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R. Razera ¹★, B. Barbuy ¹★, T. C. Moura ¹, H. Erandes ¹, A. Pérez-Villegas ², S. O. Souza ^{1,3},
C. Chiappini ³★, A. B. A. Queiroz ^{3,4}, F. Anders ⁵, J. G. Fernández-Trincado ⁶, A. C. S. Friaça ¹,
K. Cunha ^{7,8}, V. V. Smith ⁹, B. X. Santiago ¹⁰, R. P. Schiavon ¹¹, M. Valentini ³, D. Minniti ^{12,13},
M. Schultheis ¹⁴, D. Geisler ^{15,16,17}, J. Sobeck ¹⁸, V. M. Placco ⁹ and M. Zoccali ^{19,20}

¹IAG, Universidade de São Paulo, Rua do Matão 1226, Cidade Universitária, São Paulo 05508-900, Brazil

²Instituto de Astronomía, Universidad Nacional Autónoma de México, A. P. 106, C.P. 22800, Ensenada, B. C., México

³Leibniz-Institut für Astrophysik Potsdam (AIP), An der Sternwarte 16, D-14482 Potsdam, Germany

⁴Institut für Physik und Astronomie, Universität Potsdam, Haus 28 Karl-Liebknecht-Str 24/25, D-14476 Golm (Potsdam), Germany

⁵Institut de Ciències del Cosmos, Universitat de Barcelona (IEEC-UB), Martí i Franquès 1, E-08028 Barcelona, Spain

⁶Instituto de Astronomía, Universidad Católica del Norte, Av. Angamos 0610, Antofagasta, Chile

⁷University of Arizona, Tucson, AZ 85719, USA

⁸Observatório Nacional, S Cristóvão, Rio de Janeiro, Brazil

⁹NSF's NOIRLab, 950 N. Cherry Ave., Tucson, AZ 85719, USA

¹⁰Universidade Federal do Rio Grande do Sul, Caixa Postal 15051, 91501-970 Porto Alegre, Brazil

¹¹Astrophysics Research Institute, Liverpool John Moores University, Liverpool, L3 5RF, UK

¹²Instituto de Astrofísica, Facultad de Ciencias Exactas, Universidad Andres Bello, Fernández Concha 700, Las Condes, Santiago, Chile

¹³Vatican Observatory, Vatican City State I-00120, Italy

¹⁴Observatoire de la Côte d'Azur, CNRS, Laboratoire Lagrange, Université Côte d'Azur, F-06108 Nice, France

¹⁵Departamento de Astronomía, Universidad de Concepción, Casilla 160-C, Concepción, Chile

¹⁶Instituto de Investigación Multidisciplinario en Ciencia y Tecnología, Universidad de La Serena, Avenida Raúl Bitrán S/N, La Serena, Chile

¹⁷Departamento de Astronomía, Facultad de Ciencias, Universidad de La Serena, Av. Juan Cisternas 1200, La Serena, Chile

¹⁸Department of Astronomy, University of Washington, Box 351580, Seattle, WA 98195, USA

¹⁹Instituto de Astrofísica, Pontificia Universidad Católica de Chile, Casilla 306, Santiago 22, Chile

²⁰Millennium Institute of Astrophysics (MAS), Nuncio Monseñor Sótero Sanz 100, Providencia, Santiago Chile

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ABSTRACT

The central part of the Galaxy hosts a multitude of stellar populations, including the spheroidal bulge stars, stars moved to the bulge through secular evolution of the bar, inner halo, inner thick disc, inner thin disc, as well as debris from past accretion events. We identified a sample of 58 candidate stars belonging to the stellar population of the spheroidal bulge, and analyse their abundances. The present calculations of Mg, Ca, and Si lines are in agreement with the ASPCAP abundances, whereas abundances of C, N, O, and Ce are re-examined. We find normal α -element enhancements in oxygen, similar to magnesium, Si, and Ca abundances, which are typical of other bulge stars surveyed in the optical in Baade's Window. The enhancement of [O/Fe] in these stars suggests that they do not belong to accreted debris. No spread in N abundances is found, and none of the sample stars is N-rich, indicating that these stars are not second generation stars originated in globular clusters. Ce instead is enhanced in the sample stars, which points to an s-process origin such as due to enrichment from early generations of massive fast rotating stars, the so-called spinstars.

Key words: stars: abundances – stars: atmospheres – Galaxy: bulge.

1 INTRODUCTION

The stellar populations in the central part of the Galaxy can inform us about its complex formation processes. This region was recently confirmed to contain stars in a metal-poor spheroidal bulge (e.g. Babusiaux et al. 2010; Dékány et al. 2013; Babusiaux 2016; Zoccali, Valenti & Gonzalez 2018; Arentsen et al. 2020; Kunder et al. 2020;

Savino et al. 2020; Queiroz et al. 2021 and references therein), along with a metal-rich contribution from the bar and inner thin disc, thick disc, and halo interlopers. In addition, debris of past accretion events, such as *Gaia*-Enceladus-Sausage (GES) (Belokurov et al. 2018; Helmi et al. 2018), and many other dwarf galaxy remnants, and minor substructures absorbed during the early stages of the Galaxy formation (see e.g. Horta et al. 2020, 2021, 2022; Fernández-Trincado et al. 2022) are present. Therefore studies of the Galactic bulge region are important for understanding the early stages of our Galaxy's formation (e.g. Barbuy, Chiappini & Gerhard 2018a; Rojas-Arriagada et al. 2020). In particular, Queiroz et al. (2020,

* E-mail: roberta.razera@usp.br (RR); b.barbuy@iag.usp.br (BB); cristina.chiappini@aip.de (CC)

2021) combining distance derivation with proper motions from the *Gaia* Early Data Release 3 (Gaia Collaboration et al. 2021) revealed stars of large eccentricity, but with orbits confined to the bulge region – with a maximum height from the Galactic mid-plane, $|z|_{\max}$, below 3 kpc with intermediate metallicities, which are good candidates for belonging to the oldest Galactic bulge component (which we here call spheroid bulge stars).

The spheroidal metal-poor bulge can be thought of as a pressure supported structure formed through violent processes, such as hierarchical clustering via minor mergers at a very early stage of the Galaxy. Ferraro et al. (2021) finds evidence that clumps of stars and gas existed at the time of the Milky Way formation. N -body simulations assume instead that early stellar discs heat rapidly as they form, and can lead to different density distributions for metal-rich and metal-poor stars (e.g. Debattista et al. 2017). Many other options are possible to form the metal-poor spheroid such as a major merger accretion of dwarf galaxies among others (e.g. Barbuy et al. 2018a). Whatever process it leads to an observed metal-poor spheroid, and it has also to explain the very old ages of the *in situ* globular clusters such as e.g. HP 1 (Kerber et al. 2019), Djorgovski 2 (Ortolani et al. 2019), Palomar 6 (Souza et al. 2021) of ages derived to be of 12.8 ± 0.9 , 12.7 ± 0.7 , and 12.4 ± 0.9 Gyr, respectively.

The search for the earliest stars in the Galaxy is an important endeavour to try to identify the earliest chemical abundances imprinted in the oldest stars, and the nature of the supernovae that enriched them. Most of the current observational efforts in finding the chemical imprints left by the first stars have focused on the most metal-poor stars found in the Milky Way halo (Beers & Christlieb 2005; Beers et al. 2017). Very metal-poor stars were also found in ultra-faint dwarf galaxies, which are intriguing dark matter dominated objects with very low average metallicities (Ji et al. 2016). The Galactic bulge, as well as the halo, is a potential host of some of the oldest stars in our Galaxy. Tumlinson (2010) suggests that half of the oldest stars were formed in the central parts of the Galaxy. Searches for field metal-poor stars in the Galactic bulge are the target of surveys such as those by Howes et al. (2016), Casey & Schlafman (2015), the Pristine Inner Galaxy Survey (PIGS, Arentsen et al. 2020), HERBS (Duong et al. 2019a, Duong et al. 2019b), and COMBS (Lucey et al. 2019, 2021, 2022) surveys. Metal-poor stars in the Galactic bulge have been mostly traced by Globular clusters (Rossi et al. 2015; Bica, Ortolani & Barbuy 2016) and RR Lyrae stars (Minniti et al. 2017), which show a peak at $[\text{Fe}/\text{H}] \sim -1.0$ (Barbuy et al. 2018a). This metallicity peak at $[\text{Fe}/\text{H}] \sim -1.0$ has been also recently confirmed regarding field stars by Lucey et al. (2021). In fact, it is expected that a fast chemical enrichment in the Galactic bulge results in a very old population with this relatively high metallicity that would correspond to the age of stars with $[\text{Fe}/\text{H}] \sim -3.0$ in the halo (Chiappini et al. 2011; Wise et al. 2012; Barbuy et al. 2018a).

Our main interest in the present work is to analyse the abundances of stars of the spheroidal bulge with a moderate metallicity of $[\text{Fe}/\text{H}] < -0.8$, in order to try to identify the earliest supernovae of the central regions of the Galaxy, and imposing constraints on the early chemical enrichment of the Milky Way. For the selection of sample stars, we applied kinematical and dynamical criteria by combining data from Apache Point Observatory Galactic Evolution Experiment (APOGEE) and *Gaia* Early Data Release EDR3. We chose stars with azimuthal velocity $V_\phi < 0$ (this selection will avoid contamination by disc stars, but would still include accreted debris of objects such as GES) that have orbits confined within 4 kpc of the Galactic centre, a maximum height of $|z|_{\max} < 3.0$ kpc, eccentricity > 0.7 and with orbits not supporting the bar structure. With this selection, as noted above, we expect our sample to be dominated by a pressure supported,

most probably old component of the bulge. We hope to discard the contamination of our sample by accreted debris, thanks to the detailed chemical information and, in particular, the alpha-over-iron enhancement expected to be low in most of the accreted debris. Finally, given that we used a barred potential, the z -component of angular momentum (L_z) is not conserved, and most orbits are either retrograde or prograde, and a fraction among those identified as counter-rotating keep retrograde along its orbit.

In this paper, we carried out an analysis of atomic and molecular lines for the selected sample of 58 metal-poor spheroid bulge star candidates aiming at refining the APOGEE Stellar Parameter and Chemical Abundance Pipeline (ASPCAP; García-Pérez et al. 2016) results, in order to interpret the derived abundances in terms of the early chemo-dynamical evolution of the bulge. As it will be shown, this re-analysis is critical for some alpha elements, and therefore for the identification and confirmation of old spheroid bulge stars at moderately low metallicities. In the present work, we adopt the stellar parameters issued from the DR17 release of the ASPCAP code. The C, N, and O abundances are derived from CO, OH, and CN lines that are interdependent and since there are such molecular lines all over the spectra, they can affect the abundances of atomic lines. We also refine the abundances of Ce. Other elements including Na, Al, and iron-peak elements will be the topic of a future work.

