NLTE analysis of Sr lines in spectra of late-type stars with new R-matrix atomic data

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ABSTRACT

We investigate statistical equilibrium of neutral and singly-ionized strontium in late-type stellar atmospheres. Particular attention is given to the completeness of the model atom, which includes new energy levels, transition probabilities, photoionization and electron-impact excitation cross-sections computed with the R-matrix method. The NLTE model is applied to the analysis of Sr I and Sr II lines in the spectra of the Sun, Procyon, Arcturus, and HD 122563, showing a significant improvement in the ionization balance compared to LTE line formation calculations, which predict abundance discrepancies of up to 0.5 dex. The solar Sr abundance is $\log \epsilon = 2.93 \pm 0.04$ dex, in agreement with the meteorites. A grid of NLTE abundance corrections for Sr I and Sr II lines covering a large range of stellar parameters is presented.

Key words. atomic data – line: formation – radiative transfer – sun: abundances – stars: abundances

1. Introduction

Spectroscopic observations of low-mass stars have shaped our understanding of Galactic evolution and stellar nucleosynthesis. Strontium, as one of the abundant r-process elements, has been extensively investigated in the past few decades. However, its main production site has not yet been identified: the observed abundances of Sr in metal-poor stars are far too large to be explained by conventional rapid neutron-capture nucleosynthesis in SNe II, suggesting some alternative exotic scenarios, such as the light element primary process (Travaglio et al. 2004), rp-process in accretion disks around low-mass black holes (Schatz et al. 1998), black hole - neutron star mergers (Surman et al. 2008), high-entropy winds in SN II (Farouqi et al. 2010), and low-mass electron-capture supernovae (Wanajo et al. 2011) to name just a few (see Josè & Iliadis 2011, and references therein).

Until recently, determinations of Sr abundances in metal-poor stars relied almost exclusively on the two near-UV lines of Sr\textsc{ii}, which are sufficiently strong to be detected also in the spectra of moderate-to-low resolution and signal-to-noise. The drawback is that in spectra of stars typically used for studies of Galactic chemical evolution, [Fe/H] $\approx -2$, these lines saturate and develop pronounced damping wings, overlapping with various atomic and molecular blends. Thus, at higher metallicities a preference is sometimes given to the weak Sr\textsc{i} line at 4607.34 Å and/or the Sr\textsc{ii} line at 4161 Å. However, there is evidence that the Sr\textsc{i} lines may be subject to non-local thermodynamic equilibrium (hereafter, NLTE) effects (Barklem & O’Mara 2000), which has been supported by \textit{ab initio} calculations solving for radiative transfer in NLTE for a small grid or red giant model atmospheres (Short & Hauschildt 2000). In a few studies utilizing near-IR spectra, also the Sr\textsc{ii} triplet (10037, 10327, and 10915 Å) have been used (Andrievsky et al. 2011). In the majority of published studies, the preference is given to one ionization stage only (e.g., Jehin et al. 1999), and only a few studies investigated both Sr\textsc{i} and Sr\textsc{ii} lines (Gratton & Sneden 1994, Cowan et al. 2002), finding discrepant results.

In this study, we perform for the first time a NLTE analysis of the Sr\textsc{i} and Sr\textsc{ii} lines in spectra of late-type stars. The new atomic model was constructed from the state-of-art atomic data, computed specifically for this work. The NLTE model atom is tested on a number of reference stars with parameters determined by other independent methods. Furthermore, we present a large grid of NLTE abundance corrections for Sr\textsc{i} and Sr\textsc{ii} lines. The results presented in this work will be applied to the analysis of a representative sample of metal-poor stars observed at a very-high resolution and signal-to-noise in Hansen et al. (in prep.) Furthermore, we plan to undertake the NLTE Sr abundance analysis of the thick-disk and halo stars with spectroscopic parameters from Ruchti et al. (2011).

2. Methods

2.1. Model atmospheres and NLTE calculations

All calculations in this work were performed with classical 1D LTE plane-parallel model atmospheres. We used the MAFAGS-OS (Grupp 2004a,b) and MAFAGS-ODF models, which are well-adapted for the analysis of late-type stars. A brief description of these models and comparison with other models of a similar type (MARCS, Gustafsson et al. 2008) is presented in Bergemann et al. (2012).

