Influence of Inelastic Collisions with Hydrogen Atoms on the Formation of Al I and Si I Lines in Stellar Spectra

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Abstract—We have performed calculations by abandoning the assumption of local thermodynamic equilibrium (within the so-called non-LTE approach) for Al I and Si I with model atmospheres corresponding to stars of spectral types F–G–K with different metal abundances. To take into account inelastic collisions with hydrogen atoms, for the first time we have applied the cross sections calculated by Belyaev et al. using model approaches within the formalism of the Born–Oppenheimer quantum theory. We show that for Al I non-LTE leads to higher ionization (overionization) than in LTE in the spectral line formation region and to a weakening of spectral lines, which is consistent with earlier non-LTE studies. However, our results, especially for the subordinate lines, differ quantitatively from the results of predecessors. Owing to their large cross sections, the ion-pair production and mutual neutralization processes Al I(3s2) + H− ↔ Al II(3s3p) + H+ provide a close coupling of highly excited Al I levels with the Al II ground state, which causes the deviations from the equilibrium level population to decrease compared to the calculations where the collisions only with electrons are taken into account. For three moderately metal-deficient dwarf stars, the aluminum abundance has been determined from seven Al I lines in different models of their formation. Under the assumption of LTE and in non-LTE calculations including the collisions only with electrons, the Al I 3961 Å resonance line gives a systematically lower abundance than the mean abundance from the subordinate lines, by 0.25–0.45 dex. The difference for each star is removed by taking into account the collisions with hydrogen atoms, and the rms error of the abundance derived from all seven Al I lines decreases by a factor of 1.5–3 compared to the LTE analysis. We have calculated the non-LTE corrections to the abundance for six subordinate Al I lines as a function of the effective temperature (4500 K ≤ Teff ≤ 6500 K), surface gravity (3.0 ≤ log g ≤ 4.5), and metal abundance ([M/H] = 0, −1, −2, and −3). For Si I including the collisions with H I leads to the establishment of equilibrium populations in the spectral line formation region even in hot metal-deficient models and to vanishingly small departures from LTE in spectral lines.

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1. INTRODUCTION

The accuracy of calculating the profiles and equivalent widths of spectral lines when abandoning the assumption of local thermodynamic equilibrium (below referred to as non-LTE for short) depends on the completeness of allowance for all processes of the interaction of a given atom with the radiation field and the surrounding particles when calculating the statistical equilibrium (SE) and on the accuracy of atomic data: the energy levels, photoionization, photoexcitation, and inelastic collision cross sections. In the atmospheres of solar-type stars, the electron number density (N_e) is much lower than the number density of neutral hydrogen (H I) atoms; for example, in the solar atmosphere, N_e/N_H ≃ 10^{-4} in layers with log T_{5000} < −0.5. Therefore, the level excitation and the ion production can result from the collisions not only with electrons but also with H I atoms. The importance of including the latter when calculating the SE of a sodium atom was pointed out long ago by Gehren (1975). Analysis of the observed stellar spectra has shown that in the atmospheres of stars,
especially metal-deficient stars, apart from the collisions with electrons, there must exist an additional source of thermalization, which can be the collisions with H I atoms (for a review, see Lambert 1993; Holweger 1996; Mashonkina 2009). There are accurate data for inelastic collisions with H I in the literature only for a small number of atoms, and they have appeared mostly in recent years. H I atoms in the 1s2 S ground state were considered as impinging particles everywhere. Below, they will be referred to as H I for short. The atom being investigated (target) can be in both ground and excited states. Laboratory measurements were made for only one transition, namely Na I 3s–3p (Fleck et al. 1991). Detailed quantum-mechanical calculations were performed for the Na I (nl) + H I (Belyaev et al. 1999, 2010), Li I (nl) + H I (Belyaev and Barklem 2003), Mg I (nl) + H I (Belyaev et al. 2012), and He I (nl) + H I (Belyaev 2015) collisions. The Al I (nl) + H I and Si I (nl) + H I collisions were calculated by Belyaev (2013) and Belyaev et al. (2014), respectively, with the application of model approaches within the formalism of the Born–Oppenheimer quantum theory. Before the appearance of accurate data, the rates of inelastic collisions with H I were calculated using the approximate formulas derived by Steenbock and Holweger (1984) based on the theory of Drawin (1968). They are commonly called the Drawin rates in the literature. This approach continues to be used for most atoms, because there are no other, more accurate methods. It was established empirically that the Drawin rates require scaling with a factor $S_\text{H}$ whose value is different for different atoms (for references, see Mashonkina et al. 2011). For example, $S_\text{H} = 0.01$ for Na I and Ba II, but it varies from 0.1 to 30 for other atoms. Recognizing that the criticism of using the Drawin rates in non-LTE calculations is justified (see, e.g., Barklem et al. 2011), it should be said that this is just a forced necessity.

