Magnetic field topology, chemical spot distributions, and photometric variability of the Ap star $\varphi$ Draconis

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ABSTRACT

The primary component of the multiple star $\varphi$ Dra is one of the brightest magnetic chemically peculiar stars in the northern sky. Here, we report results of a comprehensive study of the rotational photometric variability, binarity, magnetic field geometry, and surface chemical spot structure for this star. We derived a precise photometric rotational period of $1.71650213(21)$ d based on 1 yr of TESS nearly continuous space observations and discovered modulation of the stellar light curve with the phase of the 127.9-d binary orbit due to the light time effect. We revised parameters of the binary orbit and detected spectroscopic contribution of the secondary. A tomographic mapping technique was applied to the average intensity and circular polarization profiles derived from Narval high-resolution spectropolarimetric observations. This analysis yielded a detailed map of the global magnetic field topology together with the surface distributions of Si, Cr, and Fe abundances. Magnetic mapping demonstrates that the surface field structure of $\varphi$ Dra is dominated by a distorted dipolar component with a peak field strength of 1.4 kG and a large asymmetry between the poles. Chemical maps show an enhancement of Cr, Fe and, to a lesser extent, Si in a series of spots encircling intersections of the magnetic and rotational equators. These chemical spot geometries do not directly correlate with either the local field strength or the field inclination.

Key words: stars: atmospheres – stars: chemically peculiar – stars: individual: $\varphi$ Dra – stars: magnetic field – starspots.

1 INTRODUCTION

The upper main sequence magnetic chemically peculiar (ApBp or mCP) stars are A- and B-type stars distinguished by large underabundances and overabundances of chemical elements in their surface layers and the presence of a strong, stable, globally organized magnetic field. This chemical and magnetic structure is frozen into the stellar plasma and rotates with the star, giving rise to a characteristic synchronous rotational variability of the brightness in most photometric passbands, spectral energy distribution, individual line profiles, and magnetic observables. This unique combination of unusual properties and pronounced variability makes ApBp stars very useful, and sometimes the only suitable objects, for detailed studies of the origin, structure and evolution of the fossil magnetic fields in radiative stellar envelopes (e.g. Braithwaite & Nordlund 2006; Landstreet et al. 2007; Kochukhov, Shultz & Neiner 2019), and an interplay between the mass loss and rotational spin-down (e.g. ud-Doula, Owoki & Townsend 2009; Shultz et al. 2019), and microscopic chemical transport processes responsible for the observed chemical anomalies and starspot formation (e.g. Michaud, Alecian & Richer 2015; Kochukhov 2017; Ryabchikova & Romanovskaya 2017).

With a V magnitude of 4.22, $\varphi$ Dra (43 Dra, HD 170000, HR 6920, HIP 89908, TIC 356822358) is one of the brightest magnetic chemically peculiar stars. It is a rapidly rotating object showing a coherent spectroscopic (Jamar 1977; Kuschnig 1998), photometric (Musielok et al. 1980; Přiká et al. 2015; Bernhard, Hämmerich & Paunzen 2020), and magnetic (Landstreet & Borra 1977; Sikora et al. 2019b) variability with a period of $\approx 1.72$ d. This star is, in fact, a triple system with the visual secondary (component B) orbiting the primary (component A) on a 308-yr orbit with a semimajor axis of $< 1$ arcsec (Liška 2016). The magnitude difference between these two components is measured to be 1.3–1.5 mag depending on wavelength (Fabricius & Makarov 2000; Rutkowski & Waniak 2005). The primary is itself a single-line eccentric spectroscopic binary (components Aa and Ab) with a period of 128 d (Liška 2016; Bischoff et al. 2017). Multiplicity of $\varphi$ Dra is commonly ignored by studies of the magnetic Ap star (component Aa), since no evidence of the spectroscopic contribution of components Ab or B has been reported in the literature.

