Lithium abundances of halo dwarfs based on excitation temperature

I. Local thermodynamic equilibrium*

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ABSTRACT

Context. The discovery of the Spite plateau in the abundances of ⁷Li for metal-poor stars led to the determination of an observationally deduced primordial lithium abundance. However, after the success of the Wilkinson Microwave Anisotropy Probe (WMAP) in determining the baryon density, Ω_b h², there was a discrepancy between observationally determined and theoretically determined abundances in the case of ⁷Li. One of the most important uncertain factors in the calculation of the stellar ⁷Li abundance is the effective temperature, T_{eff}.

Aims. We use sixteen metal-poor halo dwarfs to calculate new T_{eff} values using the excitation energy method. With this temperature scale we then calculate new Li abundances for this group of stars in an attempt to resolve the ⁷Li discrepancy.

Methods. Using high signal-to-noise (S/N ≈ 100) spectra of 16 metal-poor halo dwarfs, obtained with the UCLES spectrograph on the AAT, measurements of equivalent widths from a set of unblended Fe lines are made. These equivalent widths are then used to calculate new T_{eff} values with the use of the single line radiative transfer program WIDTH6, where we have constrained the gravity using either theoretical isochrones or the Hipparcos parallax, rather than the ionization balance. The lithium abundances of the stars are calculated with these temperatures.

Results. The physical parameters are derived for the 16 programme stars, and two standards. These include T_{eff}, log g, [Fe/H], microturbulence and ⁷Li abundances. A comparison between the temperature scale of this work and those adopted by others has been undertaken. We find good consistency with the temperatures derived from the Hα line by Asplund et al. (2006, ApJ, 644, 229), but not with the hotter scale of Meléndez & Ramírez (2004, ApJ, 615, L33). We also present results of the investigation into whether any trends between ⁷Li and metallicity or temperature are present in these metal-poor stars.

Key words. Galaxy: halo – cosmology: early universe – nuclear reactions, nucleosynthesis, abundances – stars: abundances

1. Introduction

1.1. The lithium problem

The development of the Big Bang theory brought about predictions for the primordial abundances of the nuclides, ²H, ³He, ⁴He, and ⁷Li (Olive et al. 2000; Fields & Sarkar 2006). These values, determined from calculations of Big Bang Nucleosynthesis (BBN) for an assumed baryon to photon ratio η_b, are in at least partial agreement with those primordial abundances determined from observations of ²H, ³He, ⁴He, and ⁷Li relative to ¹H. Historically this has given constraints on η_b, or equivalently the baryon density fraction, Ω_b h², which is standard BBN’s one free parameter, where h is the Hubble constant in units of 100 km⁻¹ Mpc⁻¹. Precisely determining Ω_b h² by other means can narrow down the range for the calculated primordial abundances, and thus confirm whether the observationally inferred primordial abundances are consistent.

The Wilkinson Microwave Anisotropy Probe (WMAP) achieved high precision in measuring Ω_b h². Using measurements of the cosmic microwave background (CMB) radiation, the value Ω_b h² = 0.0227 ± 0.0006 was inferred (Dunkley et al. 2008). This value is in excellent agreement with that derived from BBN via the measured ²H/¹H and ⁴He/¹H ratios, although the systematic uncertainties in the latter make ⁴He a poor discriminant. However, for other primordial isotopes the results are not in such good agreement (Coc & Vangioni 2005, their Fig. 4). The baryon densities inferred from ³He and ⁷Li abundances differ significantly. ⁷Li has the greatest deviation. The deviation of the observationally inferred primordial ⁷Li

* Appendix A is only available in electronic form at the CDS via anonymous ftp to cdsarc.u-strasbg.fr (130.79.128.5) or via http://cdsweb.u-strasbg.fr/cgi-bin/qcat?J/A+A/493/601
abundance from that deduced from WMAP and from \(^3\text{H}\) via BBN has become known as the “lithium problem”.

The \(^4\text{He}\) discrepancy can be accommodated within large uncertainties concerning the mechanisms in its production and destruction in stars (Vangioni-Flam et al. 2003). It is not clear that the \(^7\text{Li}\) discrepancy can be explained as easily.

The primordial \(^7\text{Li}\) abundance was first deduced from observations of halo stars by Spite & Spite (1982) when they discovered the \(^7\text{Li}\) plateau, now known as the Spite plateau, with a mean abundance of \(n(^7\text{Li})/n(\text{H}) = 1.12^{+0.38}_{-0.72} \times 10^{-10}\). Many new observationally determined abundances have been published, the values of which can differ significantly, for example from the low value of \(n(^7\text{Li})/n(\text{H}) = (1.23^{+0.68}_{-0.12} \times 10^{-10})\) (95% confidence limits) (Ryan et al. 2000), to a decidedly larger value of \(n(^7\text{Li})/n(\text{H}) = 2.34^{+1.21}_{-0.10} \times 10^{-10}\) (Meléndez & Ramírez 2004). In comparison the abundance calculated from WMAP and from \(^2\text{H}\) via BBN is \(n(^7\text{Li})/n(\text{H}) = (4.15^{+0.49}_{-0.44} \times 10^{-10})\) (Coc et al. 2004) or \(n(^7\text{Li})/n(\text{H}) = (4.26^{+0.73}_{-0.60} \times 10^{-10})\) (Cyburt et al. 2001; Cyburt 2004). It is obvious that the theoretical and observational values are not in accord with each other. The observed \(^7\text{Li}\) abundance is a factor of 2 to 3 times lower than that given by BBN/WMAP, far beyond the stated range of systematic uncertainties from observational analysis or BBN calculations.

Several possibilities have emerged in an attempt to explain this \(^7\text{Li}\) discrepancy, but none of these, as yet, fully explains the problem. Broadly speaking these lie in the categories:

1. The discrepancy is due to unrecognised or underestimated systematic errors in the calculations of the observationally inferred primordial \(^7\text{Li}\) abundance;
2. The stars studied have destroyed some of the \(^7\text{Li}\) with which they were formed;
3. Some Galactic \(^7\text{Li}\) was destroyed before these stars formed;
4. The discrepancy is due to systematic uncertainties in the nuclear cross sections used in BBN calculation;
5. Standard BBN has failed to accurately predict the primordial abundances.

