

Chemical analysis of a triple system of A-type stars

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Abstract. Components of multiple systems generally originate from the same protostellar environment. Their similarities or differences in surface chemical composition therefore relates to their individual evolutionary paths (stellar evolution, rotation) and the possible influence of a close companion.

DG Leo is a spectroscopic hierarchical triple system with almost equal-mass late-A type components. Observations with high time and high spectral resolution were used to disentangle the spectrum of each component from the composite spectra. A detailed abundance analysis of the component spectra reveals that the wide companion has a nearly solar-like composition, while both components in the close binary show Am type peculiarities.

Keywords. Stars: abundances, stars: fundamental parameters, binaries: spectroscopic, stars: variable: δ Sct, stars: individual: (DG Leo).

1. Introduction

The analysis of the chemical composition of multiple systems having at least one pulsating component is an ideal tool to explore in an empirical way the interactions that may or may not exist between pulsation, diffusion, rotation and binarity. In the present study, we are, therefore, dealing with the chemical analysis of DG Leo (HD 85040, HR 3889, HIP 48218, Kui 44), a known multiple system with at least one pulsating component. DG Leo consists of a close binary (components Aa and Ab in a circular orbit) and one distant companion (component B). All three components were classified as late A-type stars. The orbital period of the Aa,b system is 4.15 days (Danziger & Dickens 1967; Fekel & Bopp 1977) while the orbital period of the visual pair Aab,B was estimated to be roughly 200 years (Fekel & Bopp 1977). All three components of the system are potential candidates for pulsations as well as for Am peculiarities (they are located in the cooler part of the δ Scuti instability strip). In fact multiple short-period oscillations of type δ Scuti with periodicities of about 2 hrs have been detected (Lampens *et al.* 2005).

2. Observations

Extensive spectroscopy was obtained at the Haute-Provence Observatory (OHP) with the ELODIE spectrograph on the 1.93-m telescope. Observations were collected during 7 of the 8 allocated nights in 2003 (January 3 – 7 & January 11 – 15) and cover about 50% of the close binary's (components Aab) orbital phase. The time exposure was fixed at 360 s to resolve the presumed p-mode pulsations, yet with a good signal-to-noise ratio and a high resolution (about 50000). 245 spectra covering $\lambda\lambda$ 3900 to 6800 were obtained with a signal-to-noise ratio that generally varies between 100 to 180 at λ 5500.

3. Spectral disentangling

To extract the individual contribution of the three components, we adopted the spectral disentangling technique introduced by Hadrava (1995, see references therein) and applied in the KOREL computer code (Release 21.3.99). KOREL assumes that the observed spectrum is composed by n (three in our case) time-independent intrinsic spectra moving relatively to each other. It fits the Fourier Transform of the observations at different orbital phases by a Least-Squares Method to provide the time-averaged spectrum of each component (see Fig. 1) and their relative radial velocities.

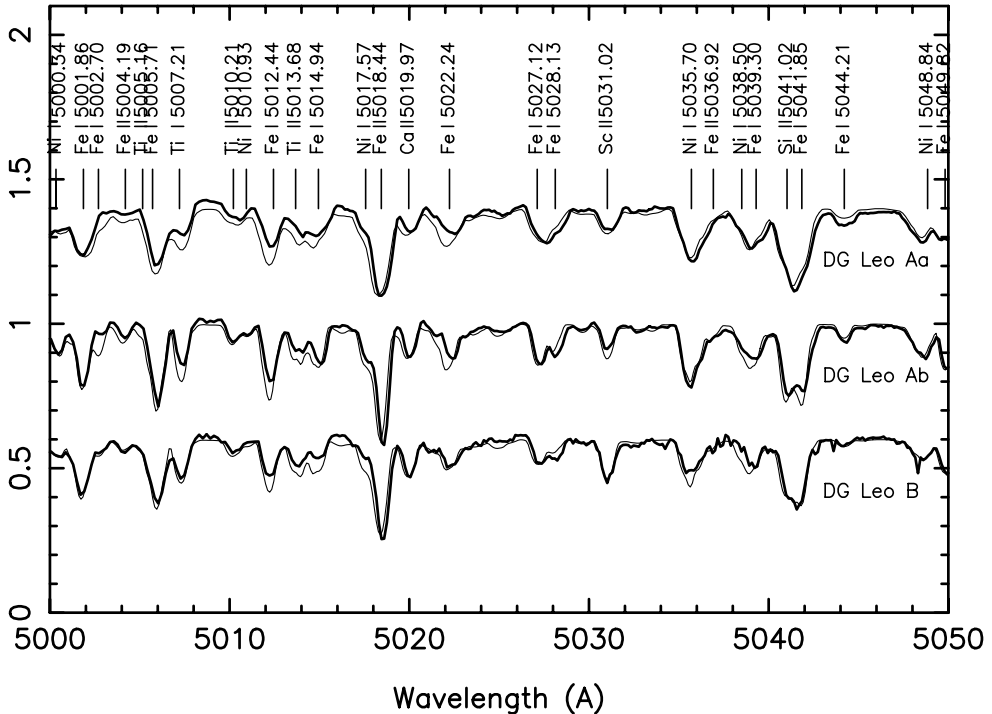


Figure 1. Component spectra. Comparison between the disentangled component spectra (thick line) and the synthetic spectra resulting from the chemical abundance analysis (thin line).