The global magnetic field of $\varphi$ Dra was first detected by Landstreet & Borra (1977) using the Balmer-line mean longitudinal magnetic field, $\langle B_{z} \rangle$, measurements with a typical uncertainty of 100 G. That study established a smooth, sinuousoidal magnetic field curve with $\langle B_{z} \rangle$ changing from about $-300$ to $+500$ G and interpreted these observations with a 3 kG-centred dipolar magnetic field topology with a low obliquity ($\beta \approx 20^\circ$) relative to the rotational axis. Considerably more precise longitudinal magnetic field measurements were obtained for $\varphi$ Dra by Sikora et al. (2019b) from metal lines in high-resolution circular polarization spectra. Although the inferred longitudinal field curve was fully compatible with the earlier hydrogen line results, Sikora et al. (2019b) derived significantly...
Table 1. Log of Narval spectropolarimetric observations of $\varphi$ Dra. The columns give the UT and heliocentric Julian dates of mid-exposure for each observation, the S/N per 1.8 km s$^{-1}$ velocity bin at 5500 Å, the rotational and orbital phases, the mean longitudinal magnetic field determined from LSD profiles, and the radial velocity measured from the hydrogen line cores. The average uncertainty of the latter measurements is 0.73 km s$^{-1}$.

| UT date      | HJD     | S/N   | Rotational phase | Orbital phase | $\langle B_\perp \rangle$ (G) | $V_H$ (km s$^{-1}$) |
|--------------|---------|-------|------------------|---------------|-------------------------------|---------------------|
| 2016-08-20   | 2457621.401 | 986   | 0.348            | 0.293         | $-236 \pm 26$                | $-24.2$            |
| 2016-08-21   | 2457622.368 | 1176  | 0.912            | 0.301         | $387 \pm 24$                 | $25.5$             |
| 2016-08-22   | 2457623.417 | 891   | 0.523            | 0.324         | $-170 \pm 33$                | $23.6$             |
| 2016-08-23   | 2457624.421 | 1069  | 0.108            | 0.332         | $407 \pm 25$                 | $24.1$             |
| 2016-08-24   | 2457625.414 | 1166  | 0.686            | 0.340         | $191 \pm 27$                 | $23.7$             |
| 2016-08-29   | 2457630.448 | 970   | 0.619            | 0.074         | $-2 \pm 32$                  | $-21.0$            |
| 2016-08-31   | 2457631.517 | 1068  | 0.242            | 0.363         | $-102 \pm 24$                | $-19.6$            |
| 2016-09-02   | 2457634.387 | 1094  | 0.914            | 0.081         | $430 \pm 26$                 | $-19.7$            |
| 2016-09-03   | 2457635.367 | 947   | 0.485            | 0.371         | $-76 \pm 30$                 | $-17.7$            |
| 2016-09-06   | 2457638.397 | 1033  | 0.250            | 0.089         | $-82 \pm 24$                 | $-16.9$            |
| 2016-09-07   | 2457639.404 | 1073  | 0.836            | 0.097         | $308 \pm 28$                 | $-17.8$            |
| 2016-09-08   | 2457640.391 | 1142  | 0.411            | 0.601         | $-286 \pm 24$                | $-15.4$            |
| 2016-09-11   | 2457643.392 | 1115  | 0.160            | 0.609         | $232 \pm 24$                 | $-15.9$            |
| 2016-09-12   | 2457644.310 | 1048  | 0.695            | 0.617         | $141 \pm 30$                 | $-16.5$            |
| 2016-12-11   | 2457734.237 | 834   | 0.084            | 0.624         | $437 \pm 33$                 | $-43.2$            |
| 2016-12-12   | 2457735.235 | 1057  | 0.666            | 0.192         | $109 \pm 30$                 | $-40.7$            |
| 2016-12-13   | 2457736.234 | 1106  | 0.248            | 0.198         | $-112 \pm 23$                | $-38.7$            |
| 2016-12-14   | 2457737.236 | 1162  | 0.831            | 0.207         | $364 \pm 26$                 | $-37.6$            |
| 2017-02-17   | 2457801.722 | 1027  | 0.400            | 0.215         | $-291 \pm 26$                | $-7.2$             |
| 2017-02-18   | 2457802.710 | 1136  | 0.976            | 0.223         | $421 \pm 25$                 | $-9.2$             |
| 2017-02-19   | 2457803.709 | 1075  | 0.557            | 0.262         | $-68 \pm 28$                 | $-8.6$             |
| 2017-02-20   | 2457804.699 | 944   | 0.134            | 0.271         | $350 \pm 28$                 | $-8.2$             |

Different values of the dipolar field strength and obliquity – 1.75 kG and 70° – compared to the dipolar model proposed by Landstreet & Borra (1977).

Surface distributions of several chemical elements on the surface of $\varphi$ Dra were studied by Kuschnig (1998) with the help of the Doppler imaging (DI) technique. Using these chemical spot maps, Prvák et al. (2015) calculated theoretically expected flux variations in different photometric bands and succeeded in reproducing the observed UV and optical photometric behaviour of this star. $\varphi$ Dra was included in the volume-limited survey of magnetic Ap stars by Sikora et al. (2019a). These authors derived atmospheric parameters, assessed evolutionary state, and measured mean abundances. They have concluded that $\varphi$ Dra is a 3.56 M$_\odot$ star that has completed 88 per cent of its life on the main sequence.

Motivated by the availability of new high-quality spectropolarimetric observations and unique long-duration, precise space-borne photometric data for $\varphi$ Dra, we have carried out a comprehensive study of the magnetic field topology, chemical abundance spots, photometric variability, and binarity of this star. Our paper is organized as follows. Section 2 presents observational data, followed by the description of analysis and results in Section 3, including a revision of the orbit of the inner spectroscopic binary (Section 3.1), time series analysis of the space photometric data (Section 3.2), assessment of the mean polarization profiles (Section 3.3), and Zeeman Doppler imaging (Section 3.4). This paper ends with conclusions and discussion in Section 4.

2 OBSERVATIONAL DATA

2.1 Spectropolarimetry

In this study, we analysed 22 high-resolution circular polarization observations of $\varphi$ Dra collected with the Narval spectropolarimeter installed at the 2 m Télescope Bernard Lyot (TBL) at the Pic du Midi observatory in France. The same data set was previously studied by Sikora et al. (2019a) and Sikora et al. (2019b). These observations provide a continuous coverage of the 3600-10000 Å wavelength interval at a resolving power of $R \approx 65000$. The spectra of $\varphi$ Dra were obtained between August 2016 and February 2017. Each observation consisted of four 360 s sub-exposures acquired with different polarimeter configurations, allowing one to derive Stokes V and diagnostic null spectra (Donati et al. 1997; Baguhl et al. 2009). The optimal extraction of 1D spectra and their polarimetric demodulation was carried out with an updated version of the ESPRIT reduction code (Donati et al. 1997), yielding data with a typical signal-to-noise ratio (S/N) of 900-1200 per 1.8 km s$^{-1}$ velocity bin at $\lambda = 5500$ Å. The spectra were post-processed with the continuum normalization routines described by Rosén et al. (2018). Further details on individual observations, including observing dates, heliocentric Julian dates, and S/N values, are given in the first three columns of Table 1.

2.2 Space photometry

The Transiting Exoplanet Survey Satellite (TESS, Ricker et al. 2015) was launched in April 2018 and is currently carrying out a survey of most of the sky in sectors measuring 24$^\circ \times 96^\circ$. Each sector is observed for 27.4 d, which corresponds to two orbits of the satellite. These observations yield continuous high-precision photometric data in a wide red (600–1000 nm) bandpass over $\approx 13$-d time span corresponding to each orbit with short gaps between the orbits and sectors. The sectors are oriented along ecliptic longitudes, resulting in repeated observations of targets in the continuous viewing zones (CVZ) around both ecliptic poles. $\varphi$ Dra is one of the few bright Ap stars located within the northern TESS CVZ. This star was therefore observed with a 2-min cadence over the entire second year of the
**3 ANALYSIS AND RESULTS**

### 3.1 Spectroscopic binary orbit

The orbital radial velocity variation of ϕ Dra has been known for a long time (e.g. Beardsley 1969; Abt & Snowden 1973). However, the period and amplitude of the radial velocity curve were poorly constrained (Liška 2016) due to the difficulty of extracting accurate velocity measurements from broad, variable spectral lines of the Ap star. This problem has been largely overcome by Bischoff et al. (2017), who provided 62 precise radial velocity measurements obtained from hydrogen Balmer lines. Since stellar atmospheres are predominantly made of hydrogen, the lines of this element are much less affected by surface chemical abundance inhomogeneities compared to metal lines (typically Mg, Fe, Cr, Si) accessible for radial velocity measurements. We have replicated the methodology used by Bischoff et al. (2017) by calculating the centre-of-gravity velocities from the cores of Hβ, Hγ, and Hδ (Hβ was avoided due to contamination by variable telluric lines). The resulting average H-line radial velocities are reported in Table 1. The scatter between different Balmer lines amounts to 0.6–0.8 km s\(^{-1}\), which is much smaller compared to 5–10 km s\(^{-1}\) rotational jitter exhibited by metal lines.

The 22 new radial velocity measurements were combined with the data from Bischoff et al. (2017) and the full set of SB1 spectroscopic orbital elements (\(P_{\text{orb}}\), HJD\(_0\), K, \(\gamma\), \(e\), \(\omega\)) was obtained by fitting observations with a non-linear least-squares optimization algorithm (Markwardt 2009). The results are presented in Table 2. The observed radial velocities are compared to the radial velocity curve corresponding to the best-fitting orbital solution in Fig. 1. The standard deviation of our H-line radial velocity measurements around the fitted curve is 0.9 km s\(^{-1}\), which is consistent with the uncertainty estimate above.

An important caveat of this orbital analysis is that it relies on the assumption that the H-line radial velocity variation is due to the orbital motion of the component Aa and that the secondary component of the inner binary, Ab, does not contribute to the spectra. But, as will be demonstrated below (Section 3.3), ϕ Dra is actually an SB2 system with the component Ab contributing to the metal line spectra and therefore to hydrogen lines as well. This means that the H-line radial velocity measurements yield a systematically underestimated orbital semi-amplitude. A corrected value of this parameter is derived in Section 3.4 as part of the Doppler imaging analysis.

### 3.2 Photometric variability

Time series analysis of the TESS observations of ϕ Dra was carried out using standard methods which combined calculation of periodograms and fitting data in the time domain (e.g. Kochukhov et al. 2021). The generalized Lomb-Scargle periodogram (Zechmeister & Kürster 2009) for the entire TESS data set is presented in Fig. 2 (upper panel). The rotational frequency and its multiple harmonics are clearly visible.

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1. https://mast.stsci.edu
Figure 2. Upper panel: periodogram for the TESS light curve of ψ Dra. Middle panel: periodogram for the residual time series after removing the rotational signal. Lower panel: TESS observations plotted as a function of the rotational phase (symbols) along with the harmonic fit (solid line).

Table 3. Photometric rotational period determinations for ψ Dra.

| Period (d) | Reference               |
|-----------|-------------------------|
| 1.71646(6) | Musielok et al. (1980)  |
| 1.71650(2) | Prvátko et al. (2015)   |
| 1.71665(9) | Bernhard et al. (2020)  |
| 1.7166(1)  | Paunzen et al. (2021)   |
| 1.71650213(21) | This work               |

In order to extract the full information about rotational modulation, we fitted the TESS light curve with a superposition of 10 harmonics of the rotational frequency. This periodic model was added to third-degree polynomial functions describing slow instrumental trend within each of the 24 continuous 10.1–13.5 d light-curve segments. All parameters of this composite model were optimized simultaneously with the help of a non-linear least-squares fitting algorithm. This analysis yields a stellar rotational period of 1.71650213(21) d, which agrees with all previous determinations within uncertainties (see Table 3) but is considerably more precise owing to a high quality and long baseline of the TESS photometric observations of ψ Dra.

Figure 3. Left-hand panel: phased TESS time series of ψ Dra averaged over 10–13 d light-curve segments. Right-hand panel: residual light curves for each segment.

Keeping the same reference Julian date as derived by Sikora et al. (2019b), we adopted the following rotational ephemeris in this work:

\[ \text{HJD} = 244263.30626 + 1.71650213(21) \times E. \]

The TESS data phased according to this ephemeris is compared to the model light curve in Fig. 2 (bottom panel). The photometric variation of ