Cumulative theoretical uncertainties in lithium depletion boundary age

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ABSTRACT

We performed a detailed analysis of the main theoretical uncertainties affecting the age at the lithium depletion boundary (LDB). To do that we computed almost 12 000 pre-main sequence models with mass in the range [0.06, 0.4] M⊙ by varying input physics (nuclear reaction cross-sections, plasma electron screening, outer boundary conditions, equation of state, and radiative opacity), initial chemical elements abundances (total metallicity, helium and deuterium abundances, and heavy elements mixture), and convection efficiency (mixing length parameter, αML). As a first step, we studied the effect of varying these quantities individually within their extreme values. Then, we analysed the impact of simultaneously perturbing the main input/parameters without an a priori assumption of independence. Such an approach allowed us to build for the first time the cumulative error stripe, which defines the edges of the maximum uncertainty region in the theoretical LDB age.

We found that the cumulative error stripe is asymmetric and dependent on the adopted mixing length value. For αML = 1.00, the positive relative age error ranges from 5 to 15%, while for solar-calibrated mixing length, the uncertainty reduces to 5 − 10%. A large fraction of such an error (≈ 40%) is due to the uncertainty in the adopted initial chemical elements abundances.

Key words: Methods: numerical – stars: abundances – stars: evolution – stars: fundamental parameters – stars: low-mass – stars: pre-main sequence

1 INTRODUCTION

Lithium burning via the proton capture 7Li(p, α)α becomes efficient in stellar conditions as temperature rises above 2.5 × 106 K. In the mass range 0.06-0.4 M⊙ (the exact values depending on the chemical composition), such a temperature is already reached during the pre-main sequence (pre-MS) phase allowing lithium to be completely destroyed in fully convective structures. Since the larger the mass and the higher the rate of central temperature increase and, consequently, the earlier the onset of lithium burning, the age at which lithium is fully depleted strongly depends on mass. Thus, in stellar associations or clusters with ages between 15 and 350 Myr, one would expect to observe a sharp transition in the very low mass regime between stars with and without photospheric lithium. Such a transition, usually called lithium depletion boundary (LDB), is an age indicator (see e.g. D’Antona & Mazzitelli 1994).

The age at the LDB can be provided only by theoretical stellar evolutionary models. The accuracy of their predictions relies on the adopted input physics (e.g. equation of state, radiative opacity, nuclear reaction cross-sections, etc.), chemical composition (initial metallicity, helium and deuterium abundance, heavy element mixture, etc.), and the numerical scheme describing macroscopic physical processes (convection, rotation, etc...). All these ingredients are affected by not negligible uncertainties, which translate into a theoretical error in the LDB age estimate.

Given the primary importance of evaluating the age of young stellar clusters, several studies have been devoted to quantify the main theoretical and observational uncertainties affecting the LDB method. A first idea of the theoretical uncertainty is provided by comparing models computed by different groups (see e.g. Jeffries & Naylor 2001). However such an approach is neither able to quantify the individual uncertainty sources nor able to fully exploit the whole theoretical uncertainty, as several input physics are in common among the different evolutionary codes.

A better and more detailed way to proceed is to vary
2 THE REFERENCE SET OF MODELS

We computed the models with the most recent release of the PROSECCO (Pisa Raphson-NewtOn Stellar Evolution Computation COdEx) stellar evolution code derived from the FRANEC code (see e.g. Degl’Innocenti et al. 2008). A detailed description of the standard models can be found in Tognelli, Prada Moroni & Degl’Innocenti (2011). Tognelli, Degl’Innocenti & Prada Moroni (2012) and Dell’Omodarme et al. (2012); here we limit the discussion only to the variations with respect to the previous version.

The reference models are standard pre-MS tracks, which do not take into account accretion, rotation, and magnetic fields. The evolution starts from a fully convective and homogeneous model on the Hayashi track, with a central temperature $T_\text{c} \approx 10^5$ K, low enough for deuterium burning to be completely inefficient. Although the initial model is not physically realistic, as it does not result from a self-consistent protostar evolution, we checked that varying the initial central temperature (hence radius) in the range $[10^5, 10^6]$ K does not affect the LDB.

We adopted the 2006 release of the OPAL equation of state (OPAL06 EOS; Rogers & Nayfonov 2002). However, in order to compute objects less massive than 0.1 $M_\odot$ in the current version of the code we included also the EOS by Saumon, Chabrier & van Horn (1995, SCVH95), for temperatures and densities not covered by the OPAL06 EOS.

We used the same radiative (OPAL 2005 and Ferguson et al. 2005) and conductive (Potekhin 1999, Shtriker & Yakovlev 2000) opacities as in Tognelli, Prada Moroni & Degl’Innocenti (2011) but for a different solar heavy elements mixture, namely the Asplund et al. (2009, AS09).

Nuclear reaction rates relevant for the pre-MS evolution have been taken from the NACRE compilation (Angulo et al. 1999), with the exception of $p(e^{-}\nu)p^{2}H$ (Marcucci, Schiavilla & Viviani 2013, Tognelli et al. 2013), $^{3}\text{H}(p,\gamma)^{3}\text{He}$ (Descouvemont et al. 2004), $^{2}\text{H}(^{3}\text{He},^{3}\alpha)^{4}\text{He}$ and $^{2}\text{H}(^{4}\text{He},^{3}\alpha)^{4}\alpha$ (Tuminino et al. 2014), $^{7}\text{Li}(p,^{3}\alpha)^{8}\alpha$ (Lamia et al. 2013), and $^{7}\text{Li}(p,^{3}\alpha)^{8}\alpha$ (Lamia et al. 2012). Nuclear reaction rates between bare nuclei have been corrected to account for plasma electron screening. We implemented the weak (Salpeter 1954), weak–intermediate–strong (Dewitt, Grabske & Cooper 1973, Grabske et al. 1973), and strong (Itoh, Totsuji & Ichimaru 1977, Itoh et al. 1979) screening. Then, an adaptive procedure determines the most suitable screening factor to be used in each mass-shell of the model. For what concerns the burning of the light elements relevant for this work (mainly $^{2}\text{H}$ and $^{7}\text{Li}$), the intermediate/strong screening is generally adopted.

Outer boundary conditions (BCs) have been obtained by the detailed atmospheric models by Brott & Hauschildt (2005, BH05) as discussed in Tognelli, Prada Moroni & Degl’Innocenti (2011). Convection is treated according to the mixing length theory (Böhm-Vitense 1958) following the formalism described in Cox & Giuli (1968). Usually the solar-calibrated mixing length value ($\alpha_{\text{ML}}$) is adopted for different stellar masses and evolutionary phases. However, there are no compelling theoretical arguments for such a choice (see e.g., Ludwig, Freytag & Steffen 1999, Trampedach 2007). In particular, there are several hints that in pre-MS phase the superadiabatic convection might be less efficient (see e.g., D’Antona & Montalbán 2003, Stassun et al. 2004, Mathieu et al. 2007, Stassun 2008, Somers & Pinsonneault 2014, and references therein).