4. Fundamental parameters

The fundamental parameters that describe the components of DG Leo were derived, using the disentangled component spectra, in three consecutive steps. **1.** The projected rotation velocity, $v \sin i$, was obtained performing a Fourier analysis of several unblended line profiles and using a procedure similar to the one intensively applied by Royer *et al.* (2002). **2.** We estimated the effective temperature, T_{eff} , of DG Leo's components by fitting the individual $H\alpha$ and $H\gamma$ lines with theoretical line profiles computed in LTE using the SYNSPEC computer code (Hubeny & Lanz 1995, and references therein) combined to ATLAS9 (Kurucz 1993; Castelli *et al.* 1997) model atmospheres. **3.** The surface gravity, $\log g$, of the stars was determined from the HIPPARCOS parallax according to a method described by North *et al.* (1997). The luminosity was calculated using the parallax and the component's V magnitude while the stellar mass and radius were then interpolated in the theoretical evolutionary tracks ($Z = 0.02$) of Schaller *et al.* (1992).

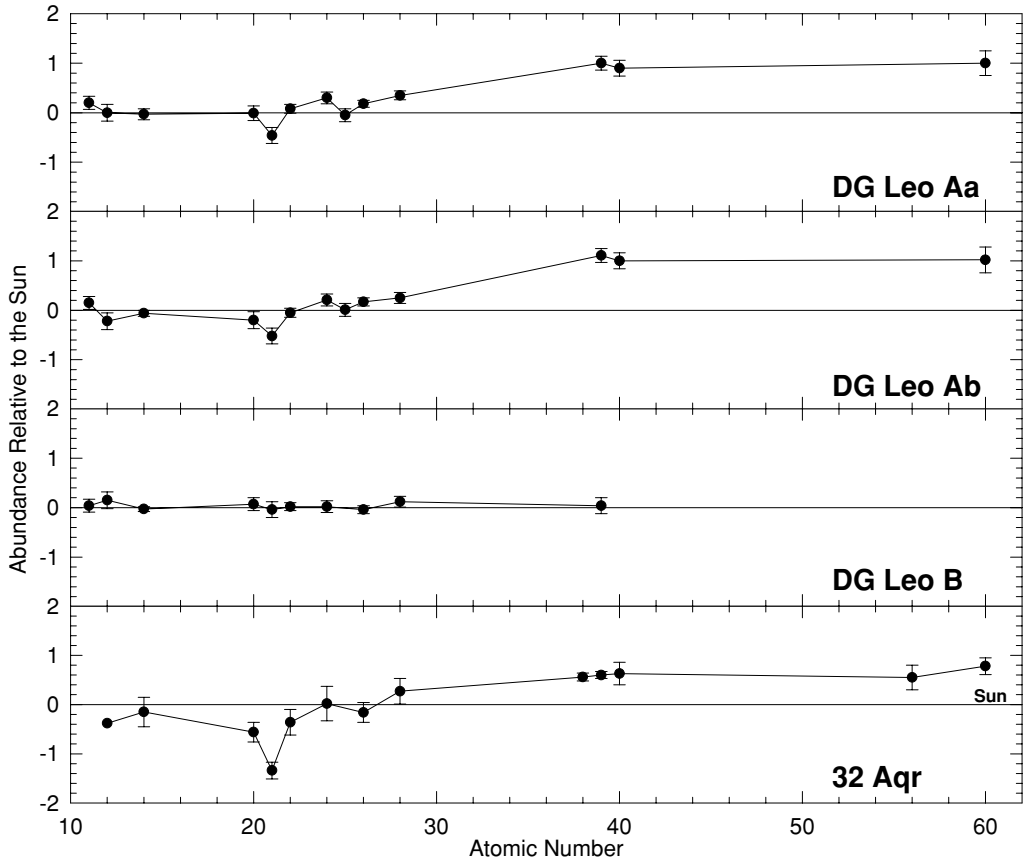


Figure 2. Chemical composition of DG Leo’s components compared to the Sun and to the Am star 32 Aqr (Kocer *et al.* 1993). Component B has a solar-like chemical composition while both A components show abundance patterns typical of Am stars.

Table 1. Component fundamental parameters. $v \sin i$, and $\log g$ were derived using, respectively, the classical Fourier transform technique. Hydrogen line modelisation and theoretical evolution tracks. ξ_{turb} is derived from the abundance analysis.

component	Aa	Ab	B
T_{eff} (K)	7470 ± 220	7390 ± 220	7590 ± 220
$\log g$	3.8 ± 0.14	3.8 ± 0.14	3.8 ± 0.12
$V \sin i$ (km s ⁻¹)	42 ± 2	28 ± 2	31 ± 3
ξ_{turb} (km s ⁻¹)	2.3 ± 0.5	2.3 ± 0.5	2.5 ± 0.5
M_{HR} (M_{\odot})	2.0 ± 0.2	2.0 ± 0.2	2.1 ± 0.2

5. Chemical composition

The atmospheric abundance analysis was then carried out by fitting theoretical line profiles to the observed component’s spectra (see Fig. 1). Oscillator strengths, energy levels and damping parameters (including Stark, van der Waals and natural broadening) used during the fitting procedure were basically those compiled in the VALD-2 database (Kupka *et al.* 1999) updated by Erspamer & North (2002). The final results are plotted in Fig. 2.

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