The NLTE statistical equilibrium calculations were performed with the revised version of the DETAIL code (Butler & Giddings 1985). The statistical equilibrium and radiative transfer equations are solved by the accelerated A-iteration method in the formulation of Rybicki & Hummel (1991, 1992). The method allows for self-consistent treatment of overlapping lines and continua. A description of the code with some recent modifications related to the treatment of background line opacity

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can be found in Bergemann et al. (2012). In the statistical equilibrium calculations, we treat the diagnostic lines investigated here with Voigt profiles, whereas all other Sr lines were computed with a Gaussian profile with 13 frequency points.

The LTE and NLTE abundances for the reference stars, as well as the NLTE abundance corrections, were determined by full spectrum synthesis with the revised version of the SIU code (Reetz 1999), which was adapted for automated NLTE abundance calculations (L. Sbordone, private communication). The line lists in SIU have been continuously updated by the members of the Munich group (Gehren et al. 2004, 2006) and they are mainly based on the Kurucz Hannover and NIST (Ralphenko et al. 2012) compilations.

2.2. Model atom

The model atom of Sr was constructed using the available atomic data from the Hannover and NIST databases and supplemented by our new calculations of atomic energy levels, dipole-permitted and forbidden transitions (see below). Thus, the model is the most complete representation of atomic system of Sr i and Sr ii, to date. In particular, we also include 135 Sr i transitions from the Coulomb approximation calculations of Lindgård & Nielsen (1977). The neutral atom is, thus, represented by 141 levels with the principle quantum numbers \( n \leq 20 \), and the uppermost level 20f \(^1\)\(^F\) is located at 5.66 eV, i.e., 0.03 eV from the 1-st ionization limit at 5.69 eV. The singly-ionized atom includes 49 levels including the highest experimentally-observed level 14d \(^2\)D, which is located at 10.684 eV. The model is closed by the Sr ii ground state 4p \(^6\)1s\(^0\). The number of dipole-allowed transitions is 336 and 214 for Sr i and Sr ii, respectively.

Our model of Sr is similar to that of Andrievsky et al. (2011) with respect to the term structure and the number of dipole-permitted transitions of Sr ii. However, we include more accurate transition probabilities for Sr ii. More importantly, our model atom includes new quantum-mechanical photoionization cross-sections for Sr i and \( e^- \) impact excitation cross-sections for dipole-permitted and forbidden transitions in Sr ii (see below). The model atoms employed by Short & Hauschildt (2006) and Mashonkina & Gehren (2001) are much simpler than that employed in this work (84 and 41 total levels, respectively).

2.2.1. New calculations of atomic data

We compute single-electron orbitals for the target ions \(^{+}\)Sr with the atomic structure code autostructure by Badnell (1986, 1997). This code is an extension of the program superstructure (Eissner & Nussbaumer 1969), computes fine-structure CI energy levels, and radiative and Auger rates in a Breit-Pauli relativistic framework. Single electron orbitals are constructed

1 kurucz.harvard.edu/
2 www.pmp.uni-hannover.de/cgi-bin/ssii/test/kurucz/sekur.html. The atomic line data used in this database are taken from Kurucz & Bell (1995).
3 The new atomic data for Sr presented in the paper and in Bautista et al. (2002), including the energy levels, \( f \)-values, photoionization and \( e^- \) impact excitation cross-sections, can be provided upon request.
4 We note that different notations are standard for atomic physics and spectroscopy, and we chose to respect this difference in our work. Ionization of an element is thus indicated by a numerical superscript in Sect. 2.2.1 describing the atomic physics calculations and by a roman numerical in astrophysical application in the further sections.
Fig. 2. Partial photoionization cross-sections for the Sr\textsc{i} ground state into the lowest six states of the Sr\textsc{ii} target. The cross-sections are dominated by prominent resonances.

get threshold), and another mesh of 0.02 Ry from 2.3 Ry to 4.3 Ry. Figure 2 shows the partial cross sections for ionization from the ground term of Sr\textsc{i} into each of the first six terms of the target Sr\textsc{ii}. We note, however, that DETAIL is not yet able to consistently treat ionization to specific states of the target ion. Therefore, total photoionization cross-sections are adopted here. Figure 3 shows the total cross sections for the lowest six excited state of Sr\textsc{i}. For comparison, we show the hydrogenic cross-sections computed with effective principal quantum number. The latter reproduce the functional dependence of the cross-section with energy. However, the quantum-mechanical photoionization cross-sections are typically larger both in the background (factor of two to five) and in the resonances (up to three orders of magnitude).