Using our models a good agreement with eclipsing binaries and surface lithium abundances in young clusters is obtained with $\alpha_{\text{ML}} \approx 1$. (Germano, Prada Moroni & Tognelli 2012 Tognelli, Degl’Innocenti & Prada Moroni 2012). We adopted $\alpha_{\text{ML}}=1.00$ in our reference models. For the sake of completeness, in the following we will show also the effect of using our solar-calibrated value $\alpha_{\text{ML}}=1.74$.

The reference set of models have been computed adopting the solar heavy element mixture by Asplund et al. (2004) and a helium-to-metal enrichment ratio $\Delta Y/\Delta Z = 2$ (Casagrande 2004). With this choices, [$Fe/H$] = +0.0 translates into an initial helium abundance $Y = 0.274$ and a global metallicity $Z = 0.013$ (see Sect. 4). The initial

\[ \text{The solar model has been computed using an iterative procedure to adjust the initial helium abundance, metallicity, and mixing length parameter in our 1 M_\odot stellar model, in order to reproduce, within a given numerical tolerance (< 10^{-4}), at the age of the Sun (4.57 Gyr) its observed radius, luminosity, and (Z/X)_\text{ph}.} \]
2.1 Lithium depletion boundary

We computed 36 pre-MS tracks in the mass range [0.06, 0.40] \( M_\odot \), in which lithium is completely destroyed in a fully convective object. We focused on this mass range because for the chosen value [Fe/H] = +0.0 objects less massive than about 0.06 \( M_\odot \) never reach the temperature required to destroy lithium, while for \( M \geq 0.4 \ M_\odot \) a radiative core develops before lithium is completely depleted.

We defined the LDB as the model at which lithium abundance in mass is reduced by a factor of 100 with respect to the original one. In order to compute a reliable lithium depletion evolution and, consequently LDB ages, the time-step adopted in model computations must be chosen with care. A time-step simply tuned to obtain an accurate pre-MS evolution would result in a too crude time resolution and in an underestimate of the lithium depletion at a given age, which in turn would lead to an overestimate of the LDB age \citep{Pla2002, Bur2004}. This is the consequence of \(^7\text{Li}\)-burning occurring over temporal-scales much shorter than the pre-MS evolutionary ones. Being aware of this, in the routine that provides the time-step for each iteration we introduced an additional condition that constraints the time-step to be short enough to lead to a relative lithium abundance variation lower than 0.1\%. Moreover, such a choice allows us to obtain also a good numerical resolution in the \(^7\text{Li}\) abundance to precisely identify the LDB point.

Fig. 1 shows our reference LDB curve in the (log \( L/L_\odot \), age) plane obtained from the reference set of models. The complete set is listed in Table 1 with the corresponding age, log \( L/L_\odot \), \( R/R_\odot \), radius, and surface gravity values at the LDB for each mass.

We also provide a fit of the log age[yr] - log \( L/L_\odot \) curve

\[ \log \text{age[yr]} = a_0 + a_1 \times \log L/L_\odot + a_2 \times (\log L/L_\odot)^2 + a_3 \times (\log L/L_\odot)^3 \]

for the reference set of models using a third-order polynomial fit. The fit parameters \( a_0, a_1, a_2, \) and \( a_3 \) are given in Table 2 the mean accuracy of the fit is better than 1% over the whole selected luminosity range.

\[ \text{Table 2. Coefficients of the third-order polynomial fit of the LDB age versus luminosity.} \]

\| \( a_0 \) \| \( a_1 \) \| \( a_2 \) \| \( a_3 \) \\
\hline
6.6408 & -0.46379 & -0.031392 & -0.014906
\hline

\footnote{Table 1 along with LDB ages for several [Fe/H] values is available in electronic form at the url: \url{http://astro.df.unipi.it/stellar-models/ldb}}

\footnote{‘log’ stands for the base-10 logarithm.}
3 INDIVIDUAL INPUT PHYSICS UNCERTAINTIES

The results of stellar model computations depend on the adopted input physics, such as nuclear reaction cross-sections, plasma electron screening, radiative opacity, equation of state (EOS), and BCs. All these ingredients are still affected by a not negligible uncertainty which directly translates into an uncertainty in stellar model outcomes [Valle et al. 2013]. In this section, we focus only on the input physics that might affect the lithium burning in pre-MS of very low mass stars.

As a first estimate of the physical uncertainty impact on the LDB age, we varied a single input physics at a time keeping all the others fixed (see Burke, Pinsonneault & Sills 2004, for a similar investigation). More in detail, when the error on a given input physics was available, we computed two additional sets of perturbed models by adopting respectively the highest and lowest value of the input physics given by its uncertainty. Then, the LDB ages provided by the perturbed models have been compared with the reference one. Unfortunately, not for all the analysed input physics an uncertainty evaluation is available; this is for example the case of the EOS and outer BCs. Each of these cases has been conveniently treated, as discussed in the following sub-sections, by substituting the adopted tables. Table 3 lists the analyzed input physics with the related assumed uncertainty/range of variation, when present, or the alternative input physics. Note that, only in the case of the nuclear reaction cross-sections the listed errors are 1σ, while in the other cases they represent the extreme values of the variability region.

Where not explicitly stated, all the models have been computed for the reference chemical composition and mixing length parameter, as described in Sect. 2.

Some of the cases analysed in the following sub-sections have been already discussed in Burke, Pinsonneault & Sills (2004). However, a detailed comparison is difficult because they adopted input physics different from ours and they recalibrated the mixing length parameter and initial helium abundance on the Sun for each perturbed set of models. A variation of the initial Y and/or αML in the perturbed models due to the solar re-calibration might partially counterbalance or increase the effect induced by the sole variation of the analysed quantity. For this reason, we preferred to show the contribution on the LDB age of the sole perturbed quantity with all the other parameters fixed. Moreover, the use of a solar calibrated αML does not guarantee a better agreement with radii of low pre-MS stars, where the lithium is actually depleted. Indeed, several papers have shown that in these stars the superadiabatic convection is much less efficient than in the Sun and that a proper value of αML is of the order of 1 (see e.g. Ventura et al. 1998, D’Antona & Montalbán 2003, Landin et al. 2004, Gennaro, Prada Moroni & Tognelli 2012, Tognelli, Degl’Innocenti & Prada Moroni 2012, and references therein).

3.1 Nuclear cross-sections

Since we are interested in LDB, the only nuclear reactions that might have an effect are those that take place before or during the 7Li-burning. Such reactions are: 2H(p,γ)3He, 3H(p,p)4H, 4H(p,p)5He, and 7Li(p,α)α.

We adopted the following uncertainties in the quoted reactions: ±3% for the p+3H reaction [Descouvemont et al. 2004, for temperatures of about 10^6 K, typical of d-burning], ±5% for the 4H+3H channels [Tumino et al. 2014], and ±10% for the 7Li+p reaction [Lamia et al. 2012].

