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EXPERIMENTAL \( f \)-VALUE AND ISOTOPIC STRUCTURE FOR THE \( \text{Ni}^\text{i} \) LINE BLENDED WITH [\( \text{O} \text{i} \)] AT 6300 Å

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ABSTRACT

We have measured the oscillator strength of the \( \text{Ni}^\text{i} \) line at 6300.34 Å, which is known to be blended with the forbidden line [\( \text{O} \text{i} \)] \( \lambda \)6300 used for the determination of the oxygen abundance in cool stars. We also give wavelengths of the two isotopic line components of \( ^{58}\text{Ni} \) and \( ^{60}\text{Ni} \) derived from the asymmetric laboratory line profile. These two line components of \( \text{Ni}^\text{i} \) have to be considered when calculating a line profile of the 6300 Å feature observed in stellar and solar spectra. We also discuss the labeling of the energy levels involved in the \( \text{Ni}^\text{i} \) line since level mixing makes the theoretical predictions uncertain.

Subject headings: atomic data — stars: abundances — Sun: abundances

1. INTRODUCTION

Allende Prieto, Lambert, & Asplund (2001, hereafter APLA) have recently reviewed the various ways of determining the oxygen abundance in the Sun and other late-type stars and discussed the different problems connected with the spectral oxygen lines used. Stellar conditions limit the number of appropriate lines to a few allowed and forbidden lines of neutral oxygen ([\( \text{O} \text{i} \)] and [\( \text{O} \text{i} \)]) in the optical wavelength region and of molecular OH in the UV and IR regions. A frequently used method for determining the oxygen abundance in cool stars and the Sun is to use the forbidden [\( \text{O} \text{i} \)] line at 6300 Å, which is particularly studied and analyzed in the solar spectrum by APLA. However, this line is associated with problems since it is blended with an \( \text{Ni}^\text{i} \) line in the solar spectrum. This was pointed out earlier by Lambert (1978). Since the [\( \text{O} \text{i} \)] line (6300,31 Å) and the blending \( \text{Ni}^\text{i} \) line (6300,34 Å) appear as a totally unresolved feature in the solar spectrum, APLA constructed a spectral profile from laboratory data (wavelengths and \( f \)-values) and compared it with the observed solar feature. Three of four crucial atomic parameters are known to a satisfactory accuracy, viz., wavelength (Eriksson 1965) and \( f \)-value (Storey & Zeippen 2000) for the [\( \text{O} \text{i} \)] line and wavelength for the [\( \text{Ni} \text{i} \)] line (Litzén, Brault, & Thorne 1993, hereafter LBT). APLA used a three-dimensional time-dependent hydrodynamical model to simulate the solar surface and applied the same technique as previously used by Asplund et al. (2000) in the determination of the solar iron abundance. They used three free parameters, the continuum level, the oxygen abundance, and the product of \( f \)-value of the [\( \text{Ni} \text{i} \)] line times the nickel abundance, \( g_f(\text{Ni}) \), to match the predicted and observed profiles. By inserting the adopted solar abundance of nickel from Grevesse & Sauval (1998), APLA derived an “astrophysical” \( \log g_f \)-value of −2.31 for the [\( \text{Ni} \text{i} \)] line from the fitted value of \( g_f(\text{Ni}) \). We quote from the APLA paper: “The \( \log g_f \) for the [\( \text{Ni} \text{i} \)] line is uncertain. There are seemingly no laboratory measurements for this line.”

We have now determined the \( g_f \)-value of the [\( \text{Ni} \text{i} \)] line at 6300.34 Å by combining two-step laser-induced fluorescence (LIF) measurements of the radiative lifetime of the upper level with branching fraction measurements using Fourier transform spectroscopy. We have also fitted two isotopic line components (\( ^{58}\text{Ni} \) and \( ^{60}\text{Ni} \)) to the laboratory [\( \text{Ni} \text{i} \)] line and derived wavelengths and absolute intensities for both components. These two Ni isotopes account for 94% of the solar nickel abundance. We have reexamined the \( \log \) \( \epsilon \) composition of the upper energy level of the [\( \text{Ni} \text{i} \)] transition since it is severely mixed and has no clear \( \log \) \( \epsilon \) signature. The level was discussed and reassigned in the extensive work on the [\( \text{Ni} \text{i} \)] spectrum by LBT, and its identity has been further discussed by APLA. Adopting one or another of the “old” \((4d^4f^4)^P\) or the “new” \((4s^2S)^P\) assignment makes a difference in the \( f \)-value of a factor of 400 (a difference in \( \log g_f \) of 2.6) if we consult the database of R. L. Kurucz, which is often used in abundance work using the spectrum synthesis technique. Thus, an experimental value of the oscillator strength will help in understanding the real \( \log \) \( \epsilon \) composition of the energy level.