2.2.2. Collision rates

Electron impact excitation rate coefficients for the 49 levels of Sr\textsc{n} levels were taken from Bautista et al. (2002). The data were computed with the same technique, i.e., the close-coupling approximation with the R-matrix method. For all other levels of Sr\textsc{i} and Sr\textsc{n} the electron collision rates were computed using the formulae of van Regemorter (1962) for dipole-allowed and Allen (1973) for dipole-forbidden transitions. Cross-sections for transitions caused by inelastic collisions with H\textsc{i} atoms are basically unknown for any atom heavier than Mg. The only available formula developed originally for collisions between equal H-like atoms (Drawin 1968, 1969) and later slightly modified by Steenbock & Holweger (1984) and Lambert (1993) was shown to over-estimate the rates of bound-bound transitions by two to seven orders of magnitude (Belyaev & Barklem 2003, Barklem et al. 2012). Furthermore, the charge transfer processes can not be described by this simple classical formula at all. Our tests with various scaling factors to the Drawin's H\textsc{i} inelastic collision cross-sections demonstrated that excitation and ionization balance of Sr\textsc{i}/Sr\textsc{ii} in the reference stars can be satisfied simultaneously only if the efficiency of H\textsc{i} collisions is very low (Sect. 4), few orders of magnitude lower than that given by the Drawin's recipe, which is consistent with the quantum-mechanical results mentioned above. On these grounds we do not include inelastic H\textsc{i} collisions in our reference model atom.

2.3. Line selection and atomic data

In spectra of FGK stars, only a few Sr lines are useful for abundance determinations. These are two Sr\textsc{i} and 6 Sr\textsc{n} lines. The Sr\textsc{i} resonance line at 4607 Å is weak, and, thus, sufficiently reliable for more metal-rich stars ([Fe/H] > −1.5) observed at high resolution. The two resonance Sr\textsc{n} lines, in contrast, remain strong even in spectra of very metal-poor stars, however, severe blends in the inner and outer wings lead to systematic errors in abundances, when the blends are not properly accounted for. The
Table 1. Lines of Sr\textsc{i} and Sr\textsc{ii} selected for abundance calculations.

| Isotope | $E_{low}$ | $E_{up}$ | $\log g f$ | $\log (\gamma/N_1)$ | $\log C_6$ |
|---------|---------|---------|----------|-----------------|----------|
| Sr\textsc{i} |         |         |          |                 |          |
| 4607.33 | 0.00 5s$^1$S$_1$ | 2.69 5p$^1$P$_1$ | 0.283 | $-7.53$ | $-31.2$ |
| 7070.01 | 1.85 5p$^2$P$_2$ | 3.60 6s$^1$S$_1$ | $-0.020$ | $-7.15$ | $-30.2$ |
| 4077.71 | 0.00 5s$^1$S$_1$/2 | 3.04 5p$^3$P$_0$ | 0.158 | $-7.81$ | $-32.0$ |
| 4167.80 | 2.94 5p$^3$P$_1$ | 5.92 6s$^1$S$_1$ | $-0.502$ | $-7.81$ | $-32.0$ |
| 4215.52 | 0.00 5s$^1$S$_1$/2 | 2.94 5p$^3$P$_0$ | $-0.155$ | $-7.81$ | $-32.0$ |
| 10036.65 | 1.81 4d$^3$D$_3$/2 | 3.04 5p$^3$P$_0$ | $-1.194$ | $-7.64$ | $-31.5$ |
| 10327.31 | 1.84 4d$^3$D$_3$/2 | 3.04 5p$^3$P$_0$ | $-0.240$ | $-7.64$ | $-31.5$ |
| 10914.89 | 1.80 4d$^3$D$_3$/2 | 2.94 5p$^3$P$_0$ | $-0.474$ | $-7.64$ | $-31.5$ |

Notes.
1. (blend). 2. Log of the FWHM per H atom at 10000 K. 3. $C_6$ are in the units of cm$^8$ s$^{-1}$.