Notwithstanding the crucial role played by deuterium burning during the early pre-MS evolution, due to the slight variation of the reaction rates within the current uncertainties, the effect on the LDB age is completely negligible. For this reason we do not show the related plots. Fig. 2 shows the relative age difference, i.e. (perturbed model – reference model)/reference model versus luminosity, due to the uncertainty in the 7Li(p,α)α cross-section. The relative age difference, about ±1%, is independent of the mass over the whole selected mass range. An increase of the cross section results
in a higher $^7\text{Li}$-burning efficiency at a given temperature, which leads to a more rapid $^7\text{Li}$ depletion and to a lower LDB age at a fixed luminosity. This effect, analysed here for the first time, is small but systematic. We computed also reference and perturbed models adopting the solar-calibrated mixing length parameter (i.e. $\alpha_{\text{ML}}=1.74$), verifying that the relative LDB age differences are completely unaffected by a variation of $\alpha_{\text{ML}}$.

### 3.2 Plasma electron screening

An important point to discuss is the effect of the electron screening on $^7\text{Li}$-burning. Plasma electrons around the interacting nuclei reduce the effective Coulombian repulsion enhancing the reaction rate by a factor $f_{\text{el}}$ (see e.g. Salpeter 1954, Dewitt, Graboske & Cooper 1973, Graboske et al. 1973). A similar effect, due to atomic electrons, is present in the measurements performed in the laboratory. To this regard, there are hints that atomic electron screening measured in laboratory is systematically lower ($\sim 1/2$) than the theoretical expectations (see e.g. Pizzzone et al. 2010, and references therein).

It is not yet clear whether theoretical computations of plasma electron screening are affected by a similar problem (see e.g. Castellani et al. 1996). Moreover, at the moment a reliable estimate of such an uncertainty source is lacking. However, the effect of modifying the plasma screening directly affects the reaction rate. For what concerns the LDB age, the uncertainty in plasma screening introduces a systematic effect on the central temperature value at which a reaction becomes efficient, thus reflecting eventually on the LDB age. Given such a situation, it might be useful to check the effect of a variation of $f_{\text{el}}$ in the $^7\text{Li}(p,\alpha)\alpha$ reaction. Given the lack of a solid estimate of the uncertainty in $f_{\text{el}}(\text{Li})$, the best thing we can do is assuming a maximum discrepancy between the predicted and real electron plasma screening efficiency similar to that observed in the laboratory; thus, we computed two sets of perturbed models with $f_{\text{el}}(\text{Li})$ increased by a factor 1.5 and 2.0.

Fig. 3 shows the relative age difference versus luminosity due to the variation of the $^7\text{Li}(p,\alpha)\alpha$ plasma electron screening. Increasing the electron screening leads to higher reaction rates and hence to lower LDB ages at a given luminosity. An enhancement by a factor of 1.5 and 2.0 produces, respectively, a relative age decrease of the order of 3% and 5.5% at a fixed luminosity. These results do not change if $\alpha_{\text{ML}}=1.74$ is adopted in place of $\alpha_{\text{ML}}=1.00$ to compute both the reference and perturbed sets of models.

The plasma electron screening variation, although never analysed before, has a significant effect on the LDB age. However, there are not yet robust estimates of the uncertainty in the $^7\text{Li}(p,\alpha)\alpha$ plasma screening factor, so present computations are intended to only give an idea of this effect on the LDB. For this reason, we decided to not include this quantity in the cumulative uncertainty computed in Sect. 5 with the warning that a possible uncertainty source has been neglected.

### 3.3 Outer BCs

To solve the differential equations which describe stellar interiors, suitable outer BCs are required. The common approach consists in specifying pressure $P(\tau_{\text{ph}})$ and temperature $T(\tau_{\text{ph}})$ provided by a detailed atmospheric model at a given optical depth $\tau_{\text{ph}}$ (the matching point between atmosphere and interior). To estimate the effect of BCs on LDB age, two aspects must be considered: (a) the adopted atmospheric model and (b) the choice of $\tau_{\text{ph}}$.

The adoption of a detailed non-grey atmospheric structure becomes important when the star has a thick convective envelope and the temperature profile is sensitive to the heat transport in the thin atmospheric layers (see e.g. Auman 1969, Dorman, Nelson & Chat 1989, Allard et al. 1997, Baraffe et al. 1998, Montalbán et al. 2004).

Unfortunately, detailed atmospheric model tables do not contain the related uncertainties. Given such a situation, one possibility to address the impact of the adopted BCs on LDB age is to compute models with BCs obtained from different atmospheric calculations. In the following we analysed the effect of the adoption of two detailed non-grey atmospheric tables, namely the Brott & Hauschildt (2000, BH05), our reference, and the Allard, Homeier & Freytag (2011, AHF11), computed by means of the same hydrostatic atmospheric code (PHOENIX, Hauschildt & Baron 1999). In both cases we used $\tau_{\text{ph}}=10$. We computed also models with the grey Krishna Swamy (1963, K366) $T-\tau$ profile, and $\tau_{\text{ph}}=2/3$. We emphasize that the last set should be considered as an extreme case, since it is well known that the adoption of grey atmospheric model is only a rough approximation of the atmospheric structure of extremely cold low-mass stars (see e.g. Baraffe et al. 1995, and references therein).

A further complication is given by the fact that the AHF11 tables for low-mass stars are available for only one metallicity ($Z=0.0134$). We decided to perform the comparison at a fixed $Z$, so even the BH05 (our reference) and the K366 models have been computed for such a metallicity, which, only in this case, is different from the reference one.

Fig. 4 shows the relative age difference versus luminosity due to the adoption of the quoted atmospheric models, for two $\alpha_{\text{ML}}$ values. As expected given the similarity of the two atmospheric models, the effect of using AHF11 in place of BH05 is relatively small, being of the order of 1% over a large interval of luminosity and reaching a maximum value...
of about 3% only at very low luminosities. Note also that, in this case, the relative LDB age differences are essentially the same using $\alpha_{\text{ML}}=1.00$ (upper panel) or the solar-calibrated value $\alpha_{\text{ML}}=1.74$ (bottom panel). A much stronger effect is due to the adoption of the grey KS66, which leads to relative age differences as large as 8-15% for $\alpha_{\text{ML}}=1.00$. These differences slightly reduce in the case of $\alpha_{\text{ML}}=1.74$ towards the upper bound of the luminosity for those masses where superadiabatic convection occurs in a progressively larger and larger region.

Besides the chosen atmospheric model, Montalbán et al. (2004) showed that the choice of $\tau_{ph}$ is important when atmospheric and interior computations do not use the same input physics/prescriptions. In this sense, $\tau_{ph}$ is another free-parameter that usually lies in the range $[2/3, 100]$ (see e.g. Table 2 in Tognelli, Prada Moroni & Degl’Innocenti 2011).

Fig. 5 shows the relative age difference versus luminosity obtained varying $\tau_{ph}$ and keeping fixed the atmospheric model, i.e. BH05, for two mixing length parameter values. Models are computed for $\tau_{ph}=2/3$, 10 (reference), and 100. Passing from $\tau_{ph}=10$ to 100 at the LDB decreases whereas from $\tau_{ph}=10$ to 2/3 it increases; the largest difference occurs when $\tau_{ph}=2/3$ is adopted. The extent of the relative age difference depends on the used mixing length parameter, the effect for $\alpha_{\text{ML}}=1.00$ being roughly twice that for $\alpha_{\text{ML}}=1.74$. Note that all the atmospheric models adopt the same mixing length value (i.e. $\alpha_{\text{ML, atm}}=2.00$, for $\tau < \tau_{ph}$). We are forced to keep it fixed because we do not have the code used to compute the atmospheric structure.