A possible solution in the second category is diffusion. Several studies have been conducted on this subject recently, e.g. Richard et al. (2005) and Korn et al. (2006). There it is suggested that the primordial \(^7\text{Li}\) is depleted in the star by diffusion, through gravitational settling, to layers in which the \(^7\text{Li}\) can not be destroyed. These models used in this process are manipulated to minimize destruction of \(^7\text{Li}\) due to nuclear burning. However, they do not completely manage to eliminate this destruction. Although these predictions do lead to a possible solution to the \(^7\text{Li}\) problem, they are still subject to questions and uncertainties. Using atomic diffusion alone does not recreate the plateau, but causes a drop in \(^7\text{Li}\) abundance in stars with \(T<_{\text{eff}} 6000\) K. It is also often hard to explain the smallness of the star to star scatter in \(^7\text{Li}\) using these processes. To solve these problems some form of turbulence at the bottom of the convective zone has to be invoked. The source of this turbulence is still uncertain, as is the amount needed, which can vary from star to star. The turbulence is also dependent on the temperature scale used. If a cooler \(T_{\text{eff}}\) scale is used then more turbulence is needed. However, observations of \(^6\text{Li}\) in the atmospheres of metal-poor turn off stars (Asplund et al. 2006; García Peréz et al. 2008) rule out the use of a more turbulent scheme, as it would destroy the \(^6\text{Li}\). It would therefore seem that the success of this solution to the lithium problem rests on an accurate \(T_{\text{eff}}\) scale, and confirmation of whether \(^6\text{Li}\) is present on the surface of metal-poor stars. However, it should be noted that the presence of \(^6\text{Li}\) is still in question. It has been shown (Cayrel et al. 2007; García Peréz et al. 2008) that a reappraisal of previously derived \(^6\text{Li}\) abundance should be undertaken. The presence of \(^6\text{Li}\) also causes problems if \(^7\text{Li}\) undergoes nuclear burning. Although the models of diffusion and turbulence are tuned so as to reduce this burning it is not clearly stated whether they eliminate it completely. As \(^6\text{Li}\) is destroyed at a lower temperature than \(^7\text{Li}\), the presence of \(^6\text{Li}\) would rule out the possibility of any models in which \(^7\text{Li}\) is destroyed.

Possible explanations in the third category are relatively new. One such mechanism involving stellar processing is that of Piau et al. (2006). However, the amount of material which this requires to be cycled through stars creates other problems; the implications of such models require further investigation.

It is of course possible that certain nuclear cross sections used in BBN calculations have been poorly determined. The effect of changing the yields of certain BBN reactions was recently considered by Coc et al. (2004). They found for example, that an increase of the \(^7\text{Be}(\alpha,\gamma)^{12}\text{Be}\) reaction by a factor of 100 would reduce the \(^7\text{Li}\) abundance by a factor of about 3 in the WMAP \(\eta\) range. This reaction has since been remeasured and precludes this solution (Angulo et al. 2005). There is also the possibility that systematic errors in the \(^3\text{He}(\alpha,\gamma)^{7}\text{Be}\) reaction are the cause of the \(^7\text{Li}\) discrepancy. This channel was considered in detail in Cyburt et al. (2004). Although the absolute value of the cross section for this key reaction is known relatively poorly both experimentally and theoretically, the agreement between the standard solar model and solar neutrino data provides additional constraints on variations in this cross section. Using the standard solar model of Bahcall et al. (2001), and recent solar neutrino data (Ahmed et al. 2004), one can exclude systematic variations of the magnitude needed to resolve the BBN \(^7\text{Li}\) problem at the \(\gtrsim 95\%\) confidence level (Cyburt et al. 2004).

The fifth category represents an interesting challenge, with several cosmological and particle physics possibilities arising which could affect BBN (Coc & Vangioni 2005). For example, the variation of the fine structure constant can induce a variation in the deuterium binding energy and could yield a decrease in the predicted abundance of \(^7\text{Li}\) (Dmitriev et al. 2004; Coc et al. 2007). The modification of the expansion rate during BBN (Salati 2003) will also affect the light element abundances. Another possibility is that gravity is not described by general relativity but is instead attracted toward general relativity during the evolution of the Universe (Damour & Nordtvedt 1993a; Damour & Nordtvedt 1993b). BBN has been extensively studied in that scenario (Damour & Pichon 1999; Coc et al. 2006). Finally, it may be that physics beyond the standard model is responsible for the post BBN processing of the light elements. One possibility recently discussed is that particle decay after BBN could lower the \(^7\text{Li}\) abundance and produce some \(^6\text{Li}\) as well (Jedamzik 2004). This has been investigated in the framework of the constrained minimal supersymmetric standard model if the lightest supersymmetric particle is assumed to be the gravitino (Feng et al. 2004; Ellis et al. 2005; Hamaguchi et al. 2007). Some models have been found which accomplish these goals (Jedamzik et al. 2006; Cyburt et al. 2006). However, all of these possibilities lack confirmation at present. Also the other four categories of explanation listed above need to be ruled out in order for the last to gain significant favour.

It is the first category of explanation that we address in this work. The largest of the uncertainties arises from the uncertain effective temperature scales for metal-poor stars, which we now examine in greater detail, before going on to calculate new effective temperatures and lithium abundances for a sample of
metal-poor main-sequence stars, which is the ultimate aim of this work.

1.2. Effective temperature \( (T_{\text{eff}}) \) scale problems

In calculating the abundances of Li, \( A(\text{Li}) \), from spectral measurements, the effective temperature is the most important atmospheric parameter, as \( A(\text{Li}) \) has a high sensitivity to temperature: \( \partial A / \partial T_{\text{eff}} \sim 0.065 \, \text{dex per 100 K} \) for halo main-sequence turnoff stars. The temperature scale adopted by different authors varies considerably. A comparison of effective temperatures adopted by Ryan et al. (2001) and Ramírez & Meléndez (2005) shows differences for very low metallicities \((\text{[Fe/H]} < -3)\) by as much as 500 K, with typical differences of \( \sim 200 \, \text{K} \). Concerning the very metal-poor star G64-37 which features in the present paper, the difference in temperature between Ryan et al. (1999), at 6240 K, and Meléndez & Ramirez (2004), at 6775 K, is 535 K. In these cases Ramírez & Meléndez (2005) and Meléndez & Ramirez (2004) have the hotter temperature scale, and hence they infer higher primordial abundances as noted above.

Several methods are routinely adopted in calculating the effective temperature. These include using the strong Balmer line wings (e.g. Asplund et al. 2006), photometric methods including the Infrared Flux Method (IRFM) (e.g. Ramírez & Meléndez 2005) and a spectroscopic method that utilizes the temperature dependence of the atomic-level populations, as given in local thermodynamic equilibrium (LTE) by the Boltzmann equation.

In the first method, the Balmer line wings are used as a temperature indicator because they have strong temperature sensitivities, whilst maintaining a low sensitivity to other physical parameters, i.e. \( \log g \) (Fuhrmann et al. 1992). H\( \alpha \) in particular has a high sensitivity to \( T_{\text{eff}}, \sim 0.5 \, \text{Å} \) in the width of the wings per 100 K, and is the least sensitive to other parameters. This technique is widely used, e.g. Fuhrmann et al. (1994), Barklem (2002), and Asplund et al. (2006). It assumes that the wings are formed in LTE. This assumption makes it simpler to model, thus leading to more sturdy temperature scales.

The preferred temperature scale in many studies is that derived by the IRFM. This method uses IR photometry and model fluxes to constrain the temperature. The basic idea is to compare a theoretically calculated infrared flux with an observed infrared flux to derive an angular diameter, given by the equation:

\[
\theta = 2 \left( \frac{f_{\text{obs,IR}}}{f_{\text{theo,IR}}} \right)^{1/2},
\]

where \( f_{\text{obs,IR}} \) and \( f_{\text{theo,IR}} \) are the observed and theoretical infrared fluxes respectively. This \( \theta \) can then be used, along with the integrated observed flux, to calculate an effective temperature, using the equation:

\[
T_{\text{eff}} = \left( \frac{4F_{\text{obs}}}{\sigma^2 \Theta^2} \right)^{1/4},
\]

where \( F_{\text{obs}} \) is the observed integrated flux at the Earth and \( \sigma \) is the Stefan-Boltzmann constant. This is an iterative process where the \( T_{\text{eff}} \) is used to refine the theoretical fluxes, thus refining the angular diameter and the effective temperature. In essence the method comes down to simultaneously solving Eqs. (1) and (2) to derive values of \( \theta \) and \( T_{\text{eff}} \). Due to the lower sensitivity of the infrared flux to temperature compared to shorter wavelength, when Eqs. (1) and (2) are solved simultaneously the calculated temperature for the star is better defined, and hence more accurate (see Blackwell et al. 1979a, Fig. 1). Mean random errors for recent work using this method are \( \sim 60–75 \, \text{K} \), depending on the type of star analysed, from Ramírez & Meléndez (2005), and \( \sim 70 \, \text{K} \) from Alonso et al. (1999). Temperatures from this method have been used to calibrate other photometric indices (e.g. Magain 1987; Alonso et al. 1996).