Table 2. Hyperfine structure and isotopic shift for the 4077 Å line.

| Isotope | $\lambda$ | $log g f$ |
|---------|---------|----------|
| 87     | 4077.697 | $-1.645$ |
| 87     | 4077.699 | $-1.485$ |
| 87     | 4077.708 | $-2.094$ |
| 87     | 4077.709 | $-0.848$ |
| 84     | 4077.710 | 0.075    |
| 87     | 4077.724 | $-1.465$ |
| 87     | 4077.725 | $-1.956$ |

three near-IR Sr\textsc{ii} lines at 1$\mu$m appear to be un-blended and are sufficiently strong to be detected even at [Fe/H] $<-2$.

The lines selected in this work are given in Table 1. The transition probabilities were taken from different experimental and theoretical sources. According to NIST, the Sr\textsc{i} $g f$-values are very accurate, the uncertainty is lower than 1 percent for the 4607 Å line and better than 10 percent for the 7070 Å subordinate line. These $g f$-values are taken from the laboratory analysis of Parkinson (1976) and García & Campos (1988), respectively. The oscillator strengths of the near-IR Sr\textsc{ii} lines were adopted from the only available laboratory results of Gallagher (1967). These are fully consistent with the most recent ab initio calculations of Bautista et al. (2002), which are also adopted in the NLTE model atom.

Damping widths for the calculation of broadening due to elastic collisions with H\textsc{i} are available for the Sr\textsc{ii} lines from the quantum-mechanical calculations of Barklem et al. (2000) and for the Sr\textsc{i} lines they were kindly provided by the referee. We adopt the values from Barklem et al. for the near-IR Sr\textsc{ii} transitions. However, for the strong resonance Sr\textsc{ii} lines, which are very sensitive to this parameter, our spectrum synthesis calculations indicate somewhat lower values, by $\sim 20$ percent. The question is whether this difference can be attributed to the uncertainties of the theoretical data arising because of certain approximations in the Anstee, Barklem, and O’Mara (ABO) theory. These include representation of the interaction potential and collisional dynamics and have been investigated by Kerkani et al. (2004) for neutral atoms of Mg, Sr, Ca, and Na. They find that whereas the semi-classical description of a collision is a sufficiently-good approximation, the ABO potentials become rather inaccurate at small interatomic separations. However, the effect of the latter is rather to under-estimate the line width that clearly does not explain our finding. For the Sr\textsc{ii} 4167 line we assume the same y as for the resonance lines, but note that the line is too weak even in the solar spectrum to be sensitive to this parameter at all.

The adopted values of damping widths at 10000 K per H atom, $\gamma/N_1$, as well as the commonly-used interaction constants in the van der Waals-type interaction $C_6$ computed for $T = 5780$ K, are given in the Table 1. Here, the latter is in parameter in the Unsöld (1955, Eq. 82.48) formula needed to reproduce the correct line width $y$ at the given temperature.

For the 4077 Sr\textsc{i} line, which will be used for the abundance determinations in Sect. 4, we included the hyperfine structure and isotopic shifts. The magnetic dipole and electric quadrupole constants were taken from Buchinger et al. (1990): A (5s$^2$1S$_0$) = $-1000.5$ MHz, $A$ (5p$^3$P$_{3/2}$) = $-36$ MHz, $B$ (5p$^3$P$_{3/2}$) = $85.5$ MHz. The data are in agreement with the older experiment by Borghs et al. (1983) and theoretical calculations by Yu et al. (2004). Isotope shifts for $^{85}$Sr, $^{86}$Sr, $^{87}$Sr were taken from Borghs et al. (1983) and the solar isotopic ratios were adopted.

The wavelengths and $log g f$ values of the hyperfine structure (HFS) components are given in Table 2. We note that the 4215 Å Sr\textsc{i} line is only used for a comparative LTE to NLTE analysis for the reasons discussed in Sect. 3.3.

3. Statistical equilibrium of Sr

The departure coefficient for selected levels of Sr are shown as a function of continuum optical depth $\log \tau_c$ at 500 in Fig. 4 for the model atmospheres with parameters corresponding to dwarfs and giants with [Fe/H] = 0 and $-2.4$. Only the levels, which give rise to the spectral lines selected for the abundance analysis (Table 1), are included in the plots. These are two levels of Sr\textsc{i}, forming the resonance lines at 4607 Å, and two levels of Sr\textsc{ii}, which are connected by the 4077 Å transition commonly used in the analysis of metal-poor stars.