Many of the chemical elements in the spectra of F–G–K stars are observed in the lines of their neutral atoms, which account for only a small fraction in the total elemental abundance. These include Li I, Na I, Mg I, Al I, K I, Fe I, etc. For example, for aluminum with the ionization energy $\chi_{\text{ion}} = 6$ eV, $N(\text{Al I})/N(\text{Al}) < 5 \times 10^{-3}$ everywhere in the solar atmosphere. The number densities of such neutral atoms easily deviate from their equilibrium values when the intensity of ionizing radiation deviates from the Planck function, and this affects the spectral line formation. Indeed, calculations for many atoms, from Li I to Co I, showed significant non-LTE effects in lines and their growth with decreasing metal abundance (for a review, see Asplund 2005; Mashonkina 2014). The iron abundance $[\text{Fe/H}] = \log(N_{\text{Fe}}/N_\text{H}) - \log(N_{\text{Fe}}/N_\text{H})_\odot$ is commonly used as a metallicity indicator. At $[\text{Fe/H}] < -2$, the difference in the elemental abundances derived under the assumption of LTE and by abandoning LTE can reach large values. For example, for a star with an effective temperature $T_\text{eff} = 4680$ K, surface gravity $\log g = 1.23$, and $[\text{Fe/H}] = -3$, LTE leads to an abundance from Al I lines underestimated by 0.18 dex, from Ca I, on average, by 0.30 dex, and from Ti I by 0.33 dex, but to a sodium abundance overestimated by 0.23 dex. The number density of electrons in metal-deficient atmospheres is low, and they cannot serve as an efficient source of thermalization. Therefore, the magnitude of the non-LTE effects strongly depends on whether and how the inelastic collisions with H I atoms are taken into account. For example, Mashonkina (2013) determined the magnesium abundance for the star HD 122563 ($T_\text{eff}/\log g/[\text{Fe/H}] = 4600/1.60/-2.56$) under the assumption of LTE and in non-LTE calculations using the quantum-mechanical rates of Mg I + H I collisions from Barklem et al. (2012) and in non-LTE calculations including only the electron collisions. For the Mg I 4703 and 5528 Å lines, the non-LTE correction to the abundance, i.e., the difference between the non-LTE and LTE abundances $\Delta_{\text{NLTE}} = \log \varepsilon_{\text{NLTE}} - \log \varepsilon_{\text{LTE}}$, was $+0.27$ and $+0.16$ dex for the case of only electron collisions and decreased to 0.04 and $-0.04$ dex when the collisions with H I were taken into account.

This paper is devoted to analyzing the influence of collisions with H I atoms on the SE and lines of Al I and Si I. An accurate determination of the aluminum and silicon abundances for stars in a wide metallicity range is very important for testing the models of nucleosynthesis and evolution of the abundances of these elements in Galactic matter. Non-LTE calculations for Al I were performed previously by Baumuller and Gehren (1996), Andrievsky et al. (2008), and Menzhevtitski et al. (2012), and the importance of allowance for the departures from LTE in analyzing the stellar spectra was shown. The influence of non-LTE effects on the silicon abundance determination for the Sun and metal-deficient stars was studied by Shi et al. (2008, 2009). In all these papers, the collisions with H I were taken into account in accordance with Steenbock and Holweger (1984). For the first time we apply the results of accurate calculations from Belyaev (2013) and Belyaev et al. (2014) and investigate how this affects the abundance determination from individual Al I and Si I lines as well as the mean abundance and its error.

We describe the methods for calculating a theoretical spectrum of Al I and Si I in Section 2. Since the departures from LTE turned out to be small for Si I, the bulk of the paper is devoted to analyzing
the Al I lines. The results are presented in Section 3 for the Sun and Section 4 for stars. The non-LTE corrections to the abundance for six subordinate Al I lines as a function of stellar parameters are presented in Section 5. The conclusions and recommendations are formulated in Section 6.

2. THE METHODS AND CODES

2.1. The Al I Model Atom

We use the Al I model atom from Baumueller and Gehren (1996). It includes 59 Al I levels with a principal quantum number \( n \leq 15 \). The transition probabilities and photoionization cross sections were calculated within the OPACITY project. The electron-impact excitation is calculated using the formula of van Regemorter (1962) for allowed transitions, and the effective collision strength is assumed to be \( \Upsilon = 1 \) for forbidden ones. In contrast to Baumueller and Gehren (1996), who used the Drawin rates for collisions with H I, we apply the model rate constants calculated within the BO approach. Belyaev (2013) considered six lower Al I terms up to \( 4d^2D \) (with the excitation energy \( E_{\text{exc}} = 4.82 \text{ eV} \)) and all the transitions between them caused by collisions with H I as well as the ion-pair production from the ground and excited states and mutual neutralization \( \text{AlI}(nl) + \text{HI} \leftrightarrow \text{AlII}(3s^21S) + \text{H}^- \) (charged exchange reactions).

In Fig. 1 we compare the excitation rates due to the collisions with electrons and H I atoms as a function of the transition energy \( E_{ij} \) as well as the ion-pair production rates in the case of collisions with electrons and H I atoms as a function of the ionization energy of the initial atomic state \( \chi_i \). The collisions with electrons are more efficient than the collisions with H I atoms for bound–bound (b–b) transitions with a large energy separation between the levels \( (E_{ij} > 3 \text{ eV}) \), but the rates \( C_e \) and \( C_H \) are comparable in magnitude for \( E_{ij} < 2 \text{ eV} \). The collisions with H I produce an \( A^++\text{H}^- \) ion pair much more efficiently than the collisions with electrons ionize an atom if the atom is in a highly excited state with \( \chi_i < 3.5 \text{ eV} \).

