- (B) Entrance slit
- (C) Collimating mirror
- (D) Diffraction grating
- (E) Focusing mirror
- (F) Exit slit
- (G) PMT (950V)

- Entrance/exit slit width, density of the diffraction grating, and source-entrance slit distance.

Slit Width Dependence for Uncalibrated H_β



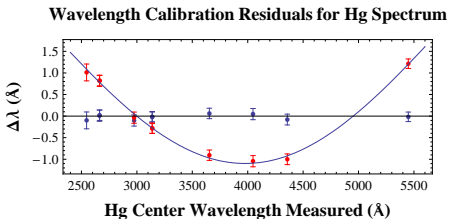
- Resolution:
 $\sigma_{1800} = 0.10 - 0.15 \text{\AA}$
 $\sigma_{3600} = 0.06 \text{\AA}$
- 100ms integration
- Slit: 4-10 μm
- 0.01 \AA step size
- 1800gvs/mm



$$\lambda_{\text{Predicted}} \sim 26.4612 + 0.9930x - 3.0940 \sin(7.4585 + 0.0009x)$$

Grating equation:

$$m\lambda = d(\sin \alpha + \sin \beta)$$



$$\sigma_{\text{Calib.}} \sim 0.06\text{\AA}$$

98% confidence level
from $\chi^2 = 1.6$.

Systematic error in our wavelength measurements are given by:

$$\sigma_{\lambda}^2 = \sigma_{\text{FWHM}}^2 + \sigma_{\text{Calib.}}^2 \quad (6)$$

Due to finite slit width, peaks approximated as Gaussians, therefore $\sigma_{\text{FWHM}} = \Delta\lambda_{\text{FWHM}}/\sqrt{2\ln 2}$

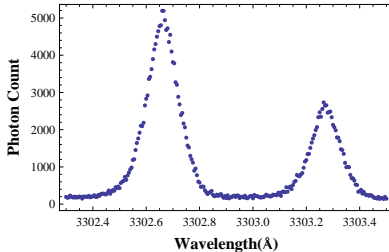
$$\sigma_{\text{FWHM}}^2 = \sigma_{\text{slit}}^2 + \sigma_{1800}^2 + \sigma_{\text{line width}}^2 + \sigma_{\text{doppler}}^2 \quad (7)$$

Doppler broadening for temperature $T = 2700\text{K}$ and $A = 11$:

$$\Delta\lambda = \lambda \cdot 10^{-6} \sqrt{\frac{2700}{11}} = \lambda \cdot 1.57 \cdot 10^{-6}$$

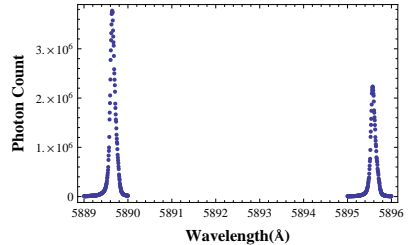
For 6160\AA , $\Delta\lambda = 0.001\text{\AA}$.

4P Doublet at 3303.26, 3302.64Å



4P Doublet at
 $\Delta\bar{\nu} = 5.60 \pm 0.94\text{cm}^{-1}$

3P Doublet at 5895.77, 5889.78Å



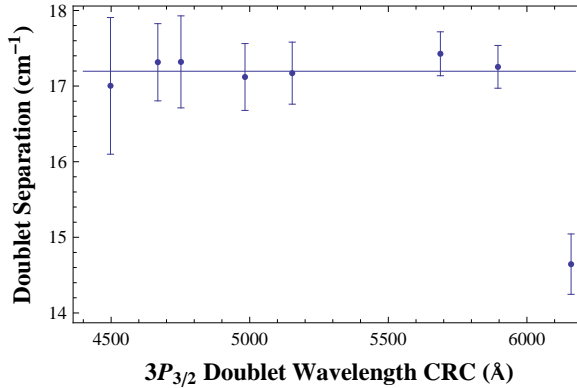
3P Doublet at
 $\Delta\bar{\nu} = 17.25 \pm 0.28\text{cm}^{-1}$

$$b = \text{Log}_{\frac{4}{3}}[\Delta\bar{\nu}_{3P}/\Delta\bar{\nu}_{4P}] = 3.88\text{Meas.} \quad 3.90\text{NIST}$$

Table: Classification of sodium fine structure (in air).