The third method usually relies on the assumption of LTE. This implies that the Boltzmann equation can be utilized in the derivation of the effective temperatures. Boltzmann’s equation contains an exponential term in \( \chi_i / T \), where \( \chi_i \) is the excitation energy of the given energy level \( i \), and \( T \) is the temperature of the gas. Each measured equivalent width (EW) of an element, in most cases FeI, is used to determine an abundance for that element. If the temperature used in the calculation is too high, then an over-population of the more excited levels will be calculated with the Boltzmann equation. This will lead to the inference that there are already more absorbers in higher levels, and hence it would then be calculated that fewer absorbers were needed to reproduce the measured EW. A dependence of abundance on excitation energy would therefore be seen due to these miscalculated populations. The correct temperature will therefore be the one that nulls any trend in the plot of abundance against excitation energy (e.g. Peterson & Carney 1979).

However, all methods have, along with their strengths, significant deficiencies. Barklem (2007) has questioned whether the assumption that the wings of the Balmer lines form under conditions of LTE is reliable. This therefore casts doubt on the accuracy of the derived temperatures, with a possible temperature rise of 100 K compared to those derived in LTE. Along with this there are also uncertainties in the way hydrogen-hydrogen collisions, so-called self-broadening, are treated. Barklem et al. (2000) found differences in \( T_{\text{eff}} \) of up to 150 K between the their theory and the Ali-Griem treatment of self-broadening. This difference was also found by Bonifacio et al. (2007) using a Ali-Griem theory which has modified broadening coefficients, which allow for a good match to solar Balmer lines. The IRFM relies heavily on photometry and modelled fluxes to derive the temperatures. It can therefore suffer from photometric errors. This method also has a dependency on model atmosphere. Finally, temperatures
derived using the excitation energy technique currently depend on the accuracy of assumptions about LTE, about the structures of 1D model atmospheres and the errors in $gf$ values, damping values and equivalent width measurements. It has been shown (Thévenin et al. 1999; Asplund et al. 1999) that the assumptions of 1D, LTE atmospheres are not always suitable for modelling radiative transfer of Fe lines in metal-poor stars, although Gratton et al. (1999) infer there is little sensitivity to non-LTE. It is thus unclear how strong these effects are.

In previous work by some of the present authors (Ryan et al. 1999), temperatures were derived using a calibration of BVRi colours and medium resolution ($\Delta \lambda \approx 1 \AA$) spectroscopy, tied principally to the Magain (1987) $b-y$ calibration of IRFM temperatures. In Ryan et al. (2000), an adjustment of $+120$ K was made to the temperature scale to make it similar to the IRFM scale of Alonso et al. (1996). Alonso et al. (1996) report similarly a mean difference of $112$ K between their IRFM values and Magain’s. The aim of the present work was to explore the repercussions of the excitation energy technique to further constrain the effective temperature scale. In other studies where excitation-energy temperatures have been derived, only limited attention has been paid to quantifying the random and systematic uncertainties on those values. Given the importance of the temperature scale to the lithium problem, we endeavor here to track the uncertainties more closely.

1.3. 1D-LTE vs. 3D/1D-non-LTE

Because the calculation of an effective temperature from the excitation-energy technique requires that the populations of the different energy levels be known, it is necessary to investigate non-LTE effects in metal-poor stars and their effect on the temperature scale calculation.

As discussed by Asplund et al. (1999) the level populations for the Fe lines are affected by departures from LTE, as radiation fields higher than the local Planck function lead to over-ionisation of the atom compared to the LTE case. That is, applying Saha’s equation with the local value of the temperature at some depth in the atmosphere will underestimate the degree of ionisation. Moreover, different energy levels will be affected differently by non-LTE, so it is not possible to apply a single ionisation correction to all levels of a given ionisation state. A second factor that could affect excitation equilibrium is that of the use of 3D atmospheres. It can be seen in Asplund (2005) (his Fig. 8) that there can be a significant $\chi$-dependent effect on the abundances produced. However, in the same paper it is stated that non-LTE corrections appear to have the opposite effect to the 3D corrections and these corrections also become larger when 3D atmospheres are used. This could lead to 3D non-LTE becoming very similar to 1D non-LTE. It is not known, however, that this is always the case. Non-LTE effects, 3D effects, and their combination relative to 1D-LTE will therefore have to be studied to assess the exact impact on the temperatures derived by this technique.

However, it is well known that non-LTE calculations for Fe I in metal-poor stars are quite uncertain, principally because of uncertainty in the role of collisions with hydrogen. Thévenin & Idiart (1999) and Gratton et al. (1999) reached opposite conclusions about the importance, or otherwise, of non-LTE in such stars. 3D calculations are also relatively rare to date. Because of these uncertainties, we proceed in two stages. For the present paper we shall concentrate on a 1D LTE analysis, and this will be the focus for the remainder of this paper. In a second paper we will examine the complexities introduced by attempting a non-LTE analysis.

2. Observations

2.1. Sample selection

The sample of stars was taken from one constructed by Ryan et al. (1999) which contains 24 metal-poor stars spread around the sky. It is a carefully selected sample at the Population II main-sequence turnoff, with as few star-to-star variations as possible. This included limiting the effective temperature to a range of $6100$ K $\leq T_{\text{eff}} \leq 6300$ K and the metallicity to $-3.5 \leq [\text{Fe/H}] \leq -2.5$. These restrictions allow the sample to have roughly the same evolutionary state, in this case around the MS turn-off, which also has the effect of limiting log $g$. The low metallicity restrictions mean that we are sampling stars that contain material that has undergone a limited amount of processing since the Big Bang. The sample was previously studied to determine not only Li abundances, but also those of magnesium (Arnone et al. 2005), based on red spectra.

The intention in the present work was to obtain blue high–resolution spectra for a subset of these stars. This was to enable us to use the excitation-energy technique to determine the $T_{\text{eff}}$, and determine what, if any, correction is required to the photometric $T_{\text{eff}}$ scale used previously. This also allows for a comparison between the excitation-energy $T_{\text{eff}}$ and those scales used by Asplund et al. (2005) and Meléndez & Ramírez (2004).