2.2. The Si I–Si II Model Atom

The Si I–Si II model atom was constructed by Shi et al. (2008). It includes 132 singlet and triplet Si I terms, 41 Si II terms, and the Si III ground state. Just as for Al I, the radiative transition probabilities and photoionization cross sections were calculated within the OPACITY project, but there are no accurate data for collisions with electrons. Therefore, we used the same theoretical approximations as those for Al I. To take into account the collisions of neutral silicon atoms with H I, we apply the model rate constants calculated by Belyaev et al. (2014) within the BO approach. For Si I we considered 25 lower terms up to \( 4d^5F^0 \) (\( E_{\text{exc}} = 7.15 \text{ eV} \)) and all transitions between them as well as the charge exchange reactions. The excitation and ion-pair production rates due to the collisions with H I atoms are compared with the electron-impact excitation and ionization rates in Fig. 1. It should be noted that, in contrast to Al I, there exist many transitions with \( E_{ij} > 3 \text{ eV} \) for Si I, where \( C_H \approx C_e \). Otherwise, the conclusions for Si I and Al I coincide.

Thus, for both Al I and Si I the inelastic collisions with H I must play a no lesser role than the collisions with electrons in establishing the SE and, consequently, must be taken into account in non-LTE calculations.
2.3. Non-LTE Calculations

We solved the system of SE and radiative transfer equations in a specified model atmosphere using the DETAIL code developed by Butler and Giddings (1985) based on the accelerated \( \Lambda \)-iteration method. The level populations obtained by solving the SE equations (non-LTE) and based on the Boltzmann–Saha formulas (LTE), or rather their ratio \( b = n_{\text{NE}} / n_{\text{LTE}} \), which we will call below the b-factor, were then used to compute a synthetic spectrum via the SIU (Spectrum Investigation Utility) code (Reetz 1991).

Everywhere in this paper, we used model atmospheres from the MARCS\(^1\) database (Gustafsson et al. 2008). For given \( T_{\text{eff}} / \log g / [\text{Fe/H}] \), the models were obtained by interpolation using the MARCS algorithm.

Figure 2 illustrates the effect of including the collisions with H I atoms on the statistical equilibrium of Al I in the 5890/4.02/−0.78 model atmosphere. Irrespective of the calculation of collision rates, all Al I levels are underpopulated relative to the equilibrium values, but the non-LTE effects do not affect the Al II number density. The latter is easy to explain, because Al II is the dominant ionization stage. As was discussed by Baumueller and Gehren (1996) and shown by other earlier non-LTE studies, the main mechanism of departures from LTE for Al I is related to an excess of the mean radiation intensity beyond the threshold of ionization from the ground level (\( \lambda_{\text{thr}} \approx 2070 \) \( \text{Å} \)) over the Planck function, which leads to overionization, i.e., excess photoionization compared to the equilibrium one. As would be expected, including the collisions with H I atoms leads to a decrease in the departures from LTE for Al I levels. This is mainly because the coupling of highly excited Al I levels with the Al II ground state is strengthened.

For Si I the ionization energy \( \chi = 8.19 \) eV is higher than that for Al I, and the departures from LTE are smaller. As an illustration, our non-LTE calculations for Si I–Si II were performed with a higher temperature, \( T_{\text{eff}} = 6115 \) K, and a lower metal abundance, \( [\text{Fe/H}] = −1.24 \), than those for Al I. Figure 3 presents the b-factors of selected Si I levels. We do not show the Si II levels, because the technique of calculating the collision rates for them did not change. When the collisions only with electrons are taken into account, Si I shows a typical picture of overionization at line formation depths (\( \log \tau_{5000} < 0 \)). The 4s\(^3\)P\(^o\) level in the outermost layers constitutes an exception. Including the collisions with H I atoms removes the departures from LTE in layers deeper than \( \log \tau_{5000} = −2 \). In higher layers, the populations of the ground and low-excitation Si I states are higher than the equilibrium ones, i.e., there is an effect opposite to overionization.

The departures from LTE in the Si I lines observed in stellar spectra are small both when the collisions with H I atoms are taken into account and without them. For example, in the 6115/4.2/−1.24 model, the non-LTE corrections for the Si I 4102 \( \text{Å} \) (3p\(^1\)S\(_0\)−4s\(^3\)P\(^o\)), 3905 \( \text{Å} \) (3p\(^1\)S\(_0\)−4s\(^1\)P\(^o\)), 5772 \( \text{Å} \) (4s\(^1\)P\(_1\)−5p\(^1\)S\(_0\)), and 6155 \( \text{Å} \) (3p\(^3\)3D\(_2\)−5f\(^3\)G\(_4\)) lines are \( \Delta_{\text{NLTE}} = 0.12, 0.03, 0.00, \) and \( −0.03 \) dex, respectively, in the case where the collisions only with electrons are taken into account and become vanishingly small when the collisions with H I atoms are included: \( \Delta_{\text{NLTE}} = 0.00, 0.00, −0.01, \) and \( −0.02 \) dex. Therefore, our subsequent discussion will be devoted to analyzing the departures from LTE in Al I lines and the role of accurate data on the Al I + H I collision.

\(^1\)http://marcs.astro.uu.se

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Fig. 2. The b-factors of selected Al I levels in the 5890/4.02/−0.78 model atmosphere calculated by taking into account the collisions only with electrons (a) and including the collisions with H I atoms (b).
rates in improving the aluminum abundance in stellar atmospheres.

3. ANALYSIS OF Al I LINES IN THE SOLAR SPECTRUM

We analyze the spectrum of the Sun as a star (Kurucz 2005) and use the abundances derived from individual lines to determine the differential abundances for stars. All our calculations were made with the 5780/4.44/0.00 model atmosphere and microturbulence $\xi_t = 0.9$ km s$^{-1}$. Our analysis is based on the reproduction of the observed line profiles rather than equivalent widths. An example of the Al I 3961 Å line is shown in Fig. 4. For comparison with observations, the theoretical profile was convolved with the rotation profile and the radial–tangential profile of macroturbulent motions. The rotation velocity was taken to be $v_{\text{rot}} = 1.8$ km s$^{-1}$, and the most probable velocity of the radial and tangential motions is $V_{\text{mac}} = 3.5$ km s$^{-1}$.

In F–G–K stars, neutral aluminum is observed in the AI 3944, 3961 Å resonance doublet and subordinate lines in the red and infrared (IR) spectral ranges. Their list together with the atomic parameters and data sources is given in Table 1. The AI 3944 Å line is blended with a strong CH molecular line and is not used to determine the abundance. Note that with the van der Waals broadening constant $\log C_6 = -30.62$ calculated for AI 3961 Å based on the data from Anstee and O’Mara (1995), the non-LTE solar abundance from this line, $\log \epsilon_{\text{AI}} = 6.26$, is too low compared to the meteoritic $\log \epsilon_{\text{AI}} = 6.45$ (Lodders et al. 2009). It is even lower, by 0.06 dex, in LTE. Here, we use the scale where $\log \epsilon_{\text{H}} = 12$ for hydrogen. Taking $\log \epsilon_{\text{AI}} = 6.45$ as the solar abundance, we found the observed AI 3961 Å profile to be reproduced in our non-LTE calculations if $\log C_6 = -31.2$ (Fig. 4). It is this value that we will use in analyzing the solar and stellar spectra.

We found the mean abundance from six subordinate lines to be $\log \epsilon_{\text{AI}} = 6.36 \pm 0.07$ in the non-LTE calculations (BO), $6.39 \pm 0.08$ in the non-LTE calculations including the collisions only with electrons, case (e), and $6.34 \pm 0.07$ in LTE. The rms deviation $\sigma_x = \sqrt{\sum(x_i-x)^2/(N-1)}$ is given everywhere as the statistical error. The abundance derived under the assumption of LTE is lower than the meteoritic one by 0.11 dex. The non-LTE calculations lead to better agreement between the photospheric and meteoritic abundances, though it should be noted that the departures from LTE for AI I lines in the solar atmosphere are small.

4. THE ALUMINUM ABUNDANCE IN SELECTED STARS

From an observational point of view, the aluminum abundance is difficult to determine in F–G–K stars, especially if the sample includes stars in a wide metallicity range. The AI 3961.52 Å resonance line is located in the Ca II K 3968.47 Å line wing and is difficult to measure in stars with $[\text{Fe/H}] < -1$. This means that for a sample of stars, say, in the range $-3 < [\text{Fe/H}] < 0$, the aluminum abundance is determined from different lines at the opposite ends of the metallicity scale, and we should be sure that using different lines does not lead to a systematic shift in the abundance. In this paper, we choose three moderately metal-deficient stars for which high-resolution spectra are available in the entire spectral range where the AI I lines are located and...
Table 1. Atomic parameters of the Al I lines under study and the solar abundance derived in different line formation models

| λ, Å  | $E_{\text{exc}}$, eV | log $gf$ | Source | log $C_6$ | Source | Solar abundance, log $\varepsilon_{\text{Al}}$ |
|-------|-------------------|----------|--------|-----------|--------|-----------------------------------|
|       |                   |          |        |           |        | LTE         | (BO)$^1$ | (e)$^2$ |
| 3961.52 | 0.01             | -0.33    | NIST   | -31.2     | S      | 6.39 | 6.45 | 6.45 |
| 6696.02 | 3.14             | -1.35    | WSM    | -30.52    | K12    | 6.27 | 6.29 | 6.32 |
| 6698.67 | 3.14             | -1.65    | WSM    | -30.52    | K12    | 6.25 | 6.27 | 6.30 |
| 7836.31 | 4.02             | -0.61    | K12    | -29.77    | K12    | 6.38 | 6.40 | 6.44 |
| 7836.13 | 4.02             | -0.46    | K12    | -29.77    | K12    | 6.45 | 6.47 | 6.51 |
| 8772.86 | 4.02             | -0.21    | K12    | -29.01    | B98    | 6.31 | 6.32 | 6.37 |
| 8773.90 | 4.02             | -0.06    | K12    | -29.01    | B98    | 6.37 | 6.38 | 6.42 |

$^1$ Non-LTE, the collisions with H I are taken into account according to Belyaev (2013).

$^2$ Non-LTE, the collisions only with electrons are taken into account.

NIST = Ralchenko et al. (2008); WSM = Wiese et al. (1969); S = analysis of the profile in the solar spectrum, B98 = Barklem et al. (1998); K12 = http://kurucz.harvard.edu/atoms.html.

Table 2. Aluminum abundance [Al/H] in stars in different Al I line formation models

| Al I lines | Mean |
|-----------|------|
|           |      |
| 3961      |      |
| 6696      |      |
| 6698      |      |
| 7835      |      |
| 7836      |      |
| 8772      |      |
| 8773      |      |

Model

LTE

(BO)$^1$

(e)$^2$

(Dr$^*0.1)^3$

Model

LTE

(BO)

(e)

(Dr$^*0.1$)

Model

LTE

(BO)

(BO)bf)$^4$

(e)

(Dr$^*0.1$)

$^1$ Non-LTE, the Al + H collisions according to Belyaev (2013).

$^2$ Non-LTE, the collisions only with electrons.

$^3$ Non-LTE, the Al + H collisions with the Drawin rates, $S_{\text{H}} = 0.1$.

$^4$ Non-LTE, only Al$^0 + H^+ \rightarrow Al^+ + H^+$. 

1. Non-LTE, the Al + H collisions according to Belyaev (2013).
2. Non-LTE, the collisions only with electrons.
3. Non-LTE, the Al + H collisions with the Drawin rates, $S_{\text{H}} = 0.1$.
4. Non-LTE, only Al$^0 + H^+ \rightarrow Al^+ + H^+$. 

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investigate how the application of accurate collisional data affects the abundance derived from individual lines and the agreement between the results from different lines.

The stellar spectra were taken by T. Gehren at the Calar Alto Observatory (Spain) using the 2.2-m telescope with the FOCES spectrograph. The spectral range is 3800–8800 Å, the spectral resolution is $R = \lambda/\Delta \lambda \simeq 40,000$, and the signal-to-noise ratio is $S/N \simeq 200$. The IR spectra were reduced by H. Zhang.

The atmospheric parameters were taken from Sitnova et al. (2015). The effective temperatures and surface gravities were determined using several methods, namely by the IR flux method ($T_{\text{eff}}$), by invoking the trigonometric parallaxes ($\log g$), and spectroscopically through a non-LTE analysis of Fe I and Fe II lines ($T_{\text{eff}}$, $\log g$). The iron abundance and microturbulence were determined simultaneously with $T_{\text{eff}}$ and $\log g$ in our non-LTE analysis of Fe I and Fe II lines. The data are given in Table 2.

The aluminum abundance in the stars was determined under the assumption of LTE and in our non-LTE calculations by taking into account the collisions only with electrons (e) and including the collisions with H I atoms with accurate data from Belyaev (2013, BO). For each line, we applied a differential approach, i.e., the abundance was determined relative to the solar one we derived: $[\text{Al/H}] = \log (g f_{\text{Al}}) - \log (g f_{\text{Si}})$. The results are presented in Table 2. Just as for the Sun, the stellar spectra were analyzed by the synthetic-spectrum method. The theoretical profile was convolved with the instrumental profile and the radial–tangential microturbulence profile. For the instrumental profile, we used a Gaussian with a dispersion of 4 km s$^{-1}$, $V_{\text{mac}}$ was varied within 10% near 4.5 km s$^{-1}$ for the subordinate lines and was larger, 6 and 8 km s$^{-1}$ for the resonance line. Since all stars rotate slowly ($v_{\text{rot}} \sin i \leq 1$ km s$^{-1}$), the rotational broadening was disregarded. An example of the observed spectra and their reproduction in models are shown in Fig. 5.

Let us analyze how the abundances from different lines agree in different Al I line formation models. The LTE calculations yield a large scatter in abundances between different lines with $\sigma_\varepsilon$ from 0.12 to 0.21 dex. This is also illustrated by Fig. 6, where the difference in abundance between each line and the mean value from the subordinate lines is presented. The subordinate lines give a systematically higher LTE abundance than does Al I 3961 Å, by 0.25–0.45 dex for different stars. This means that significant systematic shifts will arise when the aluminum abundance is determined for a sample of stars under the assumption of LTE if we use the subordinate lines for stars with [Fe/H] $> -1$ but Al I 3961 Å for highly metal-deficient stars.

Non-LTE leads to a dramatic weakening of the resonance line in all calculations and a positive non-LTE correction to the abundance with $\Delta_{\text{NLTE}}$ from 0.36 to 0.77 dex. Since the non-LTE corrections for the subordinate lines are almost equally large in non-LTE(e), the scatter in the abundances from different lines turns out to be even greater than that in LTE, and $\sigma_\varepsilon$ reaches 0.18–0.23 dex. Allowance for the collisions with H I leads to a decrease in the departures from LTE and the non-LTE corrections for the subordinate lines. For example, for Al I 8773 Å in the star HD 134169, $\Delta_{\text{NLTE}} = 0.63$ dex in non-LTE(e) and 0.09 dex in non-LTE(BO). For Al I 6696 Å in the same star, the corresponding values are 0.33 and
0.12 dex. In non-LTE(BO), the mean abundance for different stars increases compared to LTE by 0.14–0.19 dex, the scatter in the abundances from different lines decreases, and the error becomes smaller by a factor of 1.5–3. These calculations prove conclusively that inelastic collisions with H I affect the SE of aluminum and must be taken into account when calculating a theoretical spectrum. The application of accurate collisional data allows the astrophysical results to be improved.

We performed test calculations by excluding the excitation processes due to the collisions with H I. The results for HD 134169 are given in Table 2 (BObf). The departures from LTE for the subordinate lines increased, but insignificantly, so that ΔNLTE increased by 0.02–0.03 dex. In contrast, for the resonance line, ΔNLTE decreased by 0.08 dex, and the abundance turned out to be the same as in the case of allowance for the collisions only with electrons. These results suggest that all processes of the interaction between Al I and H I atoms need to be taken into account.

Since the collisions with H I for most atoms are still taken into account using the formulas from Steenbock and Holweger (1984) due to the absence of accurate data, we also determined the aluminum abundance in our non-LTE calculations with scaled Sf, 0.1 Drawin rates, (Dr°0.1) in Table 2. For the resonance line, the non-LTE effects turned out to be very close to the case where the collisions only with electrons are taken into account. For the subordinate lines, the departures from LTE decreased compared to non-LTE(e) but increased compared to non-LTE(BO). The mean abundance in non-LTE(e) turns out to be larger than that in non-LTE(BO) by 0.27–0.31 dex, while the abundance difference (Dr°0.1–BO) is from 0.07 to 0.11 dex for different stars. Thus, it is better to use the approximate formulas than to disregard the collisions with H I altogether.

5. THE NON-LTE CORRECTIONS FOR AL I LINES AS A FUNCTION OF STELLAR PARAMETERS

The non-LTE method based on accurate rates of inelastic collisions with H I was applied to calculate the non-LTE corrections to the abundance for six subordinate Al I lines in MARCS model atmospheres with Teff = 4500–6500 K, log g = 3.0–4.5, and [M/H] = 0 and −1. Even the strongest of them, Al I 8772 and 8773 Å become unmeasurable at [M/H] < −1. The corrections were computed with the LINEC code (Sakhibullin 1983), which uses the non-LTE and LTE level populations computed in the DETAIL code. By the non-LTE correction we will mean the change in abundance required to reproduce the equivalent width (EW) calculated in LTE. Everywhere in our calculations, the aluminum abundance follows the metal abundance in the model, i.e., [Al/M] = 0, and \( \xi_t = 2 \, \text{km s}^{-1} \) in accordance with the value adopted in the computations of MARCS model atmospheres. The results are presented in Table 3 and, for the three lines of different multiplets, in Fig. 7.

In the coolest models with \( T_{\text{eff}} = 4500 \) and 5000 K, the overionization of Al I is indistinct in the subordinate line formation region, while the radiative pumping in strong resonance lines leads to an overpopulation of the 4s 2S level and the 4p 2P° and 3d 2D levels, which are the lower levels of the transitions being studied. As a result, the subordinate Al I lines are strengthened compared to LTE, and the non-LTE corrections to the abundance are negative, though small in absolute value (<0.1 dex). The overionization of Al I develops and becomes stronger with
Table 3. Non-LTE corrections (dex) for the subordinate Al I lines as a function of the effective temperature and surface gravity in the models with [M/H] = 0 and −1. Non-LTE corrections are given for lines with \( EW(\text{LTE}) < 5 \text{ mÅ} \)