In convective regions, the temperature gradient becomes progressively more sensitive to the mixing length parameter as superadiabaticity gets larger. The models with $\tau_{ph}=2/3$ have a superadiabatic zone in the interior ($\tau \geq \tau_{ph}$), where we can actually change $\alpha_{\text{ML}}$, larger than the model with $\tau_{ph}=100$. The latter set of models has consequently a reduced sensitivity to the adopted mixing length parameter in the interior.

Given the difficulty in estimating the uncertainty in the adopted BCs, for the computation of the cumulative error stripe in LDB age in Sect. 5 we took into account only the effect of changing $\tau_{ph}$.

### 3.4 Radiative opacity

Although radiative opacity is one of the main ingredients in stellar computations, the current generation of tables do not contain any indication about the uncertainty. As a first step, to give an estimate of the opacity uncertainty propagation on the LDB age, we analysed the effect of an uncertainty in the Rosseland coefficients ($\kappa_{\text{rad}}$) of $\pm 5\%$ in the whole structure, as done in Valle et al. (2013a).

In convective regions, the actual temperature gradient depends on the radiative opacity only in superadiabatic
zones, while elsewhere it is mainly determined by the EOS. In the inner region of a fully convective star, given the high density, the temperature gradient is essentially the adiabatic one, while it becomes progressively more and more superadiabatic moving towards the surface, where the density drops. It is thus clear that in the interior of stars, the change of the opacity coefficients has a negligible effect on the structure, whereas the effect increases towards the surface.

For models computed with detailed non-grey BCs, we cannot modify the atmospheric opacity because we do not have the atmospheric code, thus preventing the possibility to consistently check the effect of $\tau_{\text{rad}}$. Moreover, for the selected mass range in the interior ($\tau \gg \tau_{\text{ph}}$), where we can actually modify the opacity coefficients, the density is high enough to guarantee an almost adiabatic convection. For this set of non-grey models we verified that the opacity between models computed with the reference and the ±5% perturbed radiative opacity. Grey outer BCs are adopted instead of the reference ones.

A way in which we can estimate the effect of an opacity variation extended also to the atmosphere is to compute stellar models adopting grey outer BCs. In this case we can actually calculate the atmosphere structure consistently with the $\tau_{\text{rad}}$ variation used in the interior. Although we are aware that the adoption of a grey BC is not the best choice for the mass range we are dealing with, such an approach allowed us to give an estimate of the effect on LDB age of an opacity variation in the whole structure, atmosphere included.

Fig. 6 shows the relative age difference versus luminosity caused by a variation of ±5% of the radiative opacity coefficients. Due to the progressively more extended superadiabatic zone in the outermost layers at increasing stellar mass, the relative age uncertainty increases with the luminosity from negligible values at the faint end to about 1% at the bright end. At a given luminosity, the lower the opacity and the larger the central temperature and, consequently, the lower the LDB age.

A further uncertainty source related to radiative opacity is worth to be discussed. The radiative opacity tables do not cover the entire temperature–density plane required for computing pre-MS evolution for masses in the range $0.06, 0.4] \, M_\odot$, and an extrapolation to higher densities is needed. As in Tognelli, Prada Moroni & Degl’Innocenti (2011), in order to evaluate the impact of the extrapolation on the LDB age, we analysed three different extrapolation techniques: constant, linear from the last two points, and a linear fit from the last four points (our reference). As in the previous cases, we cannot properly evaluate the effect of changing the extrapolation on the LDB age for non-grey models, because we cannot modify the atmospheric models. The only thing that we can analyse for non-grey models is the effect of changing the extrapolation in the interiors. As expected the effect of such a change is negligible.

However, to have an idea of the extrapolation impact we computed three additional sets of models with grey BCs where we can actually change the extrapolated opacities also in the atmosphere. Fig. 7 shows the relative LDB age difference between models computed with the reference, i.e. four point linear, and the constant and two points linear radiative opacity extrapolations. The last two extrapolations lead to underestimate the LDB age with respect to the reference one; however, such an effect vanishes at high luminosities. The constant extrapolation produces an effect which is about three times larger than the linear one.

A final uncertainty source related to radiative opacity is the heavy elements distribution adopted to compute Rosseland coefficients. To analyse the effect on the LDB age of changing the heavy element mixture at a fixed total metallicity $Z$, we computed models adopting the Grevesse & Sauval (1998, GS98) solar mixture. As in the previous cases, we discuss only the effect on grey models since non-grey ones are completely unaffected also by the mixture change in the opacity tables used to compute the interiors, the mixture adopted by BH05 atmospheric models being fixed. Fig. 8 shows the relative LDB age difference between models adopting the reference (AS09) and the GS98 solar mixture and computed with the grey BCs. Depending on the mass, the stellar structure crosses a region in the density–temperature plane where the opacity coefficients are particularly sensitive to different elements (see e.g. Sestito et al. 2006). Thus, the opacity variation due to the adopted mixture is a complex function of temperature.

Cumulative uncertainties in LDB age

![Figure 6](image1.png)

**Figure 6.** Relative age difference at LDB as a function of luminosity between models computed with the reference and the ±5% perturbed radiative opacity. Grey outer BCs are adopted instead of the reference ones.

![Figure 7](image2.png)

**Figure 7.** Relative age difference at LDB as a function of luminosity between the reference models and those computed with radiative opacity tables extrapolated with different procedures. Grey outer boundary conditions are adopted instead of the reference ones.
and density. However, as clearly visible in Fig. 8 the effect induced by the adopted mixture is very small, being always in the range $[-0.5\%, +0.5\%]$.

We computed sets of models with the reference and perturbed radiative opacity coefficients ($\kappa_{\text{rad}} \pm 5\%$, different high-density extrapolations, different heavy elements mixtures) also for $\alpha_{\text{ML}}=1.74$. The effect of an opacity variation on LDB age is completely insensitive to a mixing length change.

Summarizing the results of this section, we showed that changing the radiative opacity only in the interiors does not affect the LDB age. On the contrary, a sizeable effect is expected when the radiative opacity variation is extended to the atmosphere. However, since we were able to compute such an effect only in the case of grey BCS, we preferred to neglect the opacity contribution in the computation of the cumulative uncertainty in the LDB age (see Sect. 3), being aware that a possible uncertainty source is missing in the cumulative error estimate.

### 3.5 Equation of state

The structure of fully convective pre-MS stars depends on the EOS (see e.g., Mazzitelli 1989, D’Antona 1993, Tognelli, Prada Moroni & Degl’Innocenti 2011).