3.1 NLTE effects in Sr\textsc{i}

Sr\textsc{i} with the ionization potential 5.69 eV is a trace atom in the atmospheres of late-type stars. The ratio of Sr\textsc{i}/Sr\textsc{ii} falls below 10$^{-3}$ above continuum optical depth unity, and the departures from LTE in the distribution of atomic level populations are almost entirely due to overionization. The behavior of the Sr\textsc{i} departure coefficients with stellar parameters, thus, resembles that of the similar trace atoms, such as Co (Bergemann et al. 2010) or Cr (Bergemann & Cescutti 2010). The levels are underpopulated and their $b_i$-factors monotonously decrease outwards. The ground state 5s$^2$1S with the ionization edge at 2177 Å is also strongly over-ionized and it decouples from the other low-exited levels because collisions can not bridge the large energy

5. Note that $C_6$ is often computed using the Unsöld (1955, Eq. 82.55) approximation, in which the interaction constant derives from the effective principal quantum number of lower and upper levels of a transition. However, this approximation is known to yield too small damping constants (e.g. Barklem 2007 and references therein)

6. http://www.nist.gov/physRefData/ASD/comp.cfm

7. The departure coefficients are defined as the ratio of NLTE to LTE level number densities, $b_i = n_{LTE}/n_{NLTE}$.
k, log $n_F$ance line becomes optically thin, and spontaneous transitions diminishes of the levels by radiative bound-bound transitions and collisions depopulate the upper level of the transition. 

Excited levels are large enough to keep them are in detailed balance even in the models with $[\text{Fe}/H]$. For example, in the metal-rich models (Fig. 4, left panel), the departure coëfficients smooth out, slowly decreasing outwards. The NLTE effects on the formation of the Sr I line at 4607 Å are generally to decrease the line opacity, shifting the line $	au_c$ scale to the deeper hotter atmospheric layers and, thus, leading to weaker lines compared to the LTE case.

3.2. NLTE effects in Sr II

The NLTE effects on the Sr II levels are due to the non-equilibrium excitation in the line transitions. This was extensively investigated by Belyakova & Mashonkina (1997) and Short & Hauschildt (2006), and our results are qualitatively very similar to these studies. The analysis of radiative and collisional rates populating and de-populating the levels, as well as trial calculations with atomic models devoid of some key transitions, in particular the resonance and subordinate lines, allows us to draw the following conclusions.

In the solar-metallicity models (Fig. 4 left panel), the ground state, $5s^2 S$ ($\lambda_{\text{thr}} \sim 1130 \AA$) and the lowest odd level $5p^2 P^o$ level...
(Δnlte ∼ 1540 Å, E ∼ 3 eV) have nearly LTE populations out to log τc ∼ −3, and become underpopulated only in the outermost layers, which are transparent to the radiation across their ionization edges. Still, there is small net radiative excitation in the transitions from these levels upwards, which very efficiently populates the levels with excitation energy E ≥ 6 eV, such as 5d^2D, 6p^2P^0, and 7s^2S (not shown in the plots) due to their small Boltzmann factors. The high-excited levels are, however, also over-ionicized, which in turn, increases the number density of the Sr in ground state. The lowest metastable Sr n level 4d^2D (E ∼ 1.8 eV) is nearly in thermal equilibrium with 5p^2P^0 due to the dominance of collision rates in the deeper layers, however in the outer layers it gains appreciable overpopulation due to the photon losses in the near-IR transitions between the two levels. In fact, this process is a part of photon suction in a sequence of low-frequency transitions, which connect 4d^2D to a large number of high-levels close to the continuum. Many of these transitions become optically thin at log τc < −3, so that spontaneous de-excitation leads to the overpopulation 4d^2D, the lowest energy state of the cascading sequence.

The departures from LTE change with decreasing metallicity (Fig. B right panel). First, the line transitions weaken and over-population of 4d^2D occurs in the deeper layers compared to the solar metallicity models. Also, radiative pumping in the resonance transitions at 4077 and 4215 Å is more efficient due to larger radiative fluxes. This process leads to a marked over-population of the 5p^2P^0 level at −1 < log τc < 0, especially in the hotter model, T_eff = 6200 K. Outside these optical depths, over-ionicization dominates and all low-lying levels become underpopulated. The importance of strong line scattering was tested by excluding the resonance lines and subordinate lines at 1 μm from the model atom. The overpopulation of the 5p^2P^0 and 4d^2D vanished both for the metal-rich and metal-poor models.