2.2. Data acquisition and reduction

Spectra of 16 of the program stars plus two standards, out of the 22 program stars and two standards, HD140283 and HD74000, in the Ryan et al. (1999) study, were obtained using the UCLES instrument on the 3.9 m Anglo-Australian Telescope (AAO) during two separate observing runs. The first run, July 31-August 3 2006, was undertaken by Ryan & Tanner. The second run, March 9–11 2007, was undertaken by Hosford and García Pérez. The setup was the same for both observing runs: the 79 lines/mm grating was used, centred at $4174$ Å with a slit width of $1.0''$, which achieves a resolving power of $R \approx \lambda/\Delta \lambda = 43 000$. The wavelength range is from $3700 - 4900$ Å. In Table 1 we present the basic information including the photometric (and partially spectroscopic) temperature, $T_{\text{phot}}$ (Ryan et al. 1999) for each of the stars, along with the $S/N$ per 0.025 Å pixel near $4300$ Å from the present work.

The data reduction was done using standard routines within the IRAF package.

3. Data analysis

Analysis of the data proceeded in, broadly, two stages. Firstly, the equivalent widths of spectral lines of Fe were measured. For the second stage we use WIDTH6 (Kurucz & Furenlid 1978) to calculate the abundances for each line, from which the physical parameters can be deduced. The remainder of this section discusses these stages.

3.1. Equivalent width measurement and data quality checks

Equivalent widths (EW) of apparently unblended Fe I and Fe II lines in the spectrum of each star were measured by fitting Gaussian profiles using the IRAF$^\text{TM}$ "splot" program. Graphs of the
Gaussian FWHM in velocity space ($\text{FWHM}_{\lambda}$) were plotted against the measured equivalent width of the line, see Fig. 1. Due to the nature of echelle spectra, in which $\Delta \lambda / \lambda$ is constant throughout the spectrum (where $\Delta \lambda$ is the width of a pixel in wavelength units), it is possible to use these plots as checks of the quality of the measurements. The weak lines, $EW \lesssim 90 \text{mÅ}$, form a plateau ($\Delta \lambda / \lambda \sim 90 \text{mÅ}$, form a plateau and stronger lines, the plot curves upward due to the increase in their line wing strength, leading to larger FWHM. Lines whose equivalent widths fell significantly outside of these trends were re-measured or discarded as blends or noise-dominated measurements, and were rejected from further calculations. The equivalent widths measured in each star are listed in Tables A1–A3 in the appendix to this paper. (The appendix is available online only at the CDS.)

A comparison between equivalent widths of the Fe lines in the star CD-24°17504 measured by Norris et al. (2001) and re-measured by Hosford showed good precision. The standard deviation of the difference between the two sets of equivalent widths was $1.3 \text{mÅ}$ with a mean of $-0.58 \text{mÅ}$ for lines measured from the same spectra. Lines which were measured in two spectra, one obtained from Norris et al. (2001) at $S/N \approx 100$, and one obtained by A. Hosford at $S/N \approx 50$, were also compared. This comparison gave a standard deviation of $5.8 \text{mÅ}$ and a mean of $-0.39 \text{mÅ}$. To increase the accuracy of the EW measurements for the star CD-24°17504 we used a combination of EW measurements from both the Norris et al. (2001) spectra and the Hosford spectrum, weighting each measurement by the inverse of the uncertainty in the EW ($\sigma_{\text{EW}}$). This $\sigma_{\text{EW}}$ is calculated using the equation:

$$\sigma_{\text{EW}} = \frac{\Delta \lambda_{\text{pix}} \sqrt{M}}{S/N},$$

where $\Delta \lambda_{\text{pix}}$ is the width of a pixel in wavelength units, $M$ is the number of pixels comprising the full width of the spectral line and $S/N$ is the signal-to-noise of the spectra.

### 3.2. Abundance calculations

To calculate the abundances from the measured lines we use the single line radiative transfer program WIDTH96. This program calculates the EW of an unblended absorption line for a 1D model atmosphere. It varies the starting abundance, iteratively calculating an EW until that equals the measured EW for the line.

The KURUCZ06 (http://kurucz.harvard.edu/grids.html, downloaded in Nov. 2006) model atmospheres have been adopted for this work as it has been stated (Castelli et al. 1997) that these no-overshooting models are preferable in deriving accurate temperatures to earlier Kurucz overshooting models (Kurucz 1993). This is due to the effect that the different methods of treating convection have on the temperature structure of the atmosphere, and therefore on the spectral features. The no-overshooting models lead to a better fit to observable spectral features at different temperatures. These models are interpolated in $T_{\text{eff}}$, $\log g$ and [Fe/H] over the ranges $3500 \leq T_{\text{eff}} \leq 50000$, $0.0 \leq \log g \leq 5.0$ and $-0.5 \geq [\text{Fe/H}] \geq -4.0$. A comparison between temperatures derived using the different overshooting models can be seen in Sect. 3.3.

The log $gf$ values of the FeI lines were compiled from several sources: Blackwell et al. (1979b,c), Blackwell et al. (1980), Blackwell et al. (1982a,b), collectively the “Oxford FeI consortium”, Fuhr et al. (1988), Bard et al. (1991), O’Brian et al. (1991), Bard & Kock (1994) and Thévenin et al. (1989); in some cases the mean of several log $gf$ values is adopted, see Table A1 for details. The FeII log $gf$ values are compiled from: Hannaford et al. (1992), Heise & Kock (1990), Kroll & Kock (1987), Moity (1983) and Schnabel et al. (1999). Values given by Moity (1983) are corrected by $-0.06 \text{dex}$, for those lines with upper-level energy $<48000 \text{cm}^{-1}$, and $-0.24 \text{dex}$ for lines with upper-level energy greater than the above value; this follows the suggestion by Fuhr et al. (1988). Damping values have been calculated using the tabulations of Anstee & O’Mara (1995), Barklem & O’Mara (1997), and Barklem et al. (1998). Ryan (1998) noted that the damping treatment has a significant effect on the excitation energy temperatures because of the $\chi$-dependence of damping treatments.

### Table 1. Background information for the stars analysed in this work.

<table>
<thead>
<tr>
<th>Star name</th>
<th>RA (1950)</th>
<th>Dec (1950)</th>
<th>$V$</th>
<th>$B-V$</th>
<th>$S/N$@4300</th>
<th>$T_{\text{phot}}$ (K)</th>
<th>Initial [Fe/H]</th>
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<td>19°51′24″</td>
<td>11.20</td>
<td>0.40</td>
<td>95</td>
<td>6200</td>
<td>−2.64</td>
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<td>24°11′42″</td>
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<td>6170</td>
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<td>6150</td>
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<td>04°02′24″</td>
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<td>100</td>
<td>6240</td>
<td>−2.70</td>
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<tr>
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<td>08°51′24″</td>
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<td>6250</td>
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</tr>
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</tr>
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</tr>
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<td>110</td>
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</tr>
<tr>
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<td>−05°24′18″</td>
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<tr>
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<td>150</td>
<td>6040</td>
<td>−2.02</td>
</tr>
</tbody>
</table>
ties are 30 K and 0.04 dex respectively; using di
0.5 dex respectively.

3.3. Adopted approach to constraining parameters

Several other physical parameters of a star, in addition to tempera-
ture, could in principle be constrained using WIDTH6, given appro-
propriate data. These are log $g$, metallicity and microturbulence. Metallicity is determined from the calculated Fe abundance. Log $g$ could be constrained by ensuring that the abundances derived for Fe I and Fe II lines are equal, and microtur-
bulence is found by making sure there is no trend between the mea-
sured equivalent widths of the lines and the abundance.