Transition	CRC	1800gvs/mm	
	$\lambda(\text{\AA})$	$\lambda(\text{\AA})$	$\Delta\bar{\nu}(\text{cm}^{-1})$
$5S \rightarrow 3P_{3/2}$	6160.75	6159.65 ± 0.25	14.65 ± 0.40
$\rightarrow 3P_{1/2}$	6154.23	6154.09 ± 0.20	
$3P_{3/2} \rightarrow 3S$	5895.92	5895.77 ± 0.13	17.25 ± 0.28
$3P_{1/2} \rightarrow 3S$	5889.95	5889.78 ± 0.13	
$4D \rightarrow 3P_{3/2}$	5688.21	5688.25 ± 0.12	17.43 ± 0.29
$\rightarrow 3P_{1/2}$	5682.63	5682.61 ± 0.12	
$6S \rightarrow 3P_{3/2}$	5153.40	5153.58 ± 0.16	17.17 ± 0.41
$\rightarrow 3P_{1/2}$	5148.84	5149.03 ± 0.14	
$5D \rightarrow 3P_{3/2}$	4982.81	4983.11 ± 0.15	17.12 ± 0.44
$\rightarrow 3P_{1/2}$	4978.54	4978.86 ± 0.16	
$7S \rightarrow 3P_{3/2}$	4751.82	4752.09 ± 0.22	17.32 ± 0.61
$\rightarrow 3P_{1/2}$	4749.94	4748.18 ± 0.20	
$6D \rightarrow 3P_{3/2}$	4668.56	4668.77 ± 0.16	17.32 ± 0.51
$\rightarrow 3P_{1/2}$	4664.81	4665.00 ± 0.15	
$7D \rightarrow 3P_{3/2}$	4497.66	4497.89 ± 0.27	17.00 ± 0.90
$\rightarrow 3P_{1/2}$	4494.18	4494.45 ± 0.31	
$4P_{3/2} \rightarrow 3S$	3302.98	3303.26 ± 0.14	5.60 ± 0.94
$4P_{1/2} \rightarrow 3S$	3302.37	3302.64 ± 0.14	
$5P_{3/2} \rightarrow 3S$	2853.01	—	—
$5P_{1/2} \rightarrow 3S$	2852.81	—	

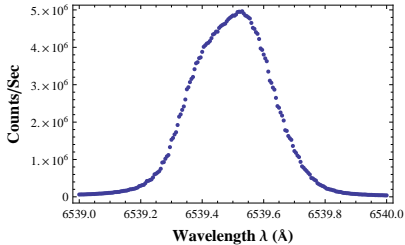
3P Doublet Separations in Sodium



$$\sigma_{\Delta\bar{\nu}}^2 = \frac{\sigma_{\lambda_1}^2}{\lambda_1^4} + \frac{\sigma_{\lambda_2}^2}{\lambda_2^4}$$

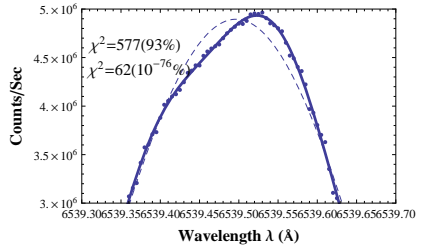
In case you were wondering...

Hydrogen Alpha Line



$$\lambda = 6539.39 \pm 0.003$$

Fitting Hydrogen Fine Structure

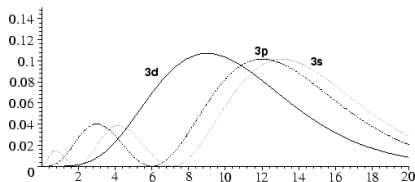


$$6539.54 \pm 0.003 \text{ Å}$$

$$\Delta\lambda_{\text{Meas.}} = 0.150 \pm 0.007 \text{ Å}$$

$$\Delta\lambda_{\text{Theory.}} = 0.1576 \pm 0.007 \text{ Å}$$

Angular momentum $L \rightarrow$ Electrons 'see' effective charge of nucleus $Z_{\text{eff}}e$