### 3.3. NLTE abundance corrections

To quantify the effect of NLTE on the abundance determinations, we use the concept of the NLTE abundance correction Δnlte, where:

\[
\Delta_{\text{NLTE}} = \log \epsilon(\text{Sr})_{\text{NLTE}} - \log \epsilon(\text{Sr})_{\text{LTE}}
\]

is the the logarithmic correction, which has to be applied to an LTE abundance determination of a specific line to obtain the exact value corresponding to the use of NLTE line formation. These values were computed for a grid of model atmospheres by equalizing the NLTE equivalent widths though varying the element abundance in the LTE calculations. The NLTE abundance corrections for the 4077, 4607 and 10036 Å are given in Tables 3 and 4 for a wide range of stellar parameters. Fig. C (top panel) also shows the values of Δnlte for selected model atmospheres as a function of effective temperature, gravity, and [Fe/H].

As evident from Fig. C (top panel) and Table 4, the NLTE abundance corrections for the 4607 Å are always positive, and they are maximum for cool giants with sub-solar metallicity, [Fe/H] < −1. This behavior can be easily understood by inspecting the plots of departure coefficients (Fig. B top right panel): at the typical depths of the resonance line formation,
log \( \tau_c \sim 0 \) to \(-0.5\), the level populations are already depleted by a factor of two compared to LTE. In the models of warmer metal-poor dwarfs, the line formation is confined to the deeper layers, log \( \tau_c \sim 0 \), where the NLTE effects in the SrI ionization balance are smaller. This is reflected in the behavior of the abundance corrections, which also decrease. \( \Delta_{\text{NLTE}} \) are comparatively small for the warmest models, corresponding to horizontal branch stars \( (T_{\text{eff}} = 6200\, \text{K} \) and log \( g = 2.2 \). Note, however, that at \( T_{\text{eff}} > 6000\, \text{K} \) and [Fe/H] \(< -1.5 \) the 4607 Å line is very weak, \( W_\lambda < 0.5\, \text{mÅ} \), and, thus, is not useful for abundance determinations. In contrast, the line could be suitable for the analysis in sufficiently good-quality (\( R > 20000\), S/N >50) spectra of giants and sub-giants down to [Fe/H] \sim -2.5.

As described in the previous section, the formation of Sr\n lines is also significantly affected by the NLTE effects. Inspection of Table 4 and Fig. 5 (middle panel) shows that for the resonance SrI line at 4077 Å the NLTE abundance corrections are typically negative for stars with metallicity [Fe/H] \sim -3, but rapidly increase for more metal-poor stars with \( T_{\text{eff}} > 5000\, \text{K} \) and/or log \( g > 3 \). The amplitude of \( \Delta_{\text{NLTE}} \) increases with \( T_{\text{eff}} \) and decreasing log \( g \). Another resonance SrI line at 4215 Å shows the same pattern and the NLTE abundance corrections are nearly identical; we caution, however, that this line is blended by an FeI line, and should be avoided in abundance analyses of cool stars. The implication is that, compared to NLTE, traditional LTE studies rely on the resonance SrI lines under-estimate the abundance of Sr in very metal-poor stars, but somewhat over-estimate it for less metal-poor objects, such as that of thick and thin disk of the Galaxy.

For the near-IR subordinate SrII lines, the NLTE abundance corrections are typically negative (Table 5 and Fig. 5, bottom panel), being as large as \(-0.2\) dex even for mildly metal-poor giants; this behavior is consistent with the discussion in Sect. 3.2 as the NLTE effects are due to strong line scattering. However, \( \Delta_{\text{NLTE}} \) are usually within \(-0.1\) dex for dwarfs and become mildly positive only at very low metallicity, [Fe/H] \(< -3 \) and \( T_{\text{eff}} > 6000\, \text{K} \). The NLTE abundance corrections presented in Table 5 are rather similar for the other two members of the near-IR triplet, i.e., the lines at 10327 and 10914 Å, and can be provided by request.