As for radiative opacity, the most commonly adopted EOS tables do not contain the uncertainty associated with the thermodynamic quantities (i.e. density, specific heat, adiabatic gradient etc.). Moreover, these quantities are strictly correlated among each other, thus preventing to follow a similar approach to estimate the uncertainty propagation into LDB age as that used for opacity. For this reason, to roughly estimate the dependence of the LDB age on the EOS, in addition to our reference set of models (OPAL06+SCVH95 EOS), we computed three additional sets adopting different EOS tables. We used the sole OPAL06 EOS, the FreeEOS (Irwin 2008, in the EOS1 configuration), and the sole SCVH95 EOS. Since both the FreeEOS and OPAL06 EOS do not cover the whole temperature–density plane suitable for computing the entire mass range $[0.06, 0.4] M_\odot$ studied in this paper, we limit the comparison to the common mass interval.

Fig. 9 shows the relative age difference versus luminosity due to the adoption of the quoted EOS. Both FreeEOS and OPAL06 induce age differences less than $\pm 1\%$. A larger effect $1-3\%$ is caused by the SCVH95 EOS.

Burke, Pinsonneault & Siils (2001) already studied the effect on LDB of the EOS, but they used tables different from those adopted here, thus preventing a detailed comparison. However, their results are, at least qualitatively, in agreement with ours. We also verified that the adoption of $\alpha_{\text{ML}}=1.74$ in the reference and perturbed EOS models leads to the same relative LDB age differences.

Due to the difficulty to properly estimate the uncertainty propagation of the adopted EOS, we did not take such an effect into account for the computation of the cumulative error bars in LDB age in Sect. 3.

### 3.6 Mixing length efficiency

Although commonly adopted, the mixing length theory is not a fully consistent treatment of the superadiabatic convection and, consequently, it introduces a relevant uncertainty source in computation of stellar structures with extended convective envelopes. The commonly adopted solar calibration of the mixing length parameter $\alpha_{\text{ML}}$ does not in principle guarantee a correct convective efficiency for stars, as in our case, of mass and/or evolutionary phase different from that of the Sun. In order to check the effect of changing the mixing length value on the LDB age, we computed two sets of models with $\alpha_{\text{ML}}=1.74$ (our solar calibrated) and the other with $\alpha_{\text{ML}}=1.00$, as suggested by some authors for pre-MS stars (see Stassun 2008, Gennaro, Prada Moroni & Tognelli 2012, Tognelli, Degl’Innocenti & Prada Moroni 2012, Somers & Pinsonneault 2014).

Fig. 10 shows the relative age difference versus luminosity due to the adoption of the quoted mixing length parameters. The effect of changing $\alpha_{\text{ML}}$ from 1.00 to 1.74 is to increase the central temperature, hence to increase the efficiency of $^7\text{Li}$-burning, and to reduce the age at the LDB. The relative age difference of the LDB is of the order of about $-1\%$ and $-6\%$ for the lowest and, respectively, highest luminosities. The effect is lower at very low luminosity (very low mass stars) as a result of the thinner superadiabatic
region. Such a result is in agreement with that shown by Burke, Pinsonneault & Sills (2004).

4 INDIVIDUAL CHEMICAL COMPOSITION UNCERTAINIES

Besides the input physics, pre-MS stellar models depend also on the adopted chemical composition, namely the initial metallicity $Z$, helium $Y$ and deuterium $X_d$ abundances. Thus, when comparing observations and theoretical models, the proper chemical composition should be adopted. However, quite often only independent measurements of the current photospheric $[\text{Fe}/\text{H}]$ of the stellar cluster/group are available, whereas direct estimates of $Y$, $X_d$, and metal abundances are missing. To overcome such a lack various assumptions are needed.

The $[\text{Fe}/\text{H}]$ value is converted into the total metallicity $Z$ adopting, for Population I stars, a solar-scaled distribution of metals, and assuming an initial helium abundance. Stellar modellers usually compute initial $Y$ using the following linear relation:

$$Y = Y_P + \frac{\Delta Y}{\Delta Z}Z,$$

where $Y_P$ is the primordial helium abundance and $\Delta Y/\Delta Z$ is the helium-to-metal enrichment ratio. The proper initial metallicity $Z$ to be adopted in stellar models follows from:

$$Z = \frac{(1 - Y_P)(Z/X)_\odot}{10^{-[\text{Fe}/\text{H}]} + (1 + \Delta Y/\Delta Z)(Z/X)_\odot},$$

where $(Z/X)_\odot$ is the current solar photospheric metal-to-hydrogen ratio. For the reference set of models we adopted $Y_P = 0.2485 \pm 0.0008$ (Cyburt 2004), $\Delta Y/\Delta Z = 2$ (Casagrande 2007), and $(Z/X)_\odot = 0.0181$ (Asplund et al. 2009).

All the quantities in equations (1) and (2) are known with an error, shown in Table 4 which directly propagates into the final $Y$ and $Z$ values. Note that these errors are intended to represent the edges of the variability region, rather than confidence intervals. Similarly to what was done in the previous sections, as a first step we computed perturbed models by varying a single parameter ($\Delta Y/\Delta Z$, $[\text{Fe}/\text{H}]$, and $(Z/X)_\odot$) at a time keeping all the others fixed to the reference value. In the following, we did not take into account the uncertainty in $Y_P$, being negligible.

To the best of our knowledge, this is the first detailed analysis of the chemical composition uncertainty in the LDB age estimates.

4.1 Initial helium abundance

Following equation (1), the initial helium abundance adopted in stellar models strongly depends on the helium-to-metal enrichment ratio, which is poorly constrained by observations (see e.g. Gennaro, Prada Moroni & Degl’Innocenti 2010). We adopted an error on $\Delta Y/\Delta Z$ of $\pm 1$ (Casagrande 2007), and we kept fixed the value of $Y_P$. In order to quantify the effect of the quoted uncertainty on the LDB age, we computed two sets of models with the reference $Z = 0.0130$ and $X_d = 2 \times 10^{-5}$, and two different values of $Y$, namely 0.261 and 0.287, which correspond to $\Delta Y/\Delta Z = 1$ and 3.

Fig. 11 shows the relative age difference versus luminosity due to the adoption of the quoted initial helium abundances. The relative age uncertainty is roughly $\pm 2\%$ over the whole mass range. The larger the helium abundance, the higher the central temperature, and the earlier the lithium depletion at a given luminosity. The results do not change if $\alpha_{ML}=1.74$ is used instead of $\alpha_{ML}=1.00$.
Relative age difference at $10$

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sets with the labelled initial metallicity $Z$

4.2 Initial metallicity

In order to analyse the effect of the metallicity on the LDB age, we computed two sets of models with the reference $Y = 0.274$ and $X_d = 2 \times 10^{-5}$, and two different metallicity values, $Z = 0.0105$ and $Z = 0.0155$, which roughly correspond to the typical observational error in $[\text{Fe}/\text{H}]$, namely $\pm 0.1$ dex.