2. ATOMIC PHYSICS BACKGROUND

Since the [\( \text{Ni} \text{i} \)] line studied in this Letter has a great influence on the determination of the oxygen abundance, it deserves special attention. Because of the level mixing, there is a need for an experimental \( g_f \)-value as well as a detailed study of the isotopic composition of the line. The effect of level mixing that makes calculated oscillator strengths very uncertain is a frequent problem in complex spectra. Very drastic cases may occur in modeling stellar spectra in which a calculated spectral line, predicted by means of theoretical atomic data, totally disagrees with the observed feature.

To illustrate the level mixing in the present [\( \text{Ni} \text{i} \)] case, we have included a small part of the [\( \text{Ni} \text{i} \)] term diagram in Figure 1, showing the relevant energy levels involved in the discussion as well as in the measurements. The upper level of the line [\( \text{Ni} \text{i} \)] \( \lambda \)6300 is located at 50276 cm\(^{-1}\), and the lower level is \( y'D_1 \), which belongs to the odd parity \( 3d^84s4p \) configuration. In the first analysis of the [\( \text{Ni} \text{i} \)] spectrum, Russell (1929) assigned the 50276 level to \( e^3P_0 \) of the even parity \( 3d^94d \) configuration, which seemingly makes the transition to \( y'D_1 \), a “two-electron jump.” The appearance of such a transition can be explained by configuration interaction between \( 3d^84s4p \) and \( 3d^84p \), making the line a participant of a regular \( 4p-4d \) transition. Theoretical calculations by LBT confirmed such a level mixing between \( y'D_1 \) and \( z'D_1 \) belonging to \( 3d^84s4p \) and \( 3d^84p \), respectively.

In the National Institute of Standards and Technology compilation of iron group elements (Corliss & Sugar 1981), issued before the LBT work, the 50276 level has been given the old

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Fig. 1.—Partial level diagram of Ni I showing the levels and transitions discussed in this Letter. The two-step excitation in the lifetime measurement is indicated to the left.

label, $4d\,e'\text{P}$, suggested by Russell (1929). However, the label of another level located at 51457 cm$^{-1}$ was changed from $4d\,^1\text{S}_0$ to $4s\,^3\text{S}_1$. This change is supported by Kurucz’s calculations (2002) since the major eigenvector component is more than 90% of $4s\,^3\text{S}_1$. The Kurucz calculation also confirms the old label of the 50276 level since it was found to contain more than 95% $e'\text{P}_0$. Strangely enough, the Cowan code calculations performed by LBT gave completely opposite results concerning these levels. The 50276 level contained a 90% component of $4s\,^3\text{S}_1$, and the 51457 level was reassigned to the $4d$ configuration. LBT give the $4d$ labels in $jK$ coupling notation, which is clearly justified by the level structure (see Fig. 3 in their paper).

According to Kurucz’s calculations of log $gf$-values, both levels (50276 and 51457) have their strongest decay to $z'\text{P}_0^-$ (see Fig. 1). However, the transitions to $y\,^3\text{D}_0^+$ differ by a factor of 400 in the log $gf$-values, which are $-1.73$ for the transition from the 50276 level and $-4.38$ for the transition from the 51457 level. The second strongest decay from the 51457 level is to $z'\text{P}_0^-$ according to Kurucz, but the corresponding line from the 50276 level should be very weak. However, the latter is observed in the laboratory spectrum by LBT. Thus, the 50276 level obviously has a significant singlet character, which might be difficult to predict with sufficient accuracy in the calculations. Therefore, we have measured the lifetime and the log $gf$-values and analyzed the isotopic structure.