| \( T_{\text{eff}} \) | \( \log g \) | [M/H] = 0 | All I lines | [M/H] = −1 | All I lines |
|---|---|---|---|---|---|
| | | 6696 | 6698 | 7835 | 7836 | 8772 | 8773 | 6696 | 6698 | 7835 | 7836 | 8772 | 8773 |
| 4500 | 3.0 | −0.06 | −0.04 | −0.02 | −0.03 | −0.06 | −0.08 | 4500 | 3.0 | −0.08 | −0.07 | −0.03 | −0.03 | −0.05 | −0.06 |
| 4500 | 3.5 | −0.05 | −0.04 | −0.02 | −0.02 | −0.05 | −0.06 | 4500 | 3.5 | −0.06 | −0.06 | −0.02 | −0.02 | −0.05 | −0.05 |
| 4500 | 4.0 | −0.04 | −0.03 | −0.01 | −0.01 | −0.04 | −0.05 | 4500 | 4.0 | −0.05 | −0.05 | −0.02 | −0.02 | −0.04 | −0.04 |
| 4500 | 4.5 | −0.03 | −0.02 | −0.01 | −0.01 | −0.02 | −0.03 | 4500 | 4.5 | −0.04 | −0.03 | −0.01 | −0.01 | −0.03 | −0.03 |
| 5000 | 3.0 | −0.05 | −0.04 | −0.02 | −0.02 | −0.05 | −0.06 | 5000 | 3.0 | −0.02 | −0.02 | 0.02 | 0.01 | −0.01 | −0.01 |
| 5000 | 3.5 | −0.04 | −0.03 | −0.01 | −0.02 | −0.04 | −0.05 | 5000 | 3.5 | −0.02 | −0.02 | 0.01 | 0.01 | −0.01 | −0.01 |
| 5000 | 4.0 | −0.03 | −0.02 | −0.01 | −0.01 | −0.03 | −0.03 | 5000 | 4.0 | −0.02 | −0.01 | 0.00 | 0.00 | −0.01 | −0.01 |
| 5000 | 4.5 | −0.02 | −0.01 | −0.01 | −0.01 | −0.02 | −0.02 | 5000 | 4.5 | −0.01 | −0.01 | 0.00 | 0.00 | −0.01 | −0.01 |
| 5500 | 3.0 | −0.03 | −0.02 | −0.00 | −0.01 | −0.03 | −0.04 | 5500 | 3.0 | 0.15 | 0.17 | 0.14 | 0.14 |
| 5500 | 3.5 | −0.02 | −0.01 | −0.00 | −0.00 | −0.02 | −0.03 | 5500 | 3.5 | 0.11 | 0.11 | 0.09 | 0.09 |
| 5500 | 4.0 | −0.01 | −0.01 | 0.00 | 0.00 | −0.01 | −0.02 | 5500 | 4.0 | 0.08 | 0.07 | 0.05 | 0.06 |
| 5500 | 4.5 | −0.01 | 0.00 | 0.00 | 0.00 | −0.01 | −0.01 | 5500 | 4.5 | 0.05 | 0.04 | 0.03 | 0.03 |
| 6000 | 3.0 | 0.06 | 0.06 | 0.08 | 0.09 | 0.07 | 0.07 | 6000 | 3.0 | 0.34 | 0.35 |
| 6000 | 3.5 | 0.06 | 0.06 | 0.07 | 0.07 | 0.07 | 0.07 | 6000 | 3.5 | 0.24 | 0.25 |
| 6000 | 4.0 | 0.05 | 0.05 | 0.06 | 0.06 | 0.06 | 0.06 | 6000 | 4.0 | 0.16 | 0.16 |
| 6000 | 4.5 | 0.04 | 0.04 | 0.04 | 0.04 | 0.04 | 0.05 | 6000 | 4.5 | 0.10 | 0.10 |
| 6500 | 3.0 | 0.16 | 0.15 | 0.20 | 0.20 | 0.19 | 0.21 | 6500 | 3.0 | 0.51 | 0.54 |
| 6500 | 3.5 | 0.15 | 0.14 | 0.17 | 0.17 | 0.18 | 0.19 | 6500 | 3.5 | 0.41 | 0.43 |
| 6500 | 4.0 | 0.13 | 0.12 | 0.14 | 0.14 | 0.14 | 0.16 | 6500 | 4.0 | 0.30 | 0.31 |
| 6500 | 4.5 | 0.10 | 0.10 | 0.10 | 0.10 | 0.11 | 0.12 | 6500 | 4.5 | 0.20 | 0.21 |
INFLUENCE OF INELASTIC COLLISIONS

rising \( T_{\text{eff}} \), which leads to a weakening of the Al I lines compared to LTE and to a positive \( \Delta_{\text{NLTE}} \). In metal-deficient models, the overionization of Al I is stronger than that in models with \([M/H] = 0\) due to the higher intensity of ionizing ultraviolet radiation and, therefore, \( \Delta_{\text{NLTE}} \) is greater.

For the subordinate lines, the non-LTE corrections were calculated as the change in abundance required to reproduce \( EW(\text{LTE}) \). However, such an approach is inapplicable for the Al I 3961 Å resonance line, which is located in the wing of the very strong Ca II 3968 Å line (see, for example, Fig. 5 for the star HD 134169 with \([\text{Fe/H}] = -0.78\)). How should \( EW \) of the Al I 3961 Å line be specified in this case? How should the calcium abundance be specified? We performed test calculations for the 5890/4.02/−0.78 model atmosphere. Our description of the observed Al I 3961 Å line profile in the spectrum of HD 134169 gives an abundance difference of 0.77 dex between non-LTE and LTE (Table 2). In this case, we found that \([\text{Ca/Fe}] = 0.3\) led to a satisfactory description of the Ca II 3968 Å wing, and such a calcium abundance is quite expectable for a star with \([\text{Fe/H}] = -0.78\). We then performed calculations with the LINEC code, in which the background opacity includes the Ca II 3933, 3968 Å lines, along with continuous absorption sources. The non-LTE correction for Al I 3961 Å turned out to depend on the Ca abundance in the model:

1. \( \Delta_{\text{NLTE}} = 0.40 \text{ dex} \) if the Ca abundance is zero, i.e., the absorption in Ca II 3968 Å is disregarded;

2. \( \Delta_{\text{NLTE}} = 0.46 \text{ dex} \) if \([\text{Ca/Fe}] = 0\);

3. \( \Delta_{\text{NLTE}} = 0.60 \text{ dex} \) if \([\text{Ca/Fe}] = 0.4\).

Thus, the synthetic-spectrum method should be used to determine the abundance from the Al I 3961 Å line in a specific star irrespective of whether this is done in LTE or non-LTE. Using the non-LTE corrections calculated from the equivalent widths and with a Ca abundance that can be far from that observed in a specific star can yield an incorrect result. In order not to mislead the potential users, we do not provide any table of non-LTE corrections for Al I 3961 Å and do not recommend to use the non-LTE corrections for this line irrespective of the source of their origin.

To illustrate the dependence of the departures from LTE in the Al I 3961 Å line on stellar atmosphere parameters, in Fig. 8 we present the corrections calculated by the method of equivalent widths and with a Ca abundance corresponding to the model atmosphere, more specifically, \([\text{Ca/Fe}] = 0.4\) for all models with \([M/H] \leq -1\). Obviously, for highly metal-deficient stars, where the aluminum abundance can be determined only from the resonance line, the theoretical spectra should be computed based on non-LTE. Using LTE can lead to an underestimation of

Fig. 6. Difference in abundance between each Al I line and the mean value obtained from the subordinate lines in different line formation scenarios: (a) LTE, (b) non-LTE including the collisions only with electrons, and (c) non-LTE including the collisions with H I atoms. Different symbols correspond to the stars HD 59374 (triangles), HD 59984 (diamonds), and HD 134169 (circles).
the abundance by 0.4–0.8 dex, depending on the stellar parameters.

6. CONCLUSIONS

We performed non-LTE calculations for Al I and Si I lines using accurate rates of collisions with H I from Belyaev (2013) and Belyaev et al. (2014) for models corresponding to the atmospheres of F–G–K stars. We showed that for b–b transitions with $E_{ij} \leq 2$ eV, the collisions with H I are equally important as the collisions with electrons. For Si I there also exist many transitions with $E_{ij} > 3$ eV, where $C_{ij} \simeq C_e$. The rates of ion-pair production $\text{Al I}(nl) + \text{H} \rightarrow \text{Al II}(3s \, ^2S) + \text{H}^-$ and $\text{Si I}(nl) + \text{H} \rightarrow \text{Si II}(3p \, ^2P^o) + \text{H}^-$ are higher than the electron-impact ionization rates by two or three orders of magnitude for levels with an ionization energy $\chi_i < 3.5$ eV. Owing to their large cross sections, the charge exchange processes provide a close coupling of highly excited levels of neutral atoms (Al I and Si I) with the ground state of ions (Al II and Si II) and a reduction of the departures from LTE in their populations compared to the calculations that take into account the collisions only with electrons. For Si I including the collisions with H I leads to the establishment of equilibrium populations in the spectral line formation region even in hot metal-deficient models. As a consequence, the departures from LTE in Si I lines are vanishingly small. However,
including the collisions with H I barely affects the population of the Al I ground level, because the radiative rates in all of the transitions that couple it with the excited levels and with Al II exceed the collisional rates.

Our calculations confirm the existence of the effects detected previously by Baumueller and Gehren (1996), Andrievsky et al. (2008), and Menzhevitski et al. (2012) for Al I, namely Al I overionization in the spectral line formation region and a weakening of lines compared to the case of LTE. However, our results, especially for the subordinate lines, quantitatively differ greatly from the results of our predecessors.

For three moderately metal-deficient dwarf stars, for which there are reliable determinations of their atmospheric parameters and high-quality observations in a wide spectral range, the aluminum abundance was determined from seven Al I lines in different models of their formation. Under the assumption of LTE and in non-LTE calculations including the collisions only with electrons, the Al I 3961 Å resonance line gives a systematically lower abundance than do the subordinate lines, by 0.25–0.45 dex. The difference for each star is removed by taking into account the collisions with hydrogen atoms, and the rms error of the abundance derived from all seven Al I lines decreases by a factor of 1.5–3 compared to the LTE analysis. Thus, the collisions with H I should be taken into account in non-LTE calculations for Al I, and using the rates from Belyaev (2013) plays a key role in increasing the accuracy of the derived aluminum abundance.

The developed method was applied to calculate the non-LTE corrections to the abundance for six subordinate lines in a grid of model atmospheres with $T_{\text{eff}} = 4500–6500$ K, $\log g = 3.0–4.5$, and $[M/H] = 0$, $-1$, $-2$, and $-3$. The departures from LTE grow with rising effective temperature, decreasing metal abundance, and decreasing surface gravity. For the Al I 3961 Å resonance line, we do not recommend to use the non-LTE corrections. The abundance from this line should be determined by the synthetic-spectrum method irrespective of which approach, LTE or non-LTE, is applied.

We hope that progress in the theory will allow accurate data to be obtained in future for inelastic collisions with H I atoms for O I, Fe I, Ti I, and other atoms that play an important role in studying the physics of stellar atmospheres.

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