Fig. 12 shows the relative age difference versus luminosity due to the adoption of the quoted initial metallicities. The effect is almost negligible at faint end of our range and it increases up to about $\pm 2\%$ at higher luminosities. Such a behaviour is qualitatively similar to that shown in Fig. 6 as a variation in metallicity mainly translates into a variation in the radiative opacity. This is essentially the consequence of the metallicity dependence of the outer BCs. We verified that varying the metallicity only in the interiors, keeping fixed the BCs, produces a very small ($\lesssim 0.5\%$) and constant effect on the LDB. In this case, a fixed luminosity, a larger $Z$ leads to a larger mass (higher central temperatures) and consequently to a lower $X_d$ age. Moreover, we verified that decreasing the luminosity, the pressure and temperature at the base of the atmosphere of models at the LDB get progressively less and less sensitive to the metallicity. At higher luminosity, the effect of $Z$ on the BCs becomes dominant and an increase of $Z$ results in a larger $LDB$ age. The comparison between reference and perturbed models computed with $\alpha_{\text{ML}} = 1.74$ gives the same results.

4.3 Heavy elements mixture

The value of $(Z/X)_\odot$ in equation 2 depends on the adopted solar heavy elements mixture. In our reference set of models we used the Asplund et al. (2009) one, which leads to $(Z/X)_\odot = 0.0181$. In order to quantify the impact of varying this quantity on LDB age, we computed two sets of models with the reference values of $\Delta Y/\Delta Z = 2$, $[\text{Fe}/\text{H}] = +0.0$, and $X_d = 2 \times 10^{-5}$, adopting an uncertainty of $\pm 15\%$ in $(Z/X)_\odot$ (Bahcall, Serenelli & Pinsonneault 2004; Bahcall & Serenelli 2003). This is also roughly representative of the range of values spanned by different, but still largely used, solar mixtures ($(Z/X)_\odot \approx 0.0165 - 0.0244$; Grevesse & Noels 1993, Grevesse & Sauval 1998, Asplund, Grevesse & Sauval 2005, Caffau et al. 2008).

Fig. 13 shows the relative age difference versus luminosity due to the adoption of the quoted initial metallicities and helium abundances. The shown behaviour is the consequence of a variation of both $Z$ and $Y$ resulting from a change of $(Z/X)_\odot$ at a fixed $[\text{Fe}/\text{H}]$ and $\Delta Y/\Delta Z$. Note that, as shown in Figs. 11 and 12, increasing the luminosity, the effect of a variation on $Z$ is opposite and partially counterbalanced by that of varying $Y$, leading to a total effect lower than $\pm 1\%$. The relative LDB age difference between reference and perturbed models is not affected by changing $\alpha_{\text{ML}}$ from 1.00 to 1.74.

4.4 Initial deuterium abundance

Although its tiny initial abundance, deuterium plays a relevant role in pre-MS evolution because the energy released during $d$-burning temporarily slows down the gravitational contraction. In order to evaluate the effect of the initial deuterium abundance uncertainty on the LDB age, we computed two sets of models with $X_d = 1 \times 10^{-5}$ and $3 \times 10^{-5}$ representative of the current $X_d$ range in the Galactic disc/local bubble/solar neighbourhood (see e.g., Sembach 2010).

Fig. 14 shows the relative age difference versus luminosity due to the adoption of the quoted initial deuterium abundances. The effect is always less than $\pm 1\%$. The larger is the $X_d$, the longer is the $d$-burning phase, and thus the later is the lithium depletion at a given luminosity. The relative LDB age difference due to the initial deuterium abundance variation does not change if $\alpha_{\text{ML}} = 1.74$ is used.

5 CUMULATIVE PHYSICAL AND CHEMICAL UNCERTAINTIES

In the previous section, we analysed the effect on LDB age of varying a single input physics at a time by keeping fixed all the others, and varying the chemical composition for the reference input physics. However, although commonly adopted (Bildsten et al. 1997, Burke, Pinsonneault & Sills 2004, 1999, 2003, 2004).
such an approach does not take into account possible interactions among the different ingredients of stellar models. In other words, it assumes that non-linear effects due to simultaneous variation of such quantities are absent or negligible.

For this reason, we decided to follow a more suitable procedure similar to that described in Valle et al. (2013a), consisting in the systematic and simultaneous variation of the main input on a fixed grid. More in detail, stellar models and hence the resulting LDB age depend on a vector of parameters \( \mathbf{p}_j \) (i.e. input physics and chemical elements abundances, as discussed in the previous sections). Each parameter \( p_j \) can assume three values, namely \( p_{j,0} - \delta p_j \), \( p_{j,0} \), and \( p_{j,0} + \delta p_j \), where \( p_{j,0} \) represents the reference value and \( \delta p_j \) the adopted uncertainty. In order to cover the whole parameter space of the simultaneously varied input, we computed a set of perturbed stellar models for each possible combination of the vector \( \mathbf{p}_j \). Such a distribution-free technique relies only on the specification of a sensible range of variation for each input and it does not require an explicit specification of the parent distributions of the varied parameters. Thus, it allows us to determine the edges of the variability region rather than a confidence interval. Such a technique is particularly useful when no information about the errors distribution on the analysed quantity is available.

This method is more robust than assuming an a priori independence of the input physics/chemical parameters impact, but it has the serious disadvantage of being much more computationally time consuming than the classic approach. So, in order to avoid a waste of time in useless computations, we did not vary all the input physics studied in Sect. 3. We did not perturb the d-burning cross-sections (\( \mathrm{p+}^4\mathrm{H} \), \( ^7\mathrm{H}+^4\mathrm{H} \)), since we showed that they have a completely negligible effect on the LDB age. Moreover, as discussed in the previous section, we did not vary the EOS, radiative opacity and atmospheric models since we can not properly take into account their effects. The list of the input physics allowed to vary is shown in the last column of Tables 3 and 4. Note that the perturbed stellar models can not be ruled out by the current observational constraints. In fact, we checked that the relative radius variation due to the perturbed input physics at fixed mass in the very low mass regime is smaller than 1%.

Regarding the chemical composition, the number of \((Y, Z)\) pairs resulting from all the possible combinations of \([\mathrm{Fe/H}]\pm0.1\) dex, \(\Delta Y/\Delta Z \pm 1\), and \((Z/X)_\odot \pm 15\%\) values is 27 (= 3\(^3\)). However, it is not needed to calculate set of models for each \((Y, Z)\) pair. Indeed, being interested only in the edges of the variability region, it is enough to compute stellar models for the 7 \((Y, Z)\) pairs listed in Table 5 which correspond to the reference set and to the sets with the extreme combinations of helium abundance and/or metallicity.

In summary, we have adopted for the cumulative error evaluation seven pairs of \((Y, Z)\), three initial deuterium abundances \(X_d\), two mixing length values \(\alpha_{\text{ML}}\), three \(\tau_{\text{ph}}\) and three \(\Delta \text{Li}(p,\alpha)\) cross-section values, for a total of 378 (= 7 \(\times\) 3 \(\times\) 2 \(\times\) 3 \(\times\) 3) sets of models. For each set we computed a grid of 31 stellar masses, for a total of 11718 stellar tracks.

From the perturbed stellar models, we then obtained the cumulative error stripe, following the algorithm of

**Table 5. Pairs of \((Y, Z)\) values adopted for the computations of the models with the perturbed chemical composition used for constructing the cumulative error stripe (see text).**

| \((Y, Z)\) | \([\mathrm{Fe/H}]\) | \(\Delta Y/\Delta Z\) | \((Z/X)_\odot\) |
|-----------|-----------------|----------------|----------------|
| \(Y = 0.2740\) | \(Z = 0.0130\) | +0.0 | 2 | \text{AS09} |
| \(Y = 0.2790\) | \(Z = 0.0100\) | −0.1 | 3 | \text{AS09} |
| \(Y = 0.2650\) | \(Z = 0.0160\) | +0.1 | 1 | \text{AS09} |
| \(Y = 0.2750\) | \(Z = 0.0090\) | −0.1 | 3 | \text{AS09−15\%} |
| \(Y = 0.2670\) | \(Z = 0.0190\) | +0.1 | 1 | \text{AS09+15\%} |
| \(Y = 0.2570\) | \(Z = 0.0090\) | −0.1 | 1 | \text{AS09−15\%} |
| \(Y = 0.3020\) | \(Z = 0.0190\) | +0.1 | 3 | \text{AS09+15\%} |

**Table 6. Pairs of \((Y, Z)\) values adopted for computing models with the individual perturbation of \([\mathrm{Fe/H}], \Delta Y/\Delta Z\), and \((Z/X)_\odot\) (see text).**

| \((Y, Z)\) | \([\mathrm{Fe/H}]\) | \(\Delta Y/\Delta Z\) | \((Z/X)_\odot\) |
|-----------|-----------------|----------------|----------------|
| \(Y = 0.2710\) | \(Z = 0.0120\) | +0.0 | 2 | \text{AS09−15\%} |
| \(Y = 0.2780\) | \(Z = 0.0150\) | +0.0 | 2 | \text{AS09+15\%} |
| \(Y = 0.2620\) | \(Z = 0.0130\) | +0.0 | 1 | \text{AS09} |
| \(Y = 0.2870\) | \(Z = 0.0130\) | +0.0 | 3 | \text{AS09} |
| \(Y = 0.2690\) | \(Z = 0.0100\) | −0.1 | 2 | \text{AS09} |
| \(Y = 0.2810\) | \(Z = 0.0160\) | +0.1 | 2 | \text{AS09} |
constructing the convex hull described in appendix A in Valle et al. (2013a). Here we briefly summarize such a procedure. Each set of perturbed stellar models provides an LDB curve in the age–luminosity plane. Thus, at a given luminosity an ensemble of perturbed LDB ages is available. The convex hull of such an ensemble provides the error in the predicted LDB age for the chosen luminosity. The error stripe is then obtained by changing the luminosity.

Our main aim is to quantify the cumulative uncertainty resulting from input physics and chemical composition, keeping fixed $\alpha_{\text{ML}}$. However, the $\alpha_{\text{ML}}$ value is itself very uncertain and it might significantly affect the estimate of the cumulative uncertainty. For this reason, Fig. 15 shows the error stripe on the relative LDB age computed by taking into account the stellar models with simultaneously perturbed input physics and chemical composition for two fixed $\alpha_{\text{ML}}$ values, namely $\alpha_{\text{ML}}= 1.00$ and 1.74.

The cumulative error stripe is asymmetric and it gets progressively broader at increasing luminosity. The case with $\alpha_{\text{ML}}=1.00$ shows the largest and most asymmetric stripe. The minimum relative LDB age uncertainty is of about $\pm 5\%$ while the maximum is of $+15\%/-8\%$. The case with $\alpha_{\text{ML}}=1.74$ (solar calibrated) shows a thinner and more symmetric error stripe, with a maximum extension of about $+10\%/-8\%$.

Both the stripe asymmetry and dependence on $\alpha_{\text{ML}}$ result essentially from the variation of $\tau_{\text{ph}}$, which assumes non-symmetric values with respect to the reference one (see Sect. 3.3). Indeed, we verified that if the contribution of $\tau_{\text{ph}}$ is removed from the cumulative uncertainty estimate, the error stripe becomes essentially symmetric and unaffected by $\alpha_{\text{ML}}$.

Fig. 15 also shows the relative LDB age uncertainty computed by linearly/quadratically adding the errors due to the individual variation of each input physics/chemical composition used to build the stripe. To do this we computed additional sets of models to account for the effect of the individual variation of each parameter used to obtain $Y$ and $Z$ (i.e. [Fe/H], $\Delta Y/\Delta Z$, and $(Z/X)_\odot$), as listed in Table 6. Such models have been obtained by keeping fixed the input physics and varying a single chemical input within its uncertainty range (i.e. [Fe/H] $\pm 0.1$, $\Delta Y/\Delta Z \pm 1$, and $(Z/X)_\odot \pm 15\%$) in equations (1) and (2). The comparison is intended to quantify the difference between the classical method (linear/quadratic individual error sum), often used in the literature, and the simultaneous variation on a fixed grid (cumulative error stripe) presented in this paper. It is evident that the adoption of a linear sum produces an uncertainty region much closer to the cumulative error stripe than that obtained by a quadratic sum. In particular the latter method gives a systematic underestimation of the uncertainty region (about two times smaller). Fig. 15 has been computed perturbing simultaneously the input physics and the chemical composition. However, to better understand the results it is worth to discuss the two effects separately.

First, we verified that the contribution of the input physics (keeping the chemical composition fixed to the reference value) sums linearly, as already obtained by Valle et al. (2013a). In other words, in the case of a constant chemical composition and small perturbations of the input physics, the uncertainty region obtained by a linear sum coincides with the edges of the cumulative error stripe, while the quadratic sum results in a significative underestimate.

Secondly, to obtain the uncertainty region edges due to the perturbed chemical composition using both the linear and the quadratic sum, we used the additional models listed in Table 6. It is important to notice that equations (1) and (2) are non-linear in the parameters [Fe/H], $\Delta Y/\Delta Z$ and $(Z/X)_\odot$. In addition, even if a symmetric perturbation of the parameters is adopted the resulting $Y$ and $Z$ variation is not necessarily symmetric. The non-linearity of eqs (1) and (2) introduces a non linear response of the LDB age variation that essentially depends on the initial helium abundance and metallicity perturbation ($\delta Y$, $\delta Z$) with respect to the reference values. In the case of the simultaneous perturbation of all the chemical parameters the initial helium abundance and metallicity variations are simply given by $\delta Y_{\text{ini}} = Y_{\text{ref}} - Y_{\text{cumulative}}$ and $\delta Z_{\text{ini}} = Z_{\text{ref}} - Z_{\text{cumulative}}$ ($Y_{\text{cumulative}}$ and $Z_{\text{cumulative}}$ are the values listed in Table 6), while in the case of the individual parameter perturbation by $\delta Y_{\text{ini}} = \sum_{\delta p_j} [Y_{\text{ref}} - Y_{\text{individual}}(\delta p_j)]$
and \( \delta Z_{\text{in}} = \sum_{p_j} [Z_{\text{ref}} - Z_{\text{individual}}(\delta p_j)] (Y_{\text{individual}}(\delta p_j) \times Z_{\text{individual}}(\delta p_j)) \) are the values listed in Table 6 and \( \delta p_j \) is the \( j \)-th parameter variation. The non-linearity of equations (1) and (2) leads to \((\delta Y_{\text{cu}}, \delta Z_{\text{cu}}) \neq (\delta Y_{\text{in}}, \delta Z_{\text{in}})\), thus to a different variation of the LDB age if the cumulative or independent linear sum models are considered. We verified that this produces the differences between the linear sum edges and the cumulative stripe.

Fig. 16 shows the cumulative error stripe obtained by varying the chemical composition parameters but keeping fixed the input physics to the reference ones, for \( \alpha_{\text{ML}} = 1.00 \) and \( \alpha_{\text{ML}} = 1.74 \). The uncertainty in the initial chemical composition produces a not negligible variation of the LDB age, which is slightly asymmetric due to the asymmetric variation of \( Y \) and \( Z \), regardless the adopted mixing length. The relative LDB age uncertainty ranges from about 3% to about 6%, increasing at larger luminosities. When compared to the total error stripe (which includes also the input physics perturbation, the contribution of the sole uncertainty on the initial chemical composition is of the order of 40%). Fig. 16 also shows the linear/quadratic sum edges. As anticipated, the uncertainty obtained with a linear sum does not coincide with the error stripe even if the differences are quite small (about 1%), while the quadratic sum drastically underestimates the uncertainty (up to 1.5–2 times smaller).

One should be aware that the global error stripe shown in Fig. 16 is probably an underestimate of the actual uncertainty in the LDB age since, as explained in the previous sections, some potential contributions have been neglected, given the difficulty to properly account for them. However, assuming a linear behaviour also for these contributions, we can give a rough estimate to the total LDB age uncertainty taking into account the additional effect of the outer BCs, the EOS, and the radiative opacity. As shown in Sect. 3, an uncertainty of 1% is due to the adoption of a different non-grey BC (Fig. 4), 1–2% to the EOS (Fig. 9), and 1% to an opacity variation (Fig. 6). Thus, adding these contributions the uncertainty gets larger, ranging from a minimum of about \( \pm 9\% \) to a maximum of \( +19\%/−12\% \), for \( \alpha_{\text{ML}} = 1.00 \), or to a maximum of \( +14\%/−12\% \), for \( \alpha_{\text{ML}} = 1.74 \). It would be worth to have an estimate of the uncertainty affecting the plasma electron screening factor, which, as shown in Sect. 3.2, might significantly affect the LDB age. Moreover, as already mentioned, we focused on standard pre-MS models that do not take into account some potentially important physical mechanism. In particular it has been recently shown that magnetic fields and star spots might significantly affect the LDB age estimate (Jackson & Jeffries 2014, Malo et al. 2014). The systematic bias due to neglecting these physical processes might be comparable or even larger than the cumulative error shown in Fig. 16.

Finally, Fig. 17 shows our reference LDB curve in the \((\log L/L_\odot, \text{age})\) plane with overplotted the cumulative error stripe, for the \( \alpha_{\text{ML}} = 1.00 \) (upper panel) and \( \alpha_{\text{ML}} = 1.74 \) (bottom panel). The cumulative error stripe is overplotted as a shaded area.
uncertainty is larger (about 50%) than that provided by Burke, Pinsonneault & Sills (2004). However, a detailed comparison of the two results is not straightforward because of both the different uncertainty sources taken into account and the adopted estimate method. As an example of the former point, we accounted for the uncertainty in the adopted chemical composition, which is lacking in Burke, Pinsonneault & Sills (2004). Concerning the latter point, they adopted as best estimate of the total uncertainty a quadratic summation rather than the cumulative stripe described here.

6 CONCLUSIONS

In this work, we performed a systematic analysis of the main theoretical uncertainties affecting the age at the LDB. With respect to other studies in the literature (e.g., Bildsten et al. 1997, Ushomirsky et al. 1998, Jeffries & Naylor 2001, Burke, Pinsonneault & Sills 2004), which analysed the effect of changing an input physics at the time by keeping fixed all the others, the main novelty of our approach consists in taking into account the simultaneous variation of the main input/parameters without an a priori assumption of independence. More in detail, we computed sets of pre-MS models for all the possible combinations of the perturbations of the ingredients allowed to vary. Such an approach is much more robust in presence of interactions between the varied quantities, but it clearly requires the computation of a huge number of models. For this work, we computed about 12,000 pre-MS evolutionary tracks.

Besides the method, we improved the analysis of the LDB age uncertainty by studying physical error sources never discussed before, such as the plasma electron screening and the nuclear cross-sections ($^7Li+p$ $^2H$, $^3H+p$). A further improvement with respect to previous studies is the detailed computation of the uncertainties propagation due to the adopted initial chemical elements abundances, i.e. the total metallicity, the helium and deuterium abundances, and heavy elements mixture.

The results of the uncertainty computations are shown in Fig. [17] where the LDB age–luminosity curve with the cumulative error stripe for both $\alpha_{ML}=1.00$ and 1.74 has been plotted. To the best of our knowledge, this is the first time that such an error stripe has been shown.

As discussed in Burke, Pinsonneault & Sills (2004), the LDB age estimate is more accurate at faint luminosity. Our detailed computations confirm that the error stripe gets progressively broader at increasing luminosities. Moreover, we also showed that the cumulative error is asymmetric and it depends on the adopted $\alpha_{ML}$ value. More in detail, the set of models with $\alpha_{ML}=1.00$ shows a larger and more asymmetric error stripe, with positive relative age errors being about twice the negative ones, ranging from $\approx 5$ to $\approx 15\%$. The uncertainty range reduces to a maximum of about 10% if the solar-calibrated mixing length value ($\alpha_{ML}=1.74$) is adopted. A further contribution of the order of 4% should be added to the global error to take into account the uncertainty in the EGS, outer BC and radiative opacity, not explicitly accounted for in the error stripe computation. Finally, an additional uncertainty source potentially sizeable and worth to be studied in more detail is the electron screening factor. We proved that increasing it of 50% (100%) leads to a 3% (5.5%) variation in the LDB age.

We showed that the cumulative error stripe computed by simultaneously perturbing the input physics is in excellent agreement with the simple linear sum of the LDB relative age differences obtained by individually perturbing the input physics themselves. Regarding the chemical composition uncertainties, the agreement is worse but always better than 1%. On the contrary, the use of a quadratic sum (as generally done in the literature) results in a systematic underestimate of the total LDB age uncertainty, which is about 1.5/2 times smaller than that obtained with the cumulative error stripe.

Another result of this work consists in quantifying, for the first time, the effect of the initial chemical elements abundances uncertainty on the LDB age. We showed that the contribution of the sole chemical composition uncertainty is not negligible ranging from $\pm 3$ to $\pm 6\%$, thus accounting for at least $\approx 40\%$ of the total error budget; this part of the error stripe is almost symmetric and independent of $\alpha_{ML}$ value.

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