It would, in principle, seem possible to simultaneously con-
strain both log $g$ and $T_{\text{eff}}$. However, as previously mentioned, there are concerns about the non-validity of the assumption of LTE for Fe I lines in metal-poor turnoff stars. This would therefore make it dangerous to try to constrain $T_{\text{eff}}$ at the same time as log $g$. This concern became evident when $T_{\text{eff}}$ and log $g$ where simultaneously constrained via a WIDTH6 analy-
sis for the star HD140283, using Fe I and Fe II lines. For this star results converged at $T_{\text{eff}}$/log$g/[\text{Fe/Fe}]/\xi = 5753/3.1/ -2.67/1.5$ using the KURUCZ06 models\(^2\). However, we know that the Hipparcos parallax gives a reliable value for log $g$ for HD140283: $\log g = 3.73 \pm 0.11$ for a mass of 0.8360 $M_\odot$ as taken from the Yonsei-Yale isochrones (http://csaweb.
yonsei.ac.kr/~kim/yyiso.html). The uncertainty is domi-
nated by the parallax error. Clearly the log $g$ value we have de-

erived is too low. This star has also appeared in many other stud-
ies leading to temperatures around 5750 K (Table 2), approxi-
mately 200 K hotter than we derived. It can then be seen that for this star, attempting to constrain $T_{\text{eff}}$ and log $g$ simultane-
ously from Fe I and Fe II lines drives the gravity and possibly the tem-
perature values down to unusually low levels. Non-LTE effects may be possible. It is, however, still viable to use WIDTH6 to constrain $T_{\text{eff}}$. This is because the main non-LTE correction in-
volved in Fe is expected to be overionisation, which may sig-
nificantly affect the abundance of Fe I and therefore have a sub-
stantial effect on log $g$ when calculated through ionisation bal-
ance. The correction on $\epsilon$ is very roughly the same for all levels within a particular star. This can be seen from Figs. 2 and 3 from Collet et al. (2005) while Asplund (2005) also states this view. Shchukina et al. (2005) do find that the effect can be different for different excitation potentials, however, this seems far more pronounced when using 3D atmospheres, compared to using 1D atmos-
pheres. Therefore overionisation will have less affect on the trend between [Fe/Fe] and $\chi$, and it is still acceptable to use this to constrain $T_{\text{eff}}$, subject to the caveats in Sect. 1.3.

It is therefore preferable to constrain log $g$ independently of deducing $T_{\text{eff}}$, rather than using Fe lines to constrain both. There are several options for doing this. In rare cases the Hipparcos parallax could be used. However a large proportion of the

![Fig. 2. Theoretical isochrones used in the determination of log $g$. Different ages are represented in the top panel, with the dashed line 13 Gyr, the solid line 12 Gyr, and the dot dashed line 11 Gyr. The bottom panel gives examples for different [Fe/H]. The dashed line is $-3.3$, the solid line $-2.5$ and the dot dashed line $-2.15$.](image_url)

<table>
<thead>
<tr>
<th>Study</th>
<th>Temperature</th>
<th>$T_{\text{eff}}$/log $g$</th>
</tr>
</thead>
<tbody>
<tr>
<td>This work</td>
<td>Simultaneous fit</td>
<td>5753/3.1/ -2.67/1.5</td>
</tr>
<tr>
<td>Asplund et al. (2006)</td>
<td>Balmer line wing</td>
<td>5753/3.7/</td>
</tr>
<tr>
<td>Ryan et al. (1996b)</td>
<td>Photometry</td>
<td>5750/3.4</td>
</tr>
<tr>
<td>Alonso et al. (1996)</td>
<td>IRFM</td>
<td>5691/4.0</td>
</tr>
</tbody>
</table>

\(^a\) The uncertainties are 75 K and 0.15 dex respectively; \(^b\) the uncertainties are 30 K and 0.04 dex respectively; \(^c\) the uncertainties are 69 K and 0.5 dex respectively.
parallax, and so we quote at most one value of \( \log g_{\text{Hipp}} \) for each star. Many of the stars’ parallaxes have extremely large errors, to the point where they would have been unsatisfactory to use, and are excluded. In the case of the HD stars, we have taken the log \( g_{\text{Hipp}} \) as the final gravity as they have smaller errors. However, for the other five stars with log \( g_{\text{Hipp}} \) in Table 3, we have taken log \( g_{\text{Hipp}} \) to be only an interim value. A theoretical gravity, log \( g_{\text{iso}} \), is then used to determine a temperature as described below. We have used \( T_{\text{phot}} \) to calculate the Hipparcos gravity, log \( g_{\text{Hipp}} \), but note that it has only a low sensitivity to \( T_{\text{phot}} \) of +0.04 dex per 100 K.

We also run WIDTH6 with a fixed \( T_{\text{phot}} \) (see Table 1) to find an interim (ionization-balance) log \( g \), which we call log \( g_{\text{phot}} \). The results of this analysis are found in Table 3. The sensitivity of log \( g_{\text{phot}} \) to \( T_{\text{phot}} \) is +0.15 dex per 100 K. The difference between the gravities, log \( g_{\text{phot}} \)-log \( g_{\text{Hipp}} \), ranges from −0.18 to −0.52, with a mean of −0.37. This suggests that this method, using the \( T_{\text{phot}} \) and the Fe ionisation balance, underestimates the gravity by ~0.4 dex. This is due to \( T_{\text{phot}} \) being too cool by ~270 K, or may be evidence of non-LTE effects (overionisation) affecting Fe I. We make no further use of the log \( g_{\text{phot}} \).

Because of the lack of reliable Hipparcos parallaxes, the unreliability of log \( g_{\text{phot}} \), and concerns about trying to constrain \( T_{\text{eff}} \) and log \( g \) simultaneously with WIDTH6, we then switched to using the log \( g_{\text{iso}} \) implied by the Yonsei-Yale theoretical isochrones. These isochrones can be seen in Fig. 2. Three different metallicities are plotted; the other available metallicities are \([\text{Fe/H}]=−2.3\) and −2.76. We use \( T_{\text{phot}} \) to obtain log \( g_{\text{iso}} \), and then null the trends of abundance with excitation energy to obtain the excitation energy temperature, \( T_{\text{e}} \). In most cases this leads to two scales for temperature: a MS value and a SGB value. Only if the Hipparcos gravity is sufficiently accurate can we be sure of the evolutionary state.

An initial comparison between results calculated using the KURUCZ93 models, with overshoot, and those for the new KURUCZ06 models, with no-overshooting, was carried out for the star HD140283. This was done using the calculated Hipparcos gravity for the star, log \( g = 3.73 \). It was found that the KURUCZ06 models gave \( T_{\text{e}} \) hotter by 41 K than the KURUCZ93 models. We decided that for the remainder of the analysis the KURUCZ06 models would be used; see Sect. 3.2 for a discussion on the reasons for this.

### 4. Results

The physical parameters for the sixteen programme stars and the two standards are given in Table 4.