In Section 2, the selection of our sample is described. The element abundances are derived in Section 3. In Section 4, the results are compared with literature data for bulge samples and chemodynamical models, and discussed. In Section 5, conclusions are drawn.

2 THE SAMPLE

APOGEE (Majewski, Schiavon & Frinchaboy 2017) is part of the Sloan Digital Sky Survey III and IV (SDSS; Blanton et al. 2017). It is a project encompassing spectroscopic programs that observe Milky Way stars at high-resolution and high signal-to-noise ratios (S/N) in the near-infrared (NIR). The project SDSS-IV technical summary, the SDSS telescope and APOGEE spectrograph are described in Blanton et al. (2017), Gunn et al. (2006), and Wilson et al. (2019), respectively, whereas Zasowski et al. (2013, 2017), Beaton et al. (2021), and Santana et al. (2021) describe the APOGEE and APOGEE-2 Target Selections. The data release 17 (DR17) contains high-resolution ($R \sim 22\,500$) NIR spectra (15140–16940 Å) for some 7×10^5 stars, covering both the northern and southern sky. While APOGEE-1 observed the Milky Way bulge/bar at $l > 0$ deg, APOGEE-2 covers the whole bulge/bar region.

Given that the central part of the Milky Way hosts members of all Galactic components, including the bulge, disc, and halo (Pérez-Villegas et al. 2020; Rojas-Arriagada et al. 2020; Queiroz et al. 2021), we have used the chemo-orbital analysis shown in Queiroz et al. (2021) to identify good candidates in the spheroidal bulge APOGEE sample. To disentangle the different stellar populations coexisting in the innermost parts of the Galaxy is not an easy task, and one of the difficulties is to compute precise distances for these stars due to the high extinction. Thanks to StarHorse (Santiago et al. 2016; Queiroz et al. 2018), precise stellar distances for the entire APOGEE sample were derived both for DR16 (Ahumada et al. 2020; Queiroz et al. 2020) and DR17¹ (Abdurro’uf et al. 2022; Queiroz et al. 2022, in preparation).

We selected stars from the reduced-proper-motion (RPM) sample of Queiroz et al. (2021). For that sample, orbits were calculated using

¹Value added catalogues are available in both releases.

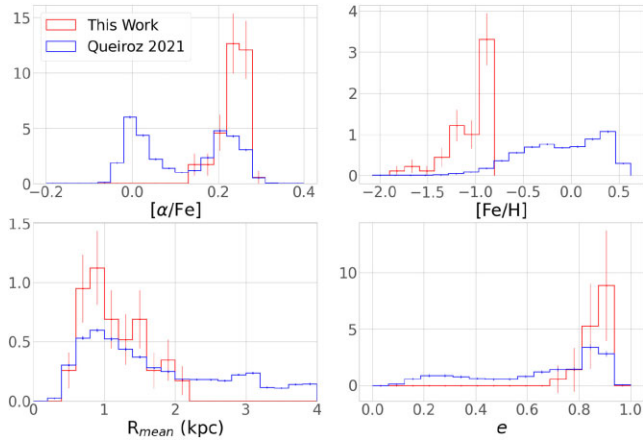


Figure 1. Comparison of the present sample of 58 selected stars (red) and the RPM sample of Queiroz et al. (2021) (blue). Upper panels: normalized distribution of metallicity and alpha-to-iron ratios from APOGEE; lower panels: mean radius $R_{\text{mean}} = (R_{\text{apocentre}} + R_{\text{pericentre}})/2$ and eccentricity of the orbits.

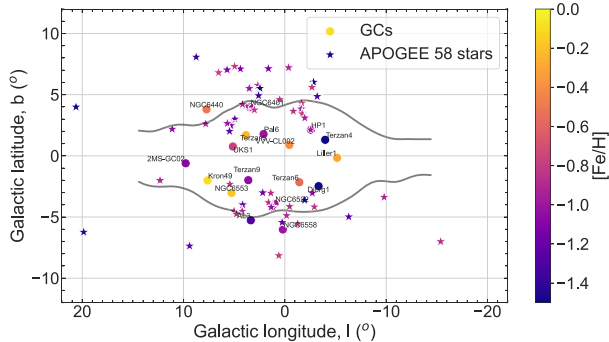


Figure 2. Projected l, b distribution of studied stars in the Galactic bulge region. Symbols: filled stars: this work; filled circles: bulge globular clusters (GCs); solid black line: contours of the bulge. The colours indicate metallicity according to the colour-bar.

bulge stars of the metal-poor spheroid, that are moderately metal-poor, α -rich and in eccentric orbits but confined to the Galactic centre region. Fig. 2 shows the projected l, b distribution of the sample in the Galactic bulge region.

As explained above, our stars were selected from the reduced-proper-motion sample of Queiroz et al. (2021), and therefore have a signal-to-noise $\text{SNR} > 50$, a good spectral fit from the ASPCAP pipeline $\text{ASPCAP_Chi2} < 25$, and a radial velocity scatter $V_{\text{scatter}} < 1.5 \text{ km s}^{-1}$. As for the renormalized unit weight error – RUWE *Gaia* EDR3 parameter, 56 out of 58 stars in our sample comply with the standard or minimal requirements to get reliable orbital elements, since astrometry from *Gaia* EDR3 has its own caveats. According to the *Gaia* consortium, the RUWE parameter is suggested to return stars astrometrically well-behaved by applying a cut with $\text{RUWE} \leq 1.4$, which is followed by the 56 stars listed in Table 1. The stars 2M17453659-2309130 and 2M18023156-2834451 have a $\text{RUWE} > 1.4$, which makes them sources with astrometric parameters that are not reliable enough.

In Fig. 3, a Kiel diagram of the sample stars is plotted with the effective temperature from ASPCAP and gravity $\log g$ coming from the StarHorse output from Queiroz et al. (2020), and compared with the reduced-proper-motion sample of Queiroz et al. (2021).

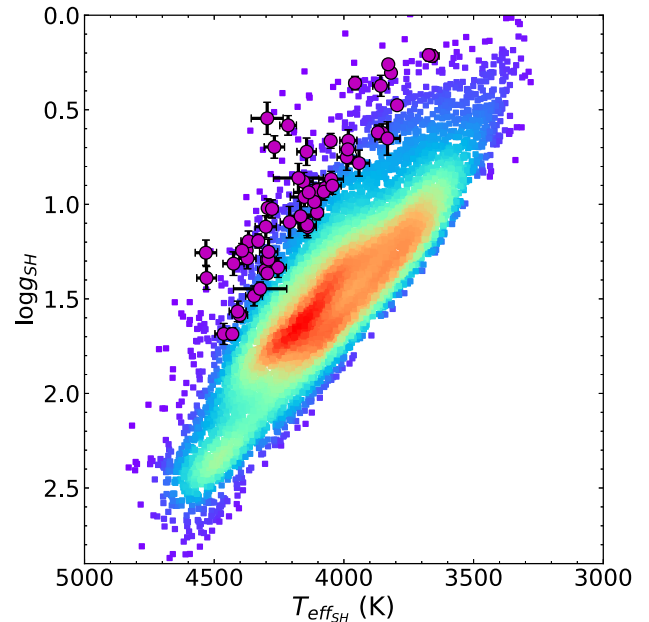


Figure 3. Kiel diagram of the 58 sample bulge stars (purple circles). In the background, we show the full reduced proper-motion sample of Queiroz et al. (2021).

3 ANALYSIS

We have initially adopted the calibrated stellar parameters effective temperature T_{eff} , gravity $\log g$, metallicity $[\text{Fe}/\text{H}]$, and microturbulence velocity v_t from APOGEE DR16 – we point out that the calibrated parameters give very different element abundances, and should not be used for such aims. In fact, the results from the ASPCAP (García-Pérez et al. 2016) are obtained for the reported non-calibrated spectroscopic stellar parameters. We then adopted these non-calibrated stellar parameters from DR17, since we became aware that these are obtained from a spectroscopic solution that minimizes the errors in seven dimensions (T_{eff} , $\log g$, $[\text{Fe}/\text{H}]$, v_t , $[\alpha/\text{Fe}]$, $[\text{C}/\text{Fe}]$, $[\text{N}/\text{Fe}]$).

For this reason, we proceeded with all the rederivation of abundances with the DR17 non-calibrated parameters. These stellar parameters are reported in Table 2, and they are the final parameters adopted.

The abundances were determined by comparing the observed spectra with the synthetic ones. The synthetic spectra calculations are carried out with the code PFANT,³ as described in Barbuy, Trevisan & de Almeida (2018b). This code is an update of the original FANTOM or ABON2 Meudon code by M. Spite. Each model atmosphere was interpolated in the MARCS grids (Gustafsson et al. 2008).

The atomic line list employed is that from the APOGEE collaboration (Smith et al. 2021). Molecular electronic transition lines of $\text{CN } A^2\Pi-X^2\Sigma$, vibration-rotation $\text{CO } X^1\Sigma^+$, $\text{OH } X^2\Pi$, and $\text{TiO } \phi$ -system $b^1\Pi-d^1\Sigma$ lines were included. The line lists for CN were made available by S. P. Davis, the CO line lists were adopted from Goorvitch (1994), and the OH are from Goldman et al. (1998). For TiO the line list is from Jorgensen (1994). More details on CN, CO, OH, and TiO molecular lines are given in Meléndez & Barbuy (1999), Meléndez, Barbuy & Spite (2001), Meléndez et al. (2003), Schiavon & Barbuy (1999) and Barbuy et al. (2018b).

³The code is available at <http://trevisanj.github.io/PFANT>.

Table 2. Selected 58 stars and corresponding DR17 non-calibrated stellar parameters.

| ID | $T_{\text{eff(nc)}}$ (K) | $\log g_{\text{(nc)}}$ | $[\text{Fe}/\text{H}]_{\text{(nc)}}$ | v_t (kms^{-1}) |
|--------------------|-----------------------------|------------------------|--------------------------------------|--------------------------------|
| 2M17153858-2759467 | 3922.7 | 0.34 | -1.62 | 2.62 |
| 2M17173693-2806495 | 3908.9 | 0.95 | -0.97 | 2.20 |
| 2M17250290-2800385 | 3796.6 | 0.91 | -0.80 | 2.39 |
| 2M17265563-2813558 | 4096.2 | 1.0 | -1.31 | 1.89 |
| 2M17281191-2831393 | 4029.1 | 0.95 | -1.17 | 1.73 |
| 2M17295481-2051262 | 4205.9 | 1.50 | -0.85 | 1.71 |
| 2M17303581-2354453 | 3863.0 | 0.77 | -0.99 | 2.13 |
| 2M17324257-2301417 | 3668.2 | 0.79 | -0.82 | 2.30 |
| 2M17330695-2302130 | 3566.6 | 0.35 | -0.93 | 2.42 |
| 2M17344841-4540171 | 3869.2 | 0.85 | -0.88 | 2.16 |
| 2M17351981-1948329 | 3553.5 | 0.44 | -1.11 | 3.06 |
| 2M17354093-1716200 | 3895.5 | 1.01 | -0.87 | 2.01 |
| 2M17390801-2331379 | 3740.4 | 0.83 | -0.81 | 2.34 |
| 2M17392719-2310311 | 3643.3 | 0.67 | -0.87 | 2.55 |
| 2M17473299-2258254 | 4018.3 | 0.47 | -1.71 | 2.12 |
| 2M17482995-2305299 | 4213.6 | 1.24 | -1.01 | 2.10 |
| 2M17483633-2242483 | 3651.5 | 0.44 | -1.09 | 2.57 |
| 2M17503263-3654102 | 3893.5 | 0.64 | -0.99 | 2.19 |
| 2M17552744-3228019 | 4018.9 | 1.0 | -1.05 | 1.99 |
| 2M18020063-1814495 | 3988.8 | 0.80 | -1.38 | 2.04 |
| 2M18050452-3249149 | 3940.8 | 0.77 | -1.16 | 2.08 |
| 2M18050663-3005419 | 3439.9 | 0.23 | -0.92 | 2.52 |
| 2M18065321-2524392 | 3893.1 | 0.95 | -0.89 | 2.02 |
| 2M18104496-2719514 | 4153.1 | 1.33 | -0.82 | 2.05 |
| 2M18125718-2732215 | 3617.2 | 0.44 | -1.31 | 2.64 |
| 2M18200365-3224168 | 3976.6 | 0.95 | -0.86 | 1.94 |
| 2M18500307-1427291 | 4076.0 | 1.23 | -0.94 | 1.73 |
| 2M17173248-2518529 | 3977.0 | 1.0 | -0.91 | 1.81 |
| 2M17285088-2855427 | 3838.0 | 0.63 | -1.20 | 2.18 |
| 2M17291778-2602468 | 3844.3 | 0.71 | -0.99 | 2.10 |
| 2M17301495-2337002 | 3814.0 | 0.69 | -1.06 | 2.22 |
| 2M17310874-2956542 | 4175.7 | 1.19 | -0.92 | 2.07 |
| 2M17382504-2424163 | 3880.4 | 0.99 | -1.05 | 1.55 |
| 2M17453659-2309130 | 4133.1 | 1.27 | -1.20 | 1.08 |
| 2M17511568-3249403 | 3921.2 | 0.98 | -0.90 | 2.04 |
| 2M17532599-2053304 | 3896.9 | 0.91 | -0.87 | 2.10 |
| 2M17552681-3334272 | 4051.0 | 1.08 | -0.89 | 1.98 |
| 2M18005152-2916576 | 4158.9 | 1.04 | -1.02 | 2.21 |
| 2M18010424-3126158 | 3773.1 | 0.68 | -0.83 | 2.20 |
| 2M18042687-2928348 | 4164.7 | 0.88 | -1.19 | 2.14 |
| 2M18044663-3132174 | 3832.6 | 0.92 | -0.90 | 2.22 |
| 2M18052388-2953056 | 4252.9 | 0.92 | -1.56 | 1.92 |
| 2M18080306-3125381 | 4310.0 | 1.57 | -0.90 | 1.48 |
| 2M18142265-0904155 | 3920.5 | 1.12 | -0.85 | 2.13 |
| 2M18195859-1912513 | 4102.0 | 1.05 | -1.22 | 1.78 |
| 2M17190320-2857321 | 4139.6 | 1.19 | -1.20 | 1.83 |
| 2M17224443-2343053 | 4058.3 | 1.02 | -0.88 | 1.97 |
| 2M17292082-2126433 | 3983.4 | 0.78 | -1.27 | 2.59 |
| 2M17293482-2741164 | 4143.5 | 1.03 | -1.25 | 1.85 |
| 2M17323787-2023013 | 3865.7 | 1.03 | -0.85 | 1.94 |
| 2M17330730-2407378 | 4042.5 | 0.25 | -1.87 | 1.88 |
| 2M17341796-3905103 | 4163.5 | 1.42 | -0.89 | 1.84 |
| 2M17342067-3902066 | 4380.4 | 1.40 | -0.90 | 1.99 |
| 2M17503065-2313234 | 3819.4 | 0.98 | -0.88 | 2.1 |
| 2M18023156-2834451 | 3617.4 | 0.42 | -1.19 | 3.02 |
| 2M18143710-2650147 | 4240.5 | 1.30 | -0.91 | 1.97 |
| 2M18150516-2708486 | 3833.4 | 1.0 | -0.82 | 2.14 |
| 2M18344461-2415140 | 4294.5 | 1.09 | -1.41 | 1.83 |

The atomic lines analysed initially were selected from Smith et al. (2021), Shetrone et al. (2015), Ce II lines identified by Cunha et al. (2017), and lines of Si I identified by Fanelli et al. (2021). Lines of Nd II (Hasselquist et al. 2016) and Yb II (Smith et al. 2021) were not studied. In Table 3 are reported the lines that we verified in the spectra of the 58 sample stars.

For the moderately metal-poor sample stars, some of the lines indicated in the articles above are not suitable, and in a few cases we have added other lines that we identified as suitable for the stellar parameters of the sample stars. The lines are discussed in detail below. In the present work, we adopt the ASPCAP abundances of Mg, Si, Ca, and revise the C, N, O, and Ce abundances; we also verified Ti lines and some comments are given, but the abundances are not used, given conflicting results from different lines. Other elements such as Na, Al, and iron-peak elements will be analysed elsewhere.

We identified and fitted the studied lines in the reference stars Arcturus and μ Leo, in order to check if the lines are well reproduced in these stars, and therefore reliable for deriving abundances in the sample stars. For the reference star Arcturus, we used the Hinkle, Wallace & Livingston (1995) atlas, and for the metal-rich reference giant star μ Leo a spectrum from APOGEE was used. The adopted stellar parameters for Arcturus and μ Leo are from Meléndez et al. (2003) and Zoccali et al. (2006) plus Lecureur et al. (2007), respectively.

Table 4 reports abundances in the Sun, Arcturus and μ Leo. For the Sun they are from (a) Grevesse, Noels & Sauval 1996; Grevesse et al. 1998, adopted, (b) Grevesse et al. (2015), Scott et al. (2015a, 2015b), (c) Lodders, Palme & Gail (2009). For Arcturus, the abundances are from Meléndez et al. (2003), Lecureur et al. (2007), Ramírez & Allende Prieto (2011), Barbuy et al. (2014), and Smith et al. (2013). For μ Leo, the abundances are from Gratton & Sneden (1990), Smith & Ruck (2000), Lecureur et al. (2007), Barbuy et al. (2015), Smith et al. (2013) or present fits, using the observed spectrum by Lecureur et al. (2007) in the optical.

According to Ashok et al. (2021), and Nidever et al. (2015) the average resolution of the APOGEE observations is $R \approx 22\,500$ based on a direct-measured FWHM of $\sim 0.7\text{ \AA}$, with 10–20 per cent variations seen across the wavelength range. We have employed a typical FWHM = 0.70 \AA , but to fit better different lines we varied the FWHM values from 0.60 to 0.75, from the lowest to the highest wavelengths. Note that the FWHM varies from fibers to fibers and with a fiber with wavelength.

3.1 C, N, O abundances

The abundances of C, N, and O are derived from CN, OH, and CO molecular lines. They are interdependent due to the molecular dissociative equilibrium. Since the molecular lines are spread all over the spectra, these abundances are derived first, and they are reported in Table 5.

Computing synthetic spectra employing the PFANT code described in Barbuy et al. (2018b), we have derived C, N, O abundances in two ways:

a) In the region 15144–16896 \AA , first we derive the O abundances by analysing the molecular lines of OH. Some of the most prominent OH lines in this region are at: 15264.60, 15266.160, 15278.516, 15281.045, 15719.687, 15893.524, 16074.151, 16662.187, 16872.265, 16895.164 \AA . These lines are the most sensitive to oxygen variation in the APOGEE

Table 3. Line list. log gf from VALD3 linelist (Piskunov et al. 1995; Ryabchikova et al. 2015), Kurucz (1993) and NIST (Martin et al. 2002). The log gf values for Ce II lines are from Cunha et al. (2017).

| Species | λ (Å) | χ_{ex} (eV) | log gf (VALD3) | log gf (Kurucz) | log gf (NIST) | Notes |
|-----------|------------------|---------------------|-------------------|--------------------|------------------|----------------------------|
| Si I | 15361.161 | 5.954 | -1.925 | -1.990 | -1.710 | |
| | 15376.831 | 6.222 | -0.649 | -0.290 | - | |
| | 15833.602 | 6.222 | -0.168 | -0.660 | -0.078 | Apogee gap |
| | 15960.063 | 5.984 | 0.107 | 0.130 | 0.197 | |
| | 16060.009 | 5.954 | -0.566 | -0.440 | -0.429 | |
| | 16094.787 | 5.964 | -0.168 | -0.110 | -0.078 | |
| | 16215.670 | 5.964 | -0.665 | -0.990 | -0.575 | |
| | 16680.770 | 5.984 | -0.140 | -0.500 | -0.090 | |
| | 16828.159 | 5.984 | -1.102 | -1.390 | -1.012 | |
| | Ca I | 16197.075 | 4.535 | 0.089 | 0.638 | - |
| 16204.087 | | 4.535 | -0.627 | -0.111 | - | |
| Ti I | 15543.756 | 1.879 | -1.120 | -1.273 | -1.080 | |
| | 15602.842 | 2.267 | -1.643 | -1.544 | - | |
| | 15698.979 | 1.887 | -2.060 | -2.218 | -2.020 | |
| | 15715.573 | 1.873 | -1.250 | -1.359 | -1.200 | |
| Ce II | 16635.161 | 2.345 | -1.807 | -2.178 | - | |
| | 15277.610 | 0.609 | -1.94 | - | - | too faint |
| | 15784.750 | 0.318 | -1.54 | - | - | |
| | 15829.830 | 0.320 | -1.80 | - | - | Apogee gap |
| | 15958.400 | 0.470 | -1.71 | - | - | |
| | 15977.120 | 0.232 | -2.10 | - | - | weak line strongly blended |
| | 16327.320 | 0.561 | -2.40 | - | - | |
| | 16376.480 | 0.122 | -1.79 | - | - | |
| | 16595.180 | 0.122 | -2.19 | - | - | |
| | 16722.510 | 0.470 | -1.65 | - | - | |