Our NLTE abundance corrections are in line with the results of Short & Hauschildt (2006), although they investigated the formation of the resonance SrI lines only. They find NLTE line strengthening for the giant models with [Fe/H] \sim -1...-2 and NLTE weakening for the models with [Fe/H] \sim -4...-5. However, we note some qualitative differences with Andrievsky et al. (2011). For example, for the near-IR SrI lines their abundance corrections at [Fe/H] = -3 and \( T_{\text{eff}} = 4800 \) \sim 5300 K are of the order of \(-0.3 \) to \(-0.5\) dex and are very sensitive to surface gravity (their Fig. 7, top panel). In contrast, our values for the 10036 Å line remain within \sim -0.1 dex for any log \( g \) value in this [Fe/H] and \( T_{\text{eff}} \) range. For the resonance SrI line at 4077 Å, our \( \Delta_{\text{NLTE}} \) are mildly negative for any log \( g \) and \( T_{\text{eff}} \) at [Fe/H] = -3, whereas Andrievsky et al. (2011) obtain large positive corrections for dwarfs and negative \( \Delta_{\text{NLTE}} \) for giants. The differences are most likely caused by the differences in the NLTE model atoms, as described in Sect. 2.2.

4. Application to the abundance analysis of late-type stars

In addition to the Sun, three other late-type stars were selected for the abundance analysis. The principle selection criteria were that stellar parameters are well-constrained by independent techniques (interferometry, parallaxes), observations cover a large wavelength interval and their quality is very high for the 4607 SrI line and the near-IR SrI triplet to be accurately measured. Stellar parameters for the warm turn-off star Procyon, moderately metal-poor giant Arcturus (HD 124897), and very metal-poor giant HD 122563 were taken from Korn et al. (2003), Triel (2009), and Mashonkina et al. (2009), respectively (Table 6).

The accuracy of these parameters was verified in our recent analysis of Fe NLTE statistical equilibrium with mean 3D model atmospheres (Bergemann et al. 2012). We stress that it is not our goal here to obtain Sr abundances for a large sample of stars, as this will be the focus of our forthcoming publications (Hansen et al. in prep., Ruchti et al. in prep.), but rather to test the performance of the new NLTE model atom.

For the spectrum synthesis (Sect. 2.1), we used the solar KPNO flux spectrum (Kurucz et al. 1984). The spectra of Arcturus (HD 124897), Procyon (HD 61421), and HD 122563 were retrieved from the UVESPOP database (Bagnulo et al. 2003). These UVES (VLT, Paranal) spectra have a slit-determined resolution of \sim 80000 and a signal-to-noise ratio S/N ~ 300 near 5000 Å.

The solar abundance was determined using five SrI lines from the Table 1 adopting \( T_{\text{eff}} = 5777\, \text{K}, \text{log}\, g = 4.44, \text{[Fe/H]} = 0 \), and \( \xi_t = 1\, \text{km/s} \). The results are largely discrepant in LTE: \( \text{log}\, \epsilon_{\text{SrI}} = 2.81 \pm 0.06 \) and \( \text{log}\, \epsilon_{\text{SrII}} = 3.06 \pm 0.1 \). The NLTE model atom successfully eliminates this problem.
The abundances from the two ionization stages are fully consistent log $\epsilon_{\text{Sr}i} = 2.91 \pm 0.03$ and log $\epsilon_{\text{Sr}ii} = 2.94 \pm 0.04$ in NLTE. Here, the error corresponds to one standard deviation of the line sample. The NLTE result is also in agreement with the meteoritic abundance of Sr, log $\epsilon = 2.90 \pm 0.03$ dex (Lodders et al. 2009). We note that the meteoritic abundance of Sr given by Asplund et al. (2009), Sr, log $\epsilon = 2.88 \pm 0.03$ dex, is slightly lower, which is a consequence of different reference Si abundances. The solar Sr abundance was also determined by Barklem & O’Mara (2000), Andrei et al. (2011), and Mashonkina & Gehren (2001). The last two references perform NLTE calculations, which are generally in agreement with our results, both in terms of NLTE abundance corrections and absolute abundances. We caution, however, that this agreement is not particularly telling, because other parameters in the modeling are different. Mashonkina & Gehren (2001) adopted $\xi_i = 0.8$ km/s and employed the older (ODF) version of the MAFAGS model atmospheres. Barklem & O’Mara (2000) recover meteoritic Sr abundance from the neutral Sr lines even in LTE, which is not unexpected as they used $\xi_i = 0.85$ km/s and the solar Holweger-Müller (1974) model atmosphere.

The LTE and NLTE Sr abundances determined for the reference stars are given in Table 6 and a line-by-line comparison is shown in Fig. 6. Few examples of synthetic and observed line profiles are shown in Fig. 7. Comparison of the [Sr/Fe] and [Sr/Fe] from Table 6 reveals that the LTE approach predicts a systematic discrepancy between the lines of two ionization stages, which is also evident from the Fig. 6. The difference between the Sr I and the near-IR Sr II lines, which are almost insensitive to damping, is of the order 0.3 to 0.5 dex. This is far beyond the uncertainties in stellar parameters, which are ~ 80 K for $T_{\text{eff}}$, 0.1 dex for log $g$, and 0.1 dex for [Fe/H]. Also, the uncertainties of the transition probabilities are far smaller than that (Sect. 3.3).

The NLTE model atom provides, in contrast, realistic ionization balance, significantly reducing the scatter between the Sr lines of different ionization stages and excitation energies (Fig. 6). In this respect, especially important is a good agreement between the NLTE abundances from the near-IR Sr II and visual resonance Sr I line, which are very sensitive to NLTE and the effects are of different nature, i.e., overionization-stipulated decrease of opacity in the 4607 line and source function depletion in the lines of the 1 $\mu$m triplet.

The analysis of a larger sample of stars supports the importance of NLTE effects in the ionization balance of Sr. This is evident from Fig. 8 which shows the difference between the mean abundances computed using the Sr I and Sr II lines for a sample of thick-disk and halo RAVE stars with high-resolution spectroscopic follow-up observations (see Ruchti et al. 2011). In LTE, the mean offset between the two ionization stages is ~ 0.4 dex and it is alleviated when the NLTE effects are taken into account. A detailed analysis of the data will be presented in Ruchti et al. (in prep.).

We thus conclude that NLTE must be taken into account in abundance analysis of resonance Sr I and subordinate Sr II lines in spectra of late-type stars. Also, at low metallicities, the resonance Sr II lines are sensitive to NLTE effects, and abundances determined in LTE are largely underestimated. It is also important to note that the NLTE effects for Sr might be larger than our estimates, because of interlocking with NLTE-affected lines of other atoms, as demonstrated by Short & Hauschildt (2006).

5. Conclusions

We investigate statistical equilibrium of Sr in the atmospheres of late-type stars. The NLTE model atom was constructed using new atomic data, computed with the R-matrix method,
which include levels, transition probabilities, photoionization and electron-impact excitation cross-sections.

The NLTE effects are significant for the Sr I resonance line at 4607 Å and are stipulated by overionization. For the resonance Sr II lines, typically observed in metal-poor FGK stars, our model predicts small deviations from LTE, although NLTE effects become important at very low metallicity, [Fe/H] < −3, reaching +0.2 dex for very metal-poor turnoff stars. The near-IR Sr II triplet shows substantial sensitivity to NLTE, and the NLTE corrections vary between +0.1 and −0.2 dex depending on the model metallicity, temperature, and surface gravity.

The NLTE model atom was applied to the analysis of the Sun and three stars with well-constrained stellar parameters. In contrast to LTE approach, the NLTE model recovers ionization and excitation equilibria of Sr for all these reference stars, thus confirming that the NLTE modeling approach developed in this work provides a solid base for future abundance determinations of late-type stars.

The NLTE abundance corrections are provided for the important diagnostic Sr I and Sr II lines for a grid of model atmospheres in the following range of stellar parameters 4400 ≤ T_eff ≤ 6000 K, 2.2 ≤ log g ≤ 4.6, −4 ≤ [Fe/H] ≤ 0; for other Sr II lines and models the abundance corrections can be also computed by request.

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Fig. 5. NLTE abundance corrections for the Sr\textit{i} and Sr\textit{ii} lines. Note different x-axis scales.
Fig. 7. NLTE and LTE line profiles of the Sr $i$ and Sr $ii$ lines for selected stars compared with observed stellar spectra.
\[
\Delta \left( [\text{SrI/Fe}] - [\text{SrII/Fe}] \right)
\]

vs.

[Fe/H]