The effective temperatures for all but the HD stars in Table 4 were found using isochronal gravities, log \( g_{\text{iso}} \), based in turn on photometric temperatures. In order to understand how reliable our \( T_{\text{e}} \) determinations are, it is essential that we quantify the various sources of possible error that contribute. As stated above, we defer discussions of non-LTE effects to Paper II. One source of error is the adopted gravity, log \( g_{\text{Hipp}} \), for the HD stars and log \( g_{\text{iso}} \) for the remainder. Choice of isochrones leads to changes in \( T_{\text{e}} \) by 12 K and 24 K respectively. The isochronal gravity is also sensitive to the choice of \( T_{\text{eff}} \), for which we use \( T_{\text{phot}} \). An error of +100 K in \( T_{\text{phot}} \) would produce an error of typically +0.06 dex in log \( g_{\text{iso}} \) in the SGB gravity, and −0.06 dex in the MS gravity, though more at the turnover. This in turn produces an error of typically +24 K in the SGB \( T_{\text{e}} \), and −24 K in the MS \( T_{\text{e}} \). Hence, and importantly, the inferred \( T_{\text{e}} \) is only weakly dependent on an imperfect value of \( T_{\text{phot}} \). Concerning microturbulence, we see that an error of 0.1 km s\(^{-1}\) leads to, on average, an error of ~60 K in \( T_{\text{e}} \). Typical errors in microturbulence for this work range from ~0.05–0.15 km s\(^{-1}\).

The major contributions to the uncertainty in \( T_{\text{e}} \) comes from the procedure of nulling the dependence of [\text{Fe/H}] on \( \chi \). This involves making a linear least squares fit to the \( \chi \) pairs of \( \chi_{[\text{Fe/H}]} \) values for a star, using \( n \) Fe lines, as may be seen in Fig. 3. As there is a spread of [\text{Fe/H}] values for a given star, the slope can only be determined to a certain statistical accuracy, which turns out to be of order \( \sigma \approx 0.01–0.015 \) dex per eV, which corresponds to a temperature uncertainty of order ~60–80 K. The range in [\text{Fe/H}] values encapsulates the impact of random (not \( \chi \)-correlated) line-to-line errors in equivalent widths, \( g_f \), and damping values. We determine this value for each star, finding a particularly large value for CD-2417504 on account of the small number of Fe I lines in this very metal-poor star. The final 1σ uncertainty in \( T_{\text{e}} \), given in Table 4, includes the quadrature sum of the (dominant) statistical error arising from nulling the [\text{Fe/H}], \( \chi \) trend, and contributions for \( \Delta \chi = 1 \) Gyr, \( \Delta \chi = 0.1 \) km s\(^{-1}\), \( \Delta [\text{Fe/H}] = 0.05 \), and an allowance of \( \Delta T_{\text{phot}} = 100 \) K.

We comment briefly on the larger scale of the uncertainties quoted here, compared to the relatively small values (30–40 K) quoted for \( T_{\text{phot}} \) by Ryan et al. (1999). The aim of the previous work was to derive temperatures that minimised the impact of possible systematic errors. The purpose was to investigate the spread of Li abundances by ensuring random errors were kept to a minimum, while at the same time acknowledging that large

![Figure 3](image-url)
zero-point errors might persist but be of approximately the same size in all stars. In the current work, however, the emphasis has been to constrain the zero-point, albeit at the expense of larger random errors for individual stars. These temperatures have been used in combination with the Li equivalent widths measured by Ryan et al. (1999), and a grid of equivalent widths versus abundance for different \( T_{\text{eff}} \)'s taken from Ryan et al. (1996a), to calculate new Li abundances. The grid was constructed by determining different \( ^7\)Li equivalent widths for a number of different \( T_{\text{eff}} \) and [Fe/H] values. This was done by producing synthetic spectra of the \( ^7\)Li region with an LTE code originating from Cottrell & Norris (1978). The log \( g\phi \) values and wavelengths of the four components of \( ^7\)Li were taken from Andersen et al. (1984) and the model atmospheres were interpolated from the Bell (1981) grid. The details of the process can be found in Norris et al. (1994). Table 5 presents these values.

### 5. Discussion

#### 5.1. Comparison with other temperature scales

With the determination of the effective temperatures using the excitation method we can now compare our temperature scale to that of others, in particular to the IRFM of Meléndez & Ramírez (2004), the Balmer line wing method of Asplund et al. (2006), and the principally photometric temperatures of Ryan et al. (1999). In Fig. 4 the comparison with the IRFM \( T_{\text{eff}} \) of Meléndez & Ramírez (2004) is plotted. That scale is clearly hotter than ours, by an average of \( \sim 250 \) K. This is not to say that either scale is right, but it does lead us to the conclusion that our temperature scale will not reconcile the Li problem; the Meléndez & Ramírez (2004) IRFM scale was on the border of doing so. Our scale will give a lower mean Li abundance than will the IRFM. The difference in temperatures could have

### Table 4.

The final physical parameters deduced in this work for the main sequence (MS) scale and the sub giant (SGB) scale. The three HD stars had gravities determined by their Hipparcos parallaxes and as such their evolutionary state can be deduced.

<table>
<thead>
<tr>
<th>Star name</th>
<th>log ( g\phi ) (MS)</th>
<th>( T_{\text{eff}} ) (MS)</th>
<th>[Fe/H] (MS)</th>
<th>( \xi ) (km s(^{-1})) (MS)</th>
<th>log ( g\phi ) (SGB)</th>
<th>( T_{\text{eff}} ) (SGB)</th>
<th>[Fe/H] (SGB)</th>
<th>( \xi ) (km s(^{-1})) (SGB)</th>
</tr>
</thead>
<tbody>
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<td>BD-13° 3442</td>
<td>4.46 ± 0.14</td>
<td>6321 ± 87</td>
<td>−2.66</td>
<td>1.20</td>
<td>3.79 ± 0.14</td>
<td>6186 ± 87</td>
<td>−2.74</td>
<td>1.30</td>
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<tr>
<td>BD+20° 2030</td>
<td>4.46 ± 0.14</td>
<td>6208 ± 104</td>
<td>−2.63</td>
<td>1.00</td>
<td>3.79 ± 0.14</td>
<td>6099 ± 104</td>
<td>−2.69</td>
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<td>−2.53</td>
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<td>−2.57</td>
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<td>6233 ± 107</td>
<td>−2.67</td>
<td>1.20</td>
<td>...</td>
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<td>6172 ± 53</td>
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<td>−3.32</td>
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<td>6304 ± 90</td>
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<td>−3.29</td>
<td>1.20</td>
<td>3.69 ± 0.14</td>
<td>6181 ± 106</td>
<td>−3.30</td>
<td>1.40</td>
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<td>LP635-14</td>
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<td>6319 ± 114</td>
<td>−2.48</td>
<td>1.50</td>
<td>3.83 ± 0.14</td>
<td>6355 ± 114</td>
<td>−2.61</td>
<td>1.50</td>
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<tr>
<td>LP815-43</td>
<td>4.40 ± 0.14</td>
<td>6529 ± 107</td>
<td>−2.61</td>
<td>1.40</td>
<td>3.87 ± 0.14</td>
<td>6400 ± 107</td>
<td>−2.68</td>
<td>1.40</td>
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<tr>
<td>HD140283</td>
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<td>...</td>
<td>...</td>
<td>...</td>
<td>3.73 ± 0.12</td>
<td>5769 ± 39</td>
<td>−2.54</td>
<td>1.50</td>
</tr>
<tr>
<td>HD74000</td>
<td>4.03 ± 0.18</td>
<td>6070 ± 127</td>
<td>−2.20</td>
<td>1.20</td>
<td>...</td>
<td>...</td>
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<td>...</td>
</tr>
</tbody>
</table>
several possible explanations. It could be due to problems with our use of LTE in the excitation method, causing it to calculate lower temperatures. An incorrect bolometric flux, F_Bol, calibration used in the IRFM could lead to that method having higher temperatures. It has also been noted (Alonso et al. 1996) that the errors in determining the absolute IR flux calibration have different effects on the derived IRFM temperatures, depending on what photometric band is used. The effect of these different errors on the derived temperatures is to move the zero point of the temperature scale. This could be another reason that the Meléndez & Ramírez (2004) temperatures are hotter.