3. LIFETIME MEASUREMENTS

In the measurements of the radiative lifetime, the 50276 level was populated by applying a two-step pulsed laser excitation with $z'\text{P}_0^-$ as the intermediate level according to the scheme in Figure 1. The radiative lifetime was derived by a time-resolved observation of the fluorescence light released when the 50276 level decays to the lower odd levels. The experimental setup was similar to the one described and illustrated in a recent paper by Nilsson et al. (2000). Free nickel atoms were created by laser ablation and excited to the level investigated by laser pulses from two Nd:YAG laser-pumped dye lasers operating in the red spectral region. The wavelength of the first step excitation (3664 Å) was reached after frequency doubling and of the second step excitation (4811 Å) after Raman shifting in hydrogen gas. The two 10 ns pulses coincided in time and space during the interaction with the nickel atoms. Fluorescent light was detected using a monochromator and a fast photomultiplier. A digital transient recorder performed the data acquisition. The monochromator was set on one of the decay channels shown in Figure 1. Most of the recordings were taken on the strong 4811 Å line. An average of 1000 single decay events was typically necessary for obtaining a good signal-to-noise ratio in the exponential decay curves. Thirty curves were recorded, and the result for the lifetime of the 50276 level is $77 \pm 7$ ns. The uncertainty of the lifetime is a combination of statistical and, although carefully checked, possible systematic errors. The standard deviation of the 30 different measurements is less than 20% of the uncertainty.

4. BRANCHING FRACTIONS AND OSCILLATOR STRENGTHS

We have recorded the Ni I spectrum with the Lund UV Fourier transform spectrometer (FTS) and have also extracted spectra from the Kitt Peak FTS database, previously used in the extensive study of the Ni spectrum by LBT. All spectra are from hollow cathode lamps at various running conditions (pressure and DC). The spectra have been intensity-calibrated by means of branching ratios for internal argon lines (Whaling, Carle, & Pitt 1993), i.e., argon lines produced by the carrier gas in the nickel hollow cathode lamp. The branching fractions have been derived from the calibrated intensities of the spectral lines corresponding to the four decay channels indicated in Figure 1.

Since the lower levels in all transitions studied have short radiative lifetimes, there is no need for any corrections for self-absorption in the light source. According to predictions and observations, the transitions measured account for more than 99% of the decay from the 50276 level, and no residual branching fraction has to be considered. The total decay rate, given by the inverse value of the measured lifetime, has therefore been distributed among the four transitions in accordance with the experimental branching fractions. The results are given in Table 1. The uncertainties in the $f$-values are determined according to a procedure suggested by Sikström et al. (2002), and they include estimated errors in lifetime measurements, intensity measurements, and the instrumental response function.

5. DISCUSSION

There are some conclusions that can be drawn from Table 1. First, the astrophysical log $gf$-value of the $\lambda$6300 line derived by APLA in the fitting of the synthesized profile of the combined [O I] and Ni I lines to the observed solar line is closer to the present measurement than to the value given in the Kurucz database. The astrophysical log $gf$-value differs by 0.20 dex ($\approx 60\%$) from the new laboratory value, which has an uncertainty of about 0.05 dex (15%). Second, the triplet content of the 50276 level is overestimated in the calculations by Kurucz and perhaps underestimated in the
uncertainties in the center of gravity (c.g.) wavelength for the Ni \( \text{i} \) \( 6300 \) line from which the oxygen abundance in the Sun and cool stars can be determined. The line is blended with an Ni \( \text{i} \) \( 6300 \) line at 6300.34 \( \AA \) reveals an unresolved isotope structure. At least two isotopic line components, separated by 20 \( \text{mA} \), have to be included when calculating a line profile to match the \( \lambda 6300 \) feature in high-resolution stellar and solar spectra. The log \( gf \)-values for the individual components should then be weighted by the relative abundances of the two isotopes.

In this Letter, we report on the wavelengths and log \( gf \)-values for the two major isotopic components of the Ni \( \text{i} \) \( 6300 \) line. We
have measured the radiative lifetime of the upper level of the Ni i line using laser techniques and combined it with branching fractions to get the absolute transition probabilities. The branching fractions are derived from calibrated line intensities in a Fourier transform spectrum, which is also used to disentangle the isotopic structure. The new experimental log gf-value differs by 0.2 dex from the “astrophysical log gf-value” derived by APLA.

An appropriate labeling of the upper level of the Ni i line has previously been discussed by LBT. The additional information obtained in the present work, i.e., the lifetime and isotopic structure, supports the label suggested in that paper.

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