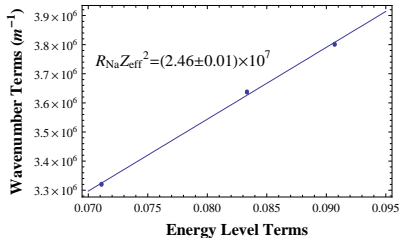
Assume final and initial states differ in Z_{eff} :

$$\frac{1}{\lambda} = -R_{\text{Na}} \left(\frac{Z_{\text{eff}_i}^2}{n_i^2} - \frac{Z_{\text{eff}_f}^2}{n_f^2} \right).$$

and equate common initial and final states:

$$\frac{1}{\lambda_2} \pm_d \frac{1}{\lambda_1} = Z_{\text{eff}}^2 R_{\text{Na}} \left(\frac{1}{n_f^2} - \frac{1}{n_i^2} \right) \quad (8)$$

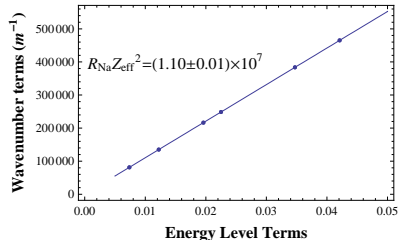
Combined Wavenumber vs. State Transitions



$$Z_s = 1.50 \pm 0.10$$

$$\lambda_{vac} = \lambda_{meas.} \cdot n_{air}$$

Combined Wavenumber vs. State Transitions



$$Z_d = 1.00 \pm 0.09$$

$$n_{air} = 1.0002739 \quad (9)$$

Typo in Mellisinos (1.14): From Slater's Rules,

$$Z_{eff} = 1.844 \sim 11 - 0.85 \cdot 8 - 2 = 2.2 \quad (10)$$

Table: Classification of sodium fine structure.

	CRC/NIST	Measured
$\Delta\tilde{\nu}_{3P}$	17.196cm^{-1}	$17.12 \pm 0.52 \pm 0.05\text{cm}^{-1}$
$\Delta\tilde{\nu}_{4P}$	5.59cm^{-1}	$5.60 \pm 0.94\text{cm}^{-1}$
Z_s	1.844	1.50 ± 0.10
Z_d	1	1.00 ± 0.09
IE_1	5.1391eV	$3.68 \pm 0.03\text{eV}$

- Correction to Bohr model from the spin-orbit perturbation, and angular momentum selection rules.
- Effect of screening for outer ‘shell’ electrons, and IE.
- Experimental confirmation of String Theory? *No*.

Acknowledgements

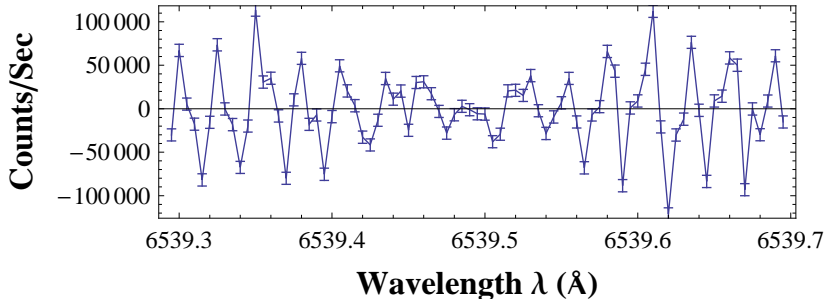
- Charles Herder: Physics appears to be easier when you're actually Course 6.
- JLab staff: For being helpful despite knowing more than we do.
- Glass: For attenuating the transmission of UV waves, thereby saving us from near certain death (by which I mean good UV data).
- Lab notebook bonfire after class? *Eh?*

8.13 T_c is finally over! :(



Extra

Fitting Hydrogen Fine Structure



$$^1I_{E_{Na}} = E_{3P-3S} + E_{\infty}^{3P} \quad (11)$$