Table 4. Solar abundances from (1) Grevesse et al. (1996), Grevesse & Sauval (1998) (adopted); (2) Steffen et al. (2015); (3) Scott et al. (2015a, 2015b); (4) Grevesse et al. (2015); (5) Lodders et al. (2009); Arcturus abundances from: (6) Ramírez & Allende Prieto (2011), (7) McWilliam et al. (2013), (8) Leclercq et al. (2007), (9) Barbay et al. (2014), (10) Smith et al. (2013); (11) Cunha et al. (2017) μ Leo abundances from: (10) Smith et al. (2013); (12) Gratton & Sneden (1990), 13: Barbay et al. (2015), (14) Van der Swaelmen et al. (2016). The last column lists values from Smith et al. (2013).

| El. | Z | $\log \epsilon(X)_{\odot}$ | | [X/Fe] | $\log \epsilon(X)$ | [X/Fe] | $\log \epsilon(X)$ | | |
|-----|----|----------------------------|---------|---------------------|--------------------|--------|----------------------|------|----------|
| | | Sun | | Arcturus adopted | | | μ Leo adopted | | |
| Fe | 26 | 7.50 | 7.50 | 7.50 | -0.54 | 6.96 | +0.30 | 7.80 | 7.76[10] |
| C | 6 | 8.55[1] | - | 8.39[5] | -0.22[8] | 7.79 | -0.3[10] | 8.55 | 8.52[10] |
| N | 7 | 7.97[1] | - | 7.86[5] | +0.22[8] | 7.65 | +0.45[10] | 8.72 | 8.71[10] |
| O | 8 | 8.76[2] | - | 8.73[5] | 0.39[9] | 8.62 | +0.0[10] | 9.06 | 9.05[10] |
| Na | 11 | 6.33[1] | 6.21[3] | 6.30[5] | 0.11[6] | 5.90 | +0.50[8] | 7.13 | - |
| Mg | 12 | 7.58[1] | 7.59[3] | 7.54[5] | 0.37[6] | 7.41 | -0.03[10] | 7.85 | 7.85[10] |
| Al | 13 | 6.47[1] | 6.43[3] | 6.47[5] | 0.37[7] | 6.30 | +0.13[10] | 6.90 | 6.90[10] |
| Si | 14 | 7.55[1] | 7.51[3] | 7.52[5] | 0.33[6] | 7.34 | -0.10[10] | 7.75 | 7.76[10] |
| Ca | 20 | 6.36[1] | 6.32[3] | 6.33[5] | 0.11[6] | 5.93 | -0.04[10] | 6.62 | 6.62[10] |
| Sc | 21 | 3.17[1] | 3.16[3] | 3.10[5] | 0.23[6] | 2.86 | +0.10[11] | 3.57 | - |
| Ti | 22 | 5.02[1] | 4.93[3] | 4.90[5] | 0.26[7] | 4.74 | +0.10[10] | 5.42 | 5.40[10] |
| V | 23 | 4.00[1] | 3.89[3] | 4.00[5] | 0.12[7] | 3.58 | +0.03[12] | 4.33 | 4.18[10] |
| Cr | 24 | 5.67[1] | 5.62[3] | 5.64[5] | -0.05[6] | 5.08 | -0.01[12] | 5.96 | 6.14[10] |
| Mn | 25 | 5.39[1] | 5.42[3] | 5.37[5] | -0.14[7] | 4.71 | +0.00[12] | 5.69 | 5.79[10] |
| Co | 27 | 4.92[1] | 4.93[3] | 4.92[5] | +0.09[7] | 4.49 | +0.00[12] | 5.22 | 5.23[10] |
| Ni | 28 | 6.25[1] | 6.20[3] | 6.23[5] | 0.06[6] | 5.77 | +0.05[10] | 6.60 | 6.60[10] |
| Cu | 29 | 4.21[1] | 4.19[4] | 4.21[5] | -0.26[10] | 3.55 | -0.10[10] | 4.41 | 4.41[10] |
| Zn | 30 | 4.60[1] | 4.56[4] | 4.62[5] | +0.18[6] | 4.26 | -0.10[13] | 4.80 | - |
| Y | 39 | 2.24[1] | 2.21[4] | 2.21[5] | -0.30[9] | 1.40 | +0.04[14] | 2.58 | - |
| Zr | 40 | 2.60[1] | 2.59[4] | 2.58[5] | -0.28[7] | 1.78 | +0.10[12] | 3.00 | - |
| Ba | 56 | 2.13[1] | 2.25[4] | 2.17[5] | -0.30[9] | 1.29 | +0.10[14] | 2.53 | - |
| La | 57 | 1.22[1] | 1.11[4] | 1.14[5] | -0.30[9] | 0.38 | -0.37[14] | 1.15 | - |
| Ce | 58 | 1.55[1] | 1.58[4] | 1.61[5] | -0.45[11] | 0.99 | -0.37[14] | 1.15 | - |
| Eu | 63 | 0.51[1] | 0.52[4] | 0.52[5] | 0.23[7] | 0.20 | -0.14[14] | 0.67 | - |

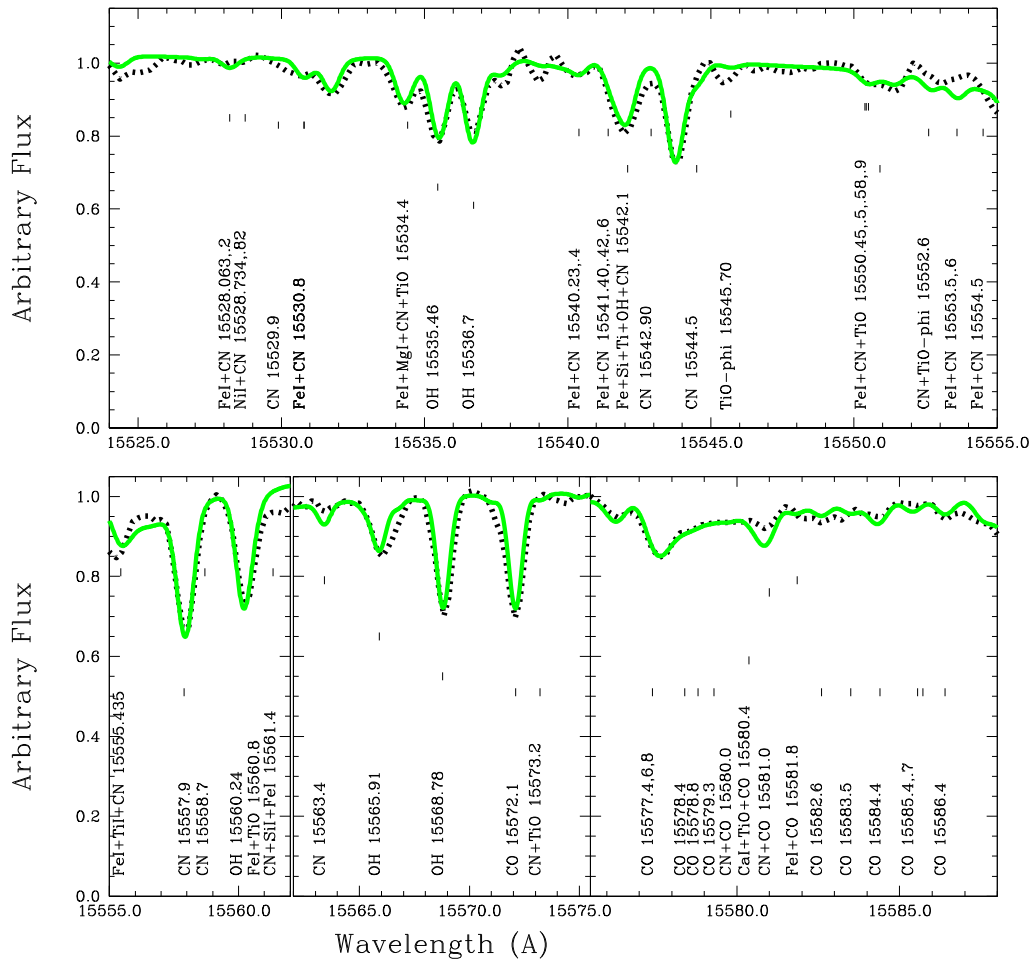


Figure 4. Star 2M17382504-2424163: Observed spectrum (black dotted) and synthetic spectrum computed with $[C, N, O/Fe] = -0.20, 0.30, 0.40$ (green).

sample. We derive C abundances by analysing the CO molecular lines, but there are not many strong CO lines in the range of 15100–17000 Å. In our sample, the strongest lines of CO, used to measure C abundances, are at 15983.214, 15985.598, 15990.420, 16016.081 Å. Next we see how the CN lines change when we modify Nitrogen. The most sensitive lines of CN are at 15162.648, 15222.382 Å. There are many CN lines in the region we are working with (especially in the range 15522–15600 Å), but most of them are too shallow to give reliable abundance measurements.

b) A derivation of CNO abundances using the region 15525–15595 Å, where there are clear lines of OH, and a clear bandhead of CO, as well as lines of CN, although less conspicuous, as done for example in Barbuy et al. (2021a) for Phoenix spectra that were observed in this region only. For these calculations a FWHM = 0.60 was adopted which is suitable for the wavelength region in question.

This is illustrated in Fig. 4 for star 2M17382504-2424163. Note the clear OH lines at 15535.46, 15565.91, 15566.78, and clear CO bandhead at 15577.4 Å.

We concluded that both methods a) and b) give very similar results within ± 0.1 dex.