The comparison between the temperatures of Asplund et al. (2006) and this work is plotted in Fig. 5. Here we see that the two scales are comparable. As discussed in Sect. 1.2 both of these methods may suffer from LTE effects. A departure from LTE of only a few percent would increase Balmer wing temperatures by of order 100 K (Barklem 2007). It is speculated that the use of LTE in the treatment of the Fe atom may be inaccurate, e.g. overionisation of the element relative to LTE may occur (Thévenin et al. 1999). If true this would also lead to the excitation temperatures changing, as they are based on an LTE treatment. As yet the value of the temperature change is not determined and the sensitivity of each method to non-LTE may be very different. This is an aspect that we intend to look into in future work. It is clearly of interest to know whether non-LTE corrections to these scales could increase both to be akin to that of the IRFM scale, and push the mean Li abundances toward that of the WMAP inferred abundances. This agreement

Table 5. The final lithium abundances calculated with the excitation energy temperatures (T_e).

<table>
<thead>
<tr>
<th>Star name</th>
<th>EW (Li)</th>
<th>T_e (K) (MS)</th>
<th>A(Li) (MS)</th>
<th>T_e (K) (SGB)</th>
<th>A(Li) (SGB)</th>
</tr>
</thead>
<tbody>
<tr>
<td>BD-13° 3442</td>
<td>21.0</td>
<td>6321 ± 87</td>
<td>2.20 ± 0.057</td>
<td>6186 ± 87</td>
<td>2.10 ± 0.057</td>
</tr>
<tr>
<td>BD+20° 2030</td>
<td>20.5</td>
<td>6208 ± 104</td>
<td>2.11 ± 0.068</td>
<td>6099 ± 104</td>
<td>2.03 ± 0.068</td>
</tr>
<tr>
<td>BD+24° 1676</td>
<td>21.1</td>
<td>6296 ± 55</td>
<td>2.18 ± 0.009</td>
<td>6220 ± 55</td>
<td>2.13 ± 0.009</td>
</tr>
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<td>BD+26° 2621</td>
<td>22.5</td>
<td>6233 ± 107</td>
<td>2.17 ± 0.070</td>
<td>6188 ± 105</td>
<td>2.10 ± 0.070</td>
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<td>BD+3° 740</td>
<td>19.5</td>
<td>6344 ± 113</td>
<td>2.18 ± 0.074</td>
<td>6188 ± 113</td>
<td>2.07 ± 0.074</td>
</tr>
<tr>
<td>BD+9° 2190</td>
<td>14.6</td>
<td>6486 ± 129</td>
<td>2.14 ± 0.084</td>
<td>6352 ± 129</td>
<td>2.05 ± 0.084</td>
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<td>CD-24° 17504</td>
<td>18.1</td>
<td>6102 ± 232</td>
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<td>CD-33° 1173</td>
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<td>2.24 ± 0.025</td>
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<tr>
<td>CD-71° 1234</td>
<td>25.9</td>
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<td>G64-12</td>
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<td>G64-37</td>
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<td>6070 ± 127</td>
<td>2.05 ± 0.083</td>
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Fig. 4. Comparison of the Meléndez & Ramírez (2004) temperatures and the temperatures from this work. Filled circles represent SGB temperatures and open circles represent MS temperatures. Those stars of unknown evolutionary state have two points on the plot.

Fig. 5. Comparison of the Asplund et al. (2006) temperatures and those from this work. Filled circles represent SGB temperatures and open circles represent MS temperatures. Those stars of unknown evolutionary state have two points on the plot.
between the two scales also gives an indirect comparison with the Balmer wing temperature scale of Bonifacio et al. (2007).

Finally, Fig. 6 presents the comparison between the photometric temperatures of Ryan et al. (1999) and \( T_\text{eff} \) of this work. At lower temperatures, \( T_\text{eff} < 6250 \) K, the two scales are in excellent agreement. For the MS scale we have seven stars cooler than 6250 K. The difference, \( T_\text{eff} - T_\text{phot} \), ranges from –126 K to 56 K, with a mean of –19 K. For the SGB scales we have eleven stars cooler than 6250 K, with the difference ranging from –99 K to 47 K and a mean of –18 K. However, at higher values, \( T_\text{eff} \geq 6250 \) K, we see that the excitation scale becomes hotter. In this case, for eight MS scale stars we have a range of 87 K to 262 K and a mean of 143 K, and for three SGB scale stars a range of 57 K to 117 K with a mean of 73 K. This suggests that the \( T_\text{phot} \) values are reliable for \( T_\text{eff} \leq 6250 \) K, but may be \( \sim 70-140 \) K too cool for \( T_\text{eff} \geq 6250 \) K. Photometric temperature errors can be induced through inaccurate dereddening, inaccurate calibration to standards, or a flawed colour-temperature calibration. The last of these is perhaps the most readily encountered; earlier in this work we cited differences of \( \sim 100 \) K between the IRFM scales of Magain (1987) and Alonso et al. (1996). Of course, the difference could also be due to inaccuracies in the model atmospheres used in this work to derive \( T_\text{eff} \).

5.2. Lithium abundances

Concerning the lithium problem, with our temperature scales we achieve a mean Li abundance of \( A(\text{Li}) = 2.16 \) dex for the MS scale and \( A(\text{Li}) = 2.10 \) dex for the SGB scale, with a scatter of 0.074 dex and 0.068 dex respectively. The mean Li abundance determined from the five stars whose evolutionary state is defined is \( A(\text{Li}) = 2.18 \) dex with a scatter of 0.038 dex. It is pleasing to note the similarity between our mean \(^7\text{Li} \) abundances and that derived by Spite et al. (1996), \( A(\text{Li}) = 2.08 \pm 0.08 \) dex, who also use a temperature scale based on iron excitation. Bonifacio et al. (2007) also found a similar result, \( A(\text{Li}) = 2.10 \pm 0.09 \) dex, with their Balmer line wing temperature scale.

These values are obviously still too low to reconcile with the WMAP value of \( A(\text{Li}) \approx 2.62 \) dex. Even if we later find that non-LTE corrections give temperature rises of roughly 150 K, we would get mean lithium abundances of \( A(\text{Li}) \approx 2.26 \) dex and \( A(\text{Li}) \approx 2.20 \) dex for the MS and SGB scales respectively. This value is still too low. However, it is not yet known what is the true effect of non-LTE and/or 3D corrections on our temperature scale.

Although the primary purpose of this analysis was to investigate possibly large systematic errors in temperature, we can also investigate whether there is evidence for dependencies of \( A(\text{Li}) \) on \( T_\text{eff} \) and \([\text{Fe}/\text{H}]\) within the sample. With the \( A(\text{Li}) \) values from Table 5 and the physical parameters calculated for each star we are able to perform a multiple regression fit, linear in \( T_\text{eff} \) and \([\text{Fe}/\text{H}]\). In these fits we include the stars with known evolutionary state, three on the MS and two on the SGB, combined with the eleven remaining stars of unknown evolutionary state for which we assess two cases: all MS and all SGB. This produces the equation

\[
A(\text{Li}) = 1.36(\pm 0.84) + 0.00019(\pm 0.00013)T_\text{eff} \\
+ 0.138(\pm 0.048)[\text{Fe}/\text{H}]
\]

for the MS case and

\[
A(\text{Li}) = 0.79(\pm 1.22) + 0.00025(\pm 0.00019)T_\text{eff} \\
+ 0.095(\pm 0.055)[\text{Fe}/\text{H}]
\]

for the SGB case. Here the bracketed values are 1σ errors in the coefficients. The large errors seen in the intercept coefficient comes from the large extrapolation to \( T = 0 \) and \([\text{Fe}/\text{H}] = 0\). We see no statistically significant trend in either \( T_\text{eff} \) or \([\text{Fe}/\text{H}]\) for the SGB as each coefficient has errors of similar size to the coefficient itself. Although there is the possibility of a trend existing in \([\text{Fe}/\text{H}]\) for the MS case (the coefficient is \( \sim 3\sigma \)) there is no certainty in this as it is not known whether all the stars belong in the MS evolutionary state.

The lack of significant trends here, in contrast to the significant \([\text{Fe}/\text{H}]\) dependence in the study by Ryan et al. (1996b) and as Asplund et al. 2006, comes about because the random uncertainties on the \( T_\text{eff} \) values are much larger than achieved by Ryan et al. (1999) using photometry. We recall that the primary purpose of the present work was to search for potentially large systematic errors in the temperatures, and for that purpose the larger random errors here are tolerable. However, the larger random errors greatly undermine the second use of the dataset.

We have also performed single parameter regression fits in \( T_\text{eff} \) and \([\text{Fe}/\text{H}]\). Figure 7 presents the results of the fits to temperatures. The values of the temperature coefficients imply

\[
A(\text{Li}) = 0.51(\pm 1.00) + 0.00026(\pm 0.00016)T_\text{eff}
\]

for the MS case, and

\[
A(\text{Li}) = 0.71(\pm 1.33) + 0.00022(\pm 0.00021)T_\text{eff}
\]

for the SGB case. There is clearly no statistically significant trend with temperature in either the MS or the SGB result, again since the temperature coefficient errors are of similar size to the coefficients themselves. For the fit to \([\text{Fe}/\text{H}]\) we obtained the equation

\[
A(\text{Li}) = 2.58(\pm 0.14) + 0.152(\pm 0.049)[\text{Fe}/\text{H}]
\]
for the MS case, and
\[ A(\text{Li}) = 2.35(\pm0.16) + 0.089(\pm0.056)[\text{Fe}/\text{H}] \]

for the SGB case. Figure 8 shows a plot of these fits. Here we confirm the similarity in values of the metallicity coefficient with Ryan et al. (1999) and Asplund et al. (2006), but again the large errors for the SGB case do not usefully constrain a metallicity trend. The metallicity coefficient for the MS case is \( >3\sigma \), but we caution again that we cannot be sure that all stars are on the MS.

We have also used the fitting form as described by Ryan et al. (2000) to deduce the primordial Li abundance. This fit is of the form:
\[ \frac{\text{Li}}{H} = a' + b'\text{Fe}/\text{Fe}_\odot \]

where \( a' \) is the intercept and measures the primordial abundance directly. From this we obtain a value of \( ^7\text{Li}/H = (1.18 \pm 0.10) \times 10^{-10} \) for the MS case and \( ^7\text{Li}/H = (1.10 \pm 0.1) \times 10^{-10} \) for the SGB case. Comparing this to the value deduced from WMAP via BBN, \( ^7\text{Li}/H = 4.15^{+0.49}_{-0.45} \times 10^{-10} \), we once again see we have not reconciled the lithium problem.

We can see from this work that it has not been possible to find a solution to the lithium problem through the use of temperatures derived using the excitation technique. The \( T_e \) values are lower than the Meléndez & Ramírez (2004) values, and more similar to the Hα and photometric temperature estimates. It would seem that one of the biggest outstanding problems is the assumption of LTE in the calculations. As has been stated in previous sections, we have begun our own investigation into the effects of non-LTE on abundance analysis, and therefore the effective temperatures, of this group of stars.

6. Conclusions

The physical parameters of 16 program stars, and two standard stars, have been calculated using WIDTH96. In particular, we have derived excitation energy temperatures by nulling the dependence of abundance on excitation potential for Fe I lines. We have compared our temperatures with those calculated in other works using different techniques, the IRFM of Meléndez & Ramírez (2004), the Balmer line wing method of Asplund et al. (2006), which is similar to that of Bonifacio et al. (2007), and the photometric method of Ryan et al. (1999).

We have shown that the IRFM scale of Meléndez & Ramírez (2004) is hotter than ours by an average of \(~250\) K. This difference may be the effect of the \( F_{\text{bol}} \) calibration used in IRFM or it could be a shift of the zero-point as an effect of the errors induced by using different photometric bands in the IRFM calculations. On the other hand it may be the effects of assuming LTE on our temperature scale.

It has been found that our methods have produced temperatures comparable with those derived using the Balmer line wing method. Both of these scales are too low to reconcile the lithium problem. It has been noted that the LTE assumptions of the Balmer line wing method may be leading to temperatures that are on order of 100 K too low. The method used in this work may also suffer from problems with LTE, but the full effect of non-LTE on the excitation method used here is unknown. This is something we will research in future work.

Finally we also see comparable results with the photometric temperature scale, although at higher temperatures, \( T_e \geq 6250\) K, there is a shift towards our scale becoming hotter. This suggests that the photometric temperatures are reliable up until \(~6250\) K. The most readily detected error in photometric calculations comes from the colour-temperature calibration and differences in values derived from different calibrations can be as much as \(~100\) K.

With our derived temperatures we have calculated the Li abundances of the program stars. Due to large uncertainties, we do not see statistically compelling trends in \( A(\text{Li}) \) with either \( T_{\text{eff}} \) or \([\text{Fe}/\text{H}]\). Due to the uncertainty in evolutionary state for the majority of the stars we have two mean lithium abundances, \( A(\text{Li}) = 2.16 \) dex for the main sequence case and \( A(\text{Li}) = 2.10 \) dex for the SGB case. Five of the stars do have defined evolutionary states. The mean Li abundance from these five stars is \( A(\text{Li}) = 2.12 \) dex. These values are lower than the WMAP determined value of \( A(\text{Li}) \approx 2.62 \) dex. For the observationally deduced value to be comparable with the WMAP value our temperatures would need to increase by approximately 700 K.

Fig. 7. Li abundances as a function of temperature. The dotted line is the fit to the MS temperatures (open circles) and the dashed line is the fit to the SGB temperatures (filled circles). Those stars of unknown evolutionary state have two points on the plot.

Fig. 8. Li abundances as a function of metallicity. The dotted line is the fit to the MS metallicities (open circles) and the dashed line is the fit to the SGB metallicities (filled circles). Those stars of unknown evolutionary state have two points on the plot.
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