A verification of these CNO abundances was carried out by fitting lines along all the spectra, in particular the lines of CO 15600.74, 15612.5, 15667.55 Å, where only for four stars the C abundance was decreased by -0.05 to -0.10 (stars 2M17330695-2302130, 2M18050663-3005419, 2M18125718-2732215, and

2M18344461-2415140), and for the others the fits were very satisfactory.

We then proceeded with the verification of the OH lines: OH 15130.921, 15266.168, 15281.052, 15409.172, 15568.78, 15651.896, 15719.696, 15755.522 Å, and CN 5181.277, 15298.487, 15308.893, 15318.74, 15337.959, 15341.508, 15432.811, 15447.095, 15466.235, 15481.868, 15530.776, 15684.088, 15737.445 Å. Only for star 2M18023156-2834451 we increased the oxygen abundance by 0.05 dex, noting that its spectra show larger lines than the others, needing a higher spectral convolution to be fitted.

Fits are shown for selected OH lines for star 2M17382504-2424163 in Fig. 5, and CO lines in Fig. 6 for stars 2M17382504-2424163 and 2M17511568-3249403.

Regions of CN lines are verified using wavelength regions indicated by Fernández-Trincado et al. (2020a, 2020b) for example. Fig. 7 are shown the fits to good CN lines. Among these, the clearest CN feature is at 15387.6 Å, and its fits are compatible with the C, N abundances from the 15283–15287, 15320–15330, and 15355–15380 Å regions. The feature at 15514 Å is blended with a Co I line and is less reliable. The N abundances derived are confirmed for about half the stars, whereas for the other half the N abundance was decreased by a mean of 0.2 dex: this is not surprising because the CN lines in the 15555 ± 50 Å from method b) are all faint and/or blended. Results from method b) above, together with these corrections, are adopted for C, N, O abundances.

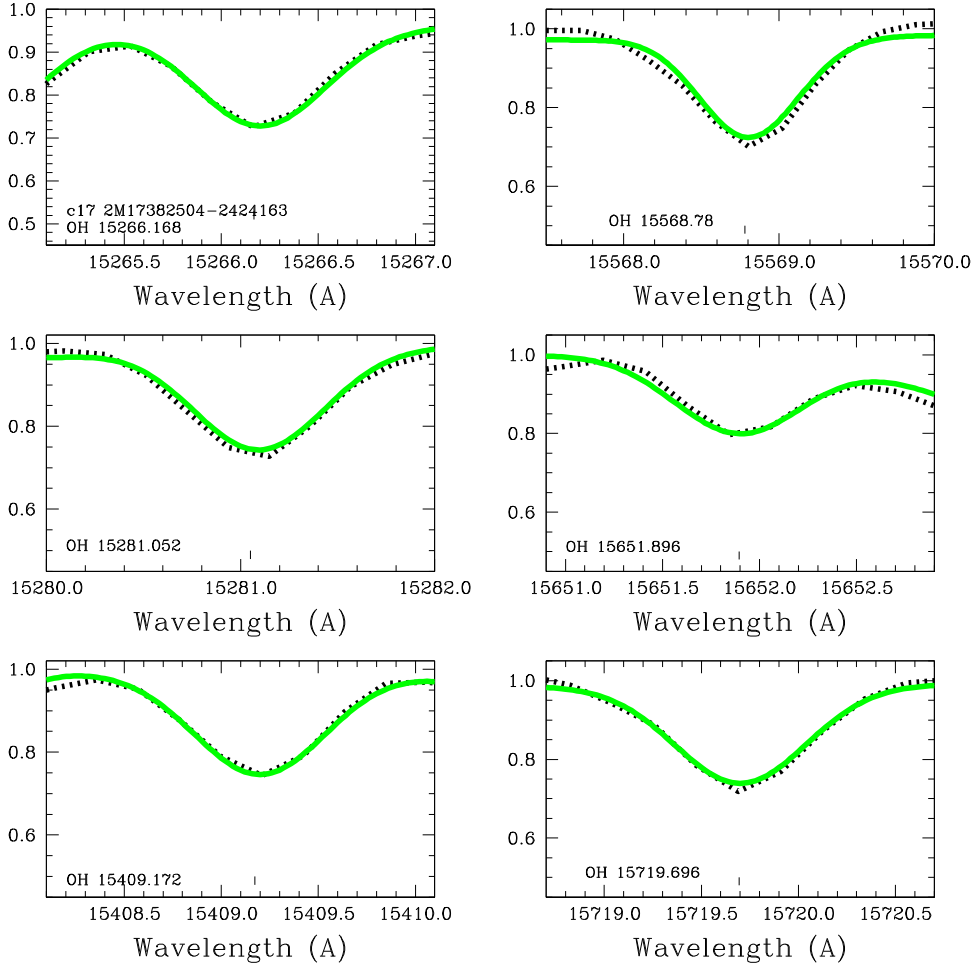


Figure 5. Star 2M17382504-2424163: Selected OH lines. Observed spectrum (black dotted) and synthetic spectrum computed with $[O/Fe] = 0.40$ (green).

The results of our manual analysis differ from the outputs of the ASPCAP pipeline for oxygen and, to a lesser degree nitrogen. Our methods a) and b) give rather similar results to each other within ± 0.05 dex, and with oxygen abundances somewhat higher than those derived with ASPCAP that appear to be too low for bulge stars. The uncertainties on the oxygen abundances were already discussed by Jönsson, Allende Prieto & Holtzman (2018), and Zasowski et al. (2019). Our oxygen abundances as compared with the DR17 ones are compatible within uncertainties, but with a trend to be higher.

In order to verify the reason for these differences, we carried out the fit to the N-rich star 2M17480576-2445000 analysed by Schiavon et al. (2017). With our method b) we have found that $[O/Fe] = 0.4$ instead of $[O/Fe] = 0.3$, and $[C/Fe] = -0.2$ instead of $[C/Fe] = 0.0$, and on the other hand the N enhancement of $[N/Fe] = 0.8$ is confirmed. Given the interplay between the CNO trio elements, it appears that the trend is to have somewhat lower C and higher O, and not much of a change in N abundances, in comparing our abundances with those from ASPCAP.

Note that none of the stars in our sample is N-enhanced, therefore they are good candidates to being similar to the first generation stars found in globular clusters.

3.2 alpha-elements Mg, Si, Ca, and Ti

We analyse the abundances of the α -elements Mg, Si, and Ca, and the iron-peak element Ti.

3.2.1 Magnesium, Silicon, Calcium and Titanium

Our fits with the original DR17 ASPCAP Mg abundances are in agreement with their results for Mg, Si, and Ca. Our calculations are in LTE with plane-parallel models, as is adopted by the original ASPCAP method. The DR17 results for Mg and Ca correspond to calculations in non-LTE (Osorio et al. 2020), and even so the compatibility is good for these elements.

The Si I lines reported in Table 3 are all suitably reproduced with the ASPCAP Si abundance with the exception of line Si I 15261.161 that is too shallow in the sample stars. Si abundance appears to be among the best determined ones by ASPCAP together with Mg.

The four Ca I lines listed in Smith et al. (2013) and that are used to ASPCAP, namely 16136.8, 16150.8, 16155.2, 16157.4 Å (see also Jönsson et al. 2018) are faint in the sample stars, and they are not fitted with the ASPCAP Ca abundance; instead they would need an extra 0.2 dex in Ca abundance to be fitted. In Table 3, we include another two lines of Ca I that we were able to identify as suitable for the metallicity of our stars: Ca I 16197.075, 16204.087 Å. The Ca I 16197.075 is well fitted in about half the stars, whereas in others it show blends, and finally the Ca I 16204.087 Å line is well fitted with the Ca abundance from ASPCAP. A FWHM = 0.65 fits better the lines. In conclusion, we adopted the ASPCAP Ca abundances, relying on the results for the Ca I 16204.087 Å line.

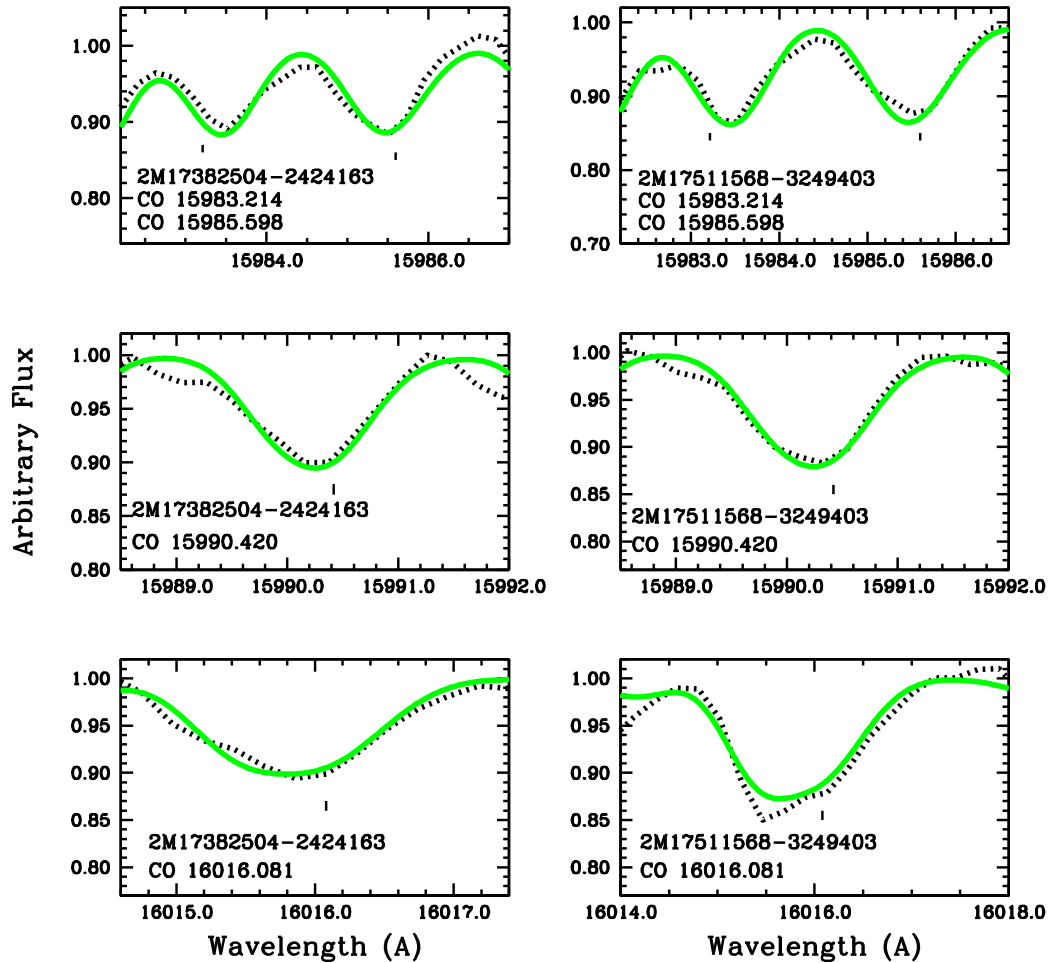


Figure 6. Stars 2M17382504-2424163 and 2M17511568-3249403: Selected CO lines. Observed spectra (black dotted) and synthetic spectra computed with $[C/Fe] = -0.2$ for both stars, and $[O/Fe] = 0.40$ and 0.38 , respectively (green).

3.2.2 Titanium

Among the five lines studied, only Ti I 15543.756 Å line is well fit in essentially all stars. Ti I 15698.979 Å tends to give the same value, but it is located in a blend with several other lines with a difficult continuum placement. Ti I 15715.753 Å tends to give either the value from ASPCAP or requires a lower Ti abundance, whereas Ti I 15602.842 and 16635.161 Å require higher values by about 0.3 ± 0.2 dex to be fitted. Because of the conflicting results from these different lines, and the fact that ASPCAP gives $[Ti/Fe] = 0.0$ for most stars, which is not compatible with the Si and Ca enhancements, we preferred not to analyse the Ti abundances in the sample stars.

Note that the lines Ti I 15602.842 and 16635.161 Å, that are only fitted with higher Ti abundances, have somewhat higher excitation potential than the other three inspected lines, and that means that there may be an effect of effective temperature.

3.3 s-process element Ce

We used six Ce II lines, among which Ce II 16722.510 Å line is well fit to almost all stars except for a few for which most of the other lines are fit with a lower value than with the best line (case of 2M17173693-2806495), followed by Ce II 15958.400 and 16595.180 Å lines, that are fit with the adopted value for almost all stars. The

Ce abundances from APOGEE DR17 and revised values are given in Table 6.

Ce II 15784.750 Å is fit for about half the stars, for a few would require lower Ce abundances and about 1/3 of them would require higher Ce abundances; 16327.320 Å is faint and is fit for about 1/3 of stars and 2/3 would require higher Ce abundances; Ce II 16376.480 Å would require higher values for about half the stars.

In eight cases all six lines can be considered well-fitted, as is the case of star 2M18500307-1427291, shown in Fig. 8.

For the fit of the Ce lines, we adopted $FWHM = 0.75$, which is suitable for the wavelength of the lines. The revised values are systematically higher than those resulting from ASPCAP (the fits to all stars are available under request).

DR17 used three Ce II windows covering the lines 15784, 16376, and 16595 Å. These are the stronger lines among the six that we used, and for all the sample stars it is clear that, from these three lines, a higher Ce abundance is needed.

We note that due to uncertainties in the Ce abundances in DR17 the APOGEE team has released internally to the collaboration a value added catalogue with revised abundances (Hayes et al. in preparation). This catalogue will be public to the community in a few months.

As for Nd, we found that the lines are not suitable for analysis, from fits to them in the reference stars Arcturus and μ Leo, therefore we disregarded this element in the present analysis.

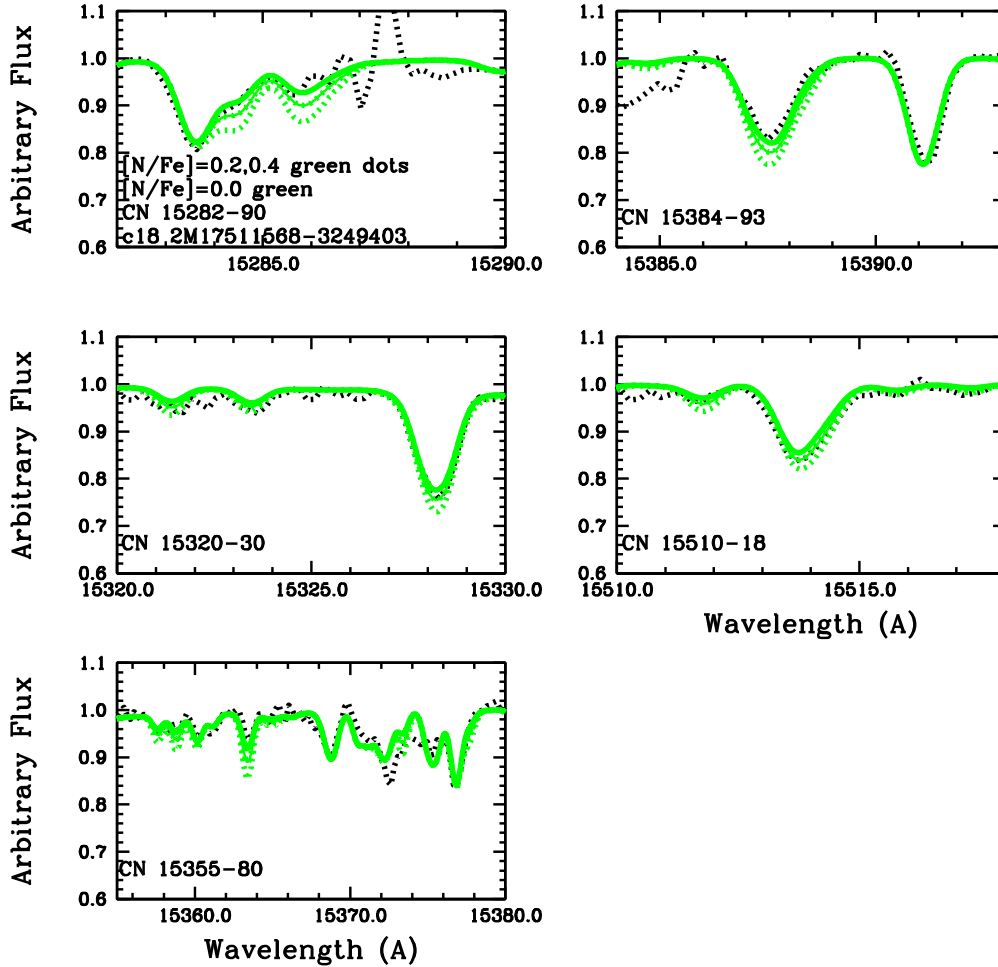


Figure 7. Star 2M17511568-3249403: regions containing CN lines. Observed spectra (black dotted) and synthetic spectra (green) computed with $[C/Fe] = -0.2$, $[N/Fe] = 0.0$, $[O/Fe] = 0.38$; green dotted lines correspond to calculations with $[N/Fe] = 0.2, 0.4$.

4 DISCUSSION

The α -element abundances in bulge stars provide us with a constraint on the formation history of its stellar populations: the formation time-scale. In other words, a mean $[\alpha/Fe] \sim 0.5$ in halo and bulge metal-poor stars of $[Fe/H] \lesssim -1.0$ indicates a fast chemical enrichment at early times, dominated by supernovae type II (SN II) (e.g. Woosley & Weaver 1995, hereafter WW95), whereas a lower $[\alpha/Fe]$ implies a slower enrichment, allowing supernovae type Ia to contribute to the enrichment of iron.

Moreover, as recently shown by Miglio et al. (2021) for a sample of Kepler stars with APOGEE spectra, stars with $[\alpha/Fe] > 0.2$ are all very old. The same probably applies to the present sample.

4.1 Oxygen and magnesium

Oxygen is produced by helium and neon burning in hydrostatic phases of the evolution of massive stars. Magnesium, together with Aluminum, are produced in hydrostatic carbon and neon burning (WW95). O and Mg are therefore the bona-fide α -elements produced by massive stars and ejected by supernova type II (SN II) event. They are enhanced relative to iron in old stellar populations, such as in bulge stars. In Fig. 9, upper panel are plotted the oxygen abundances reported by the original APOGEE DR17 release, and the revised val-

ues obtained as explained in Section 3. This Figure is readapted from that in Barbuy et al. (2018a), taking into account only the literature higher-resolution data (with a few exceptions) and data showing little abundance spread. The literature data taken into account are from Friaça & Barbuy (2017), that contains a revision of the abundances from Zoccali et al. (2006) and Lecœur et al. (2007), Cunha & Smith (2006); Alves-Brito et al. (2010), Fulbright, McWilliam & Rich (2007), only stars older than 11 Gyr from Bensby et al. (2013), Ryde et al. (2010) including a few of the same stars from Zoccali et al. (2006), Jönsson et al. (2017) including reanalysed 23 stars from Zoccali et al. (2006), Lecœur et al. (2007) and Friaça & Barbuy (2017), Siqueira-Mello et al. (2016), and metal-poor stars from García-Pérez et al. (2013), Howes et al. (2016) and Lamb et al. (2017).

Fig. 9 (lower panel) gives $[Mg/Fe]$ versus $[Fe/H]$ for metal-poor stars from García-Pérez et al. (2013), Howes et al. (2016), Lamb et al. (2017), Casey & Schlafman (2015), Koch et al. (2016), Fulbright et al. (2007), as corrected by McWilliam (2016), Alves-Brito et al. (2010), Hill et al. (2011), Bensby et al. (2017) for stars older than 8 Gyr, Johnson et al. (2014), Ryde et al. (2010, 2016), Siqueira-Mello et al. (2016), Jönsson et al. (2017) and Rojas-Arriagada et al. (2017).

Our fits show agreement with the ASPCAP Mg abundances, and they are compatible with the Mg abundances of other samples of bulge stars. The different model lines in Fig. 9 correspond to different radii from the Galactic centre.

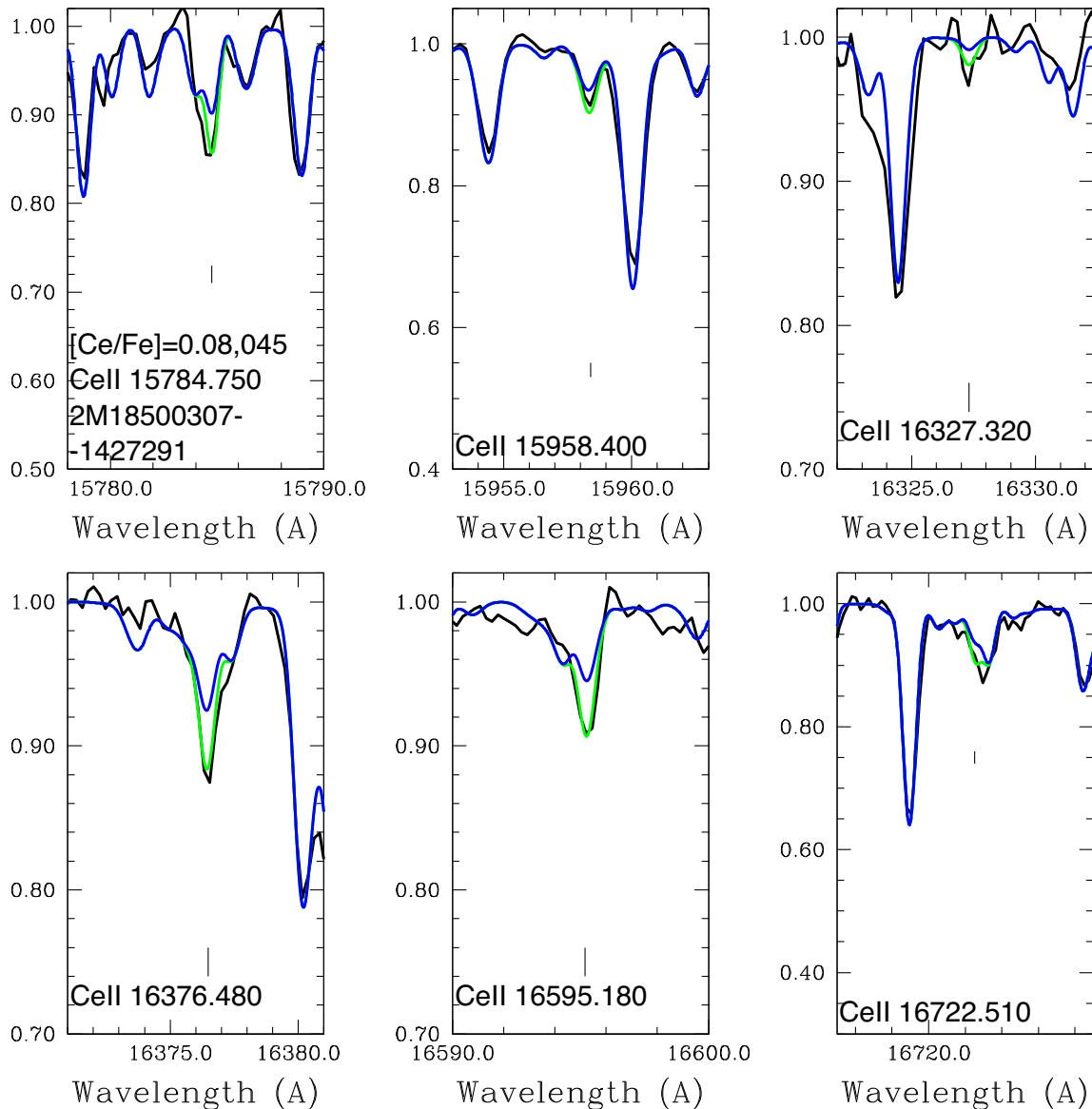


Figure 8. Star 2M18500307-1427291: fit to the six cerium lines. Observed spectrum: black; synthetic spectra are: blue with original ASPCAP $[\text{Ce}/\text{Fe}] = 0.08$ Ce abundance, green with final Ce abundance.

4.2 Chemodynamical evolution model for O and Mg

We have compared the abundances derived from observations with the predictions of chemodynamical evolution models for the bulge (Friaça & Barbuy 2017), described as a classical spheroid. It is assumed a baryonic mass of $2 \times 10^9 M_{\odot}$, and a dark halo mass $M_H = 1.3 \times 10^{10} M_{\odot}$. One central parameter of the model is the specific star formation rate ν_{SF} (i.e. the inverse of the star formation time-scale).

In the nucleosynthesis prescriptions of our model, we adopt the metallicity dependent yields from core-collapse supernovae (SNe II) from WW95 with some modifications following suggestions of Timmes, Woosley & Weaver (1995). For low metallicities ($Z < 0.01 Z_{\odot}$), we included the yields from high explosion-energy hypernovae (HNe) (Nomoto et al. 2013, and references therein). The type Ia supernovae yields are from Iwamoto et al. (1999) – their models W7 (progenitor star of initial metallicity $Z = Z_{\odot}$) and W70 (zero initial metallicity). The yields for intermediate mass stars ($0.8\text{--}8 M_{\odot}$)

with initial $Z = 0.001, 0.004, 0.008, 0.02,$ and 0.4 come from van den Hoek & Groenewegen (1997) (variable η_{AGB} case).

As we can see from Fig. 9, the abundances derived here both for the oxygen (upper panel) and for the magnesium (lower panel) are well reproduced by the chemodynamical model with $\nu_{\text{SF}} = 1 \text{ Gyr}^{-1}$ (star formation time-scale of 1 Gyr). Once more this suggests these objects to be very old.

4.3 Silicon and calcium

Si and Ca are mainly produced during the explosive nucleosynthesis of SN II events (WW95; McWilliam 2016) with smaller contributions from supernovae type Ia (SNIa).

The α -elements Si and Ca are plotted in Fig. 10 for the 58 sample stars together with literature data from García-Pérez et al. (2013), Howes et al. (2016), Lamb et al. (2017), Casey & Schlafman (2015), Koch et al. (2016), Alves-Brito et al. (2010), Bensby et al. (2017) for

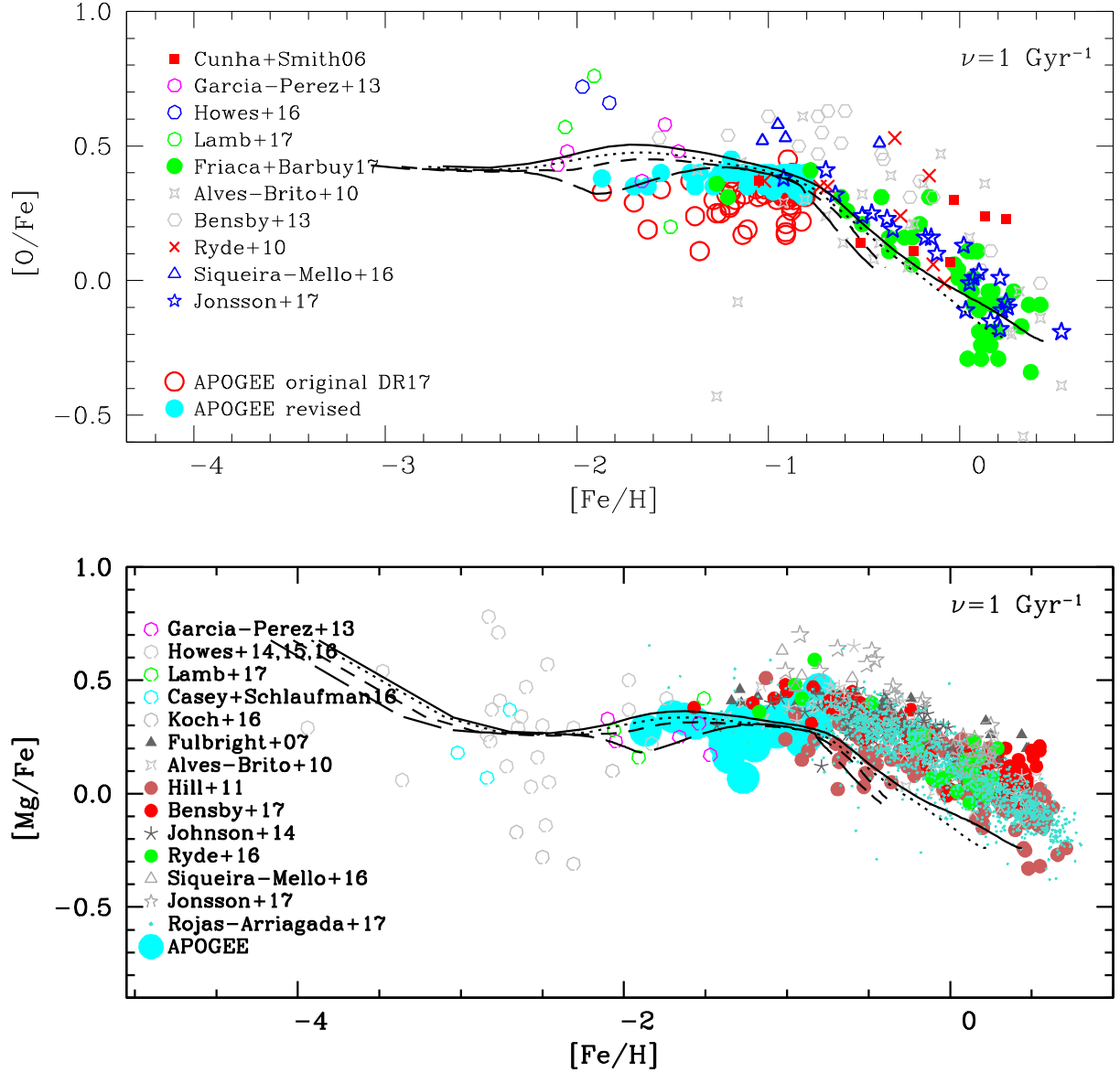


Figure 9. $[O/Fe]$ versus $[Fe/H]$ (upper panel) and $[Mg/Fe]$ versus $[Fe/H]$ (lower panel), for literature bulge field stars and the APOGEE abundances (original and revised in the case of oxygen) for the 58 sample stars. Symbols: grey 4-pointed stars: Alves-Brito et al. (2010); red filled circles: Bensby et al. (2013); red filled circles: Bensby et al. (2017); grey open pentagons: Casey & Schlaufman (2015); strong-grey filled triangles: Fulbright et al. (2007); magenta open pentagons: García-Pérez et al. (2013); red filled squares: Cunha & Smith (2006); indianred filled circles: Hill et al. (2011); grey open pentagons: Howes et al. (2016); grey stars: Johnson et al. (2014); grey 5-pointed stars: Jönsson et al. (2017); grey open pentagons: Koch et al. (2016); green open pentagons: Lamb et al. (2017); red crosses: Ryde et al. (2010); green filled circles: Ryde et al. (2016); turquoise 5-pointed stars: Rojas-Arriagada et al. (2017); grey open triangles: Siqueira-Mello et al. (2016); blue open circles: APOGEE original; cyan filled circles: final abundances for the 58 APOGEE sample stars. The oxygen abundances are normalized in terms of adopted solar abundances as explained in Friaça & Barbuy (2017). Chemodynamical evolution models from Friaça & Barbuy (2017) with formation time-scale of 1 Gyr, for several radii, are overplotted: $r < 0.5$ kpc (solid lines), $0.5 < 1$ kpc (dotted lines), $1 < r < 2$ kpc (short-dashed lines), $2 < r < 3$ kpc (long-dashed lines).

stars older than 8 Gyr, Ryde et al. (2016) and Siqueira-Mello et al. (2016).

Fig. 10 shows that a typical star formation time-scale of 1 Gyr (the chemodynamical model with $\nu_{SF} = 1 \text{ Gyr}^{-1}$) also explains the Si and Ca abundances found in the bulge.

From this figure, we can conclude that there are no differences in the Si and Ca abundances of the present confirmed *in situ* samples of bulge stars, and previous samples in bulge regions, for which no precise distances were available.

4.4 Heavy element Ce

In Fig. 11 are shown the revised Ce abundances in contrast with the lower original DR17 APOGEE abundances, together with results for M62 from Yong et al. (2014) and for field bulge stars from van der Swaelmen et al. (2016) and Lucey et al. (2022). As can be seen from the figure, we have found that the sample stars are enhanced in Cerium by about a mean value of $[Ce/Fe] \sim 0.4$. We find Ce enhancements relative to the DR17 and as well to results from the

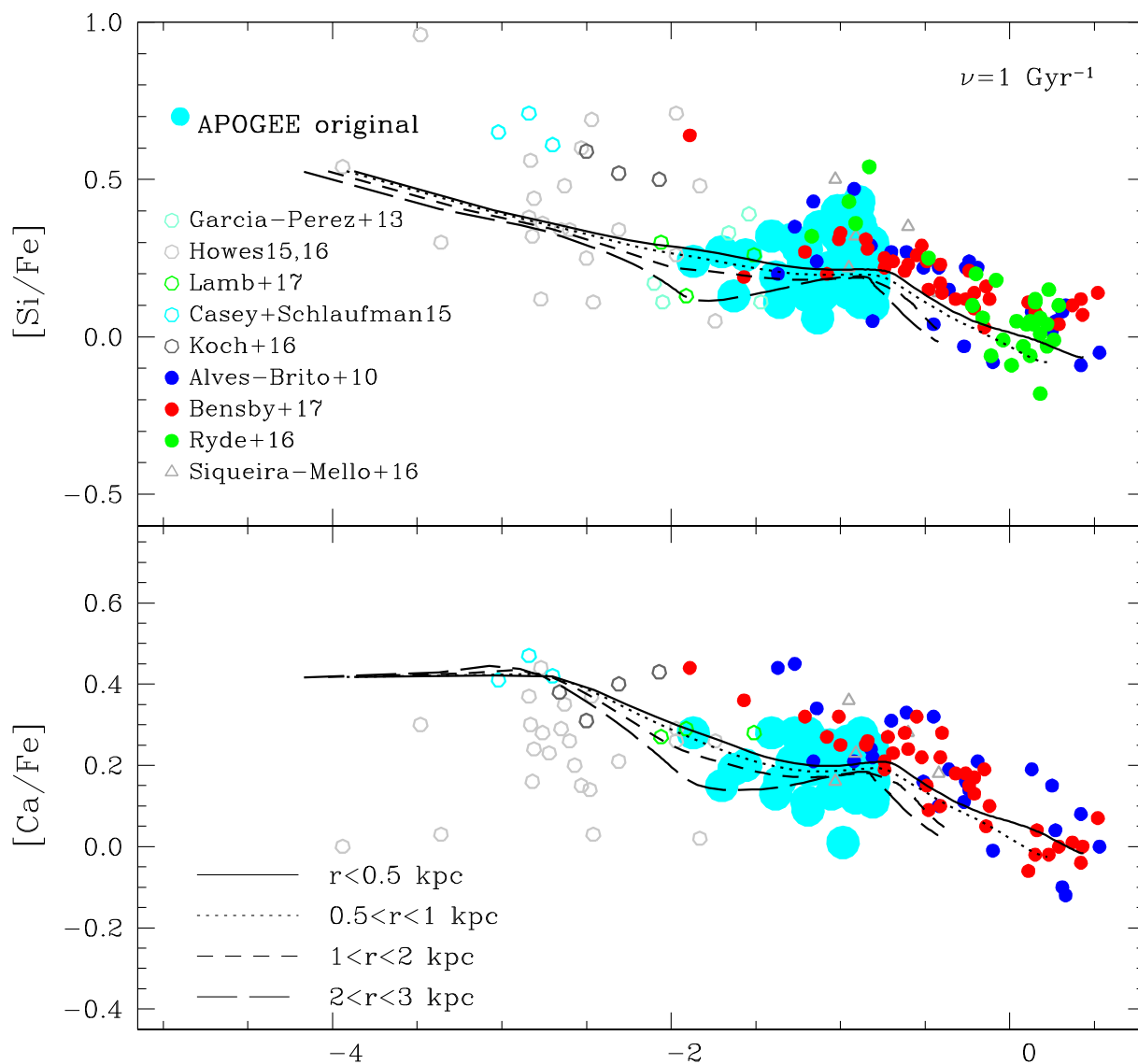


Figure 10. [Si, Ca/Fe] versus [Fe/H] for literature bulge field stars and the APOGEE abundances for the 58 sample stars. Symbols: grey 4-pointed stars: Alves-Brito et al. (2010); red filled circles: Bensby et al. (2017); grey open pentagons: Casey & Schlaufman (2015); magenta open pentagons: García-Pérez et al. (2013); grey open pentagons: Howes et al. (2016); grey open pentagons: Koch et al. (2016); green open pentagons: Lamb et al. (2017); green filled circles: Ryde et al. (2016); grey open triangles: Siqueira-Mello et al. (2016); blue open circles: APOGEE original; cyan filled circles: final abundances for the 58 APOGEE sample stars. The lines are the predictions of the chemodynamical models of Friaça & Barbuy (2017) with a formation timescale of 1 Gyr for several radii.

COMBS survey by Lucey et al. (2022). Clearly further investigation on Ce abundances in metal-poor bulge stars is needed.

Ce ($Z = 58$, $A = 140$) is essentially an element mostly formed by the s -process, with a fraction of 0.186 as r -element and 0.814 as s -element (Simmerer et al. 2004). Ce appears to be overproduced in massive spinstars (Frischknecht et al. 2016). On the other hand, since we do not have the ages of these stars, we cannot exclude that the Ce enhancement could be due to a mass transfer from a companion Asymptotic Giant Branch (AGB) star (e.g Bisterzo et al. 2011; Cristallo et al. 2015).

5 CONCLUSIONS

We have selected 58 stars from the bulge sample of Queiroz et al. (2021) based on APOGEE with characteristics to belong to

the spheroidal pressure-supported bulge. For this sample, we have analysed lines of C, N, O, alpha-elements Mg, Si, Ca, and neutron-capture element Ce. Our fits to the Mg, Si, and Ca abundances from the original APOGEE results using the ASPCAP software appeared to be in good agreement. We recomputed abundances for C, N, O, and Ce, assuming the spectroscopic non-calibrated stellar parameters from APOGEE DR17. We report differences in abundances of these elements.

We compare the abundances of these elements to literature data for bulge stars and chemodynamical models by Friaça & Barbuy (2017) – see also Barbuy et al. (2018a). These comparisons show compatibility of the abundances of the sample stars with literature and models for Mg, Si, and Ca in which a pressure supported component (spheroidal bulge) formed on a very short time-scale (below 1 Gyr). Similar results were suggested by other chemical evolution models

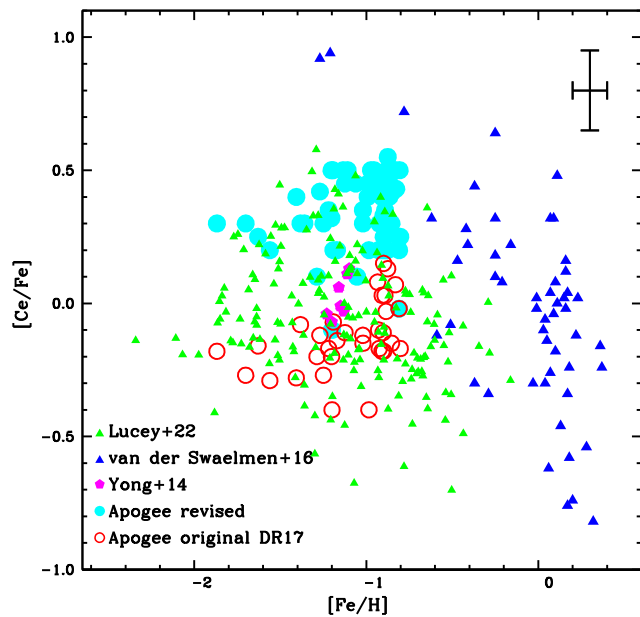


Figure 11. $[\text{Ce}/\text{Fe}]$ versus $[\text{Fe}/\text{H}]$ for literature bulge field stars and the APOGEE abundances for the 58 sample stars. Symbols: red filled triangles: van der Swaelmen et al. (2016); green filled triangles: Lucey et al. (2022); magenta pentagons: M62 from Yong et al. (2014); open blue circles: APOGEE DR17 $[\text{Ce}/\text{Fe}]$ values and cyan filled circles: revised abundances for the 58 APOGEE sample stars. Error bars are indicated in the right upper corner.

(see Matteucci 2021 for a review), and for stars with similar alpha element enhancements with asteroseismic ages (Miglio et al. 2021).

Nitrogen abundances show no exceptional enhancement for any of the sample stars therefore, there is no evidence for these stars to be a result of multiple stellar populations in dissolved globular clusters. The Ce abundance is enhanced in all stars, which would point out to a s-process origin of this element already in the very early phases of chemical enrichment. This could have been achieved with spinstars (e.g. Chiappini et al. 2011; Frischknecht et al. 2016), or alternatively due to mass transfer from a companion AGB star (e.g. Cristallo et al. 2015). This same conclusion was reached by Barbuy et al. (2009, 2014, 2021b) regarding the globular cluster NGC 6522, but here, since all the present sample stars are enhanced in Ce, all of them would have to be binaries with an AGB companion. Therefore the spinstars seem to be a more plausible explanation.

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DATA AVAILABILITY

The observed data are from the APOGEE survey. The calculations and fits to the lines can be requested to the authors. The code for spectrum synthesis is available to be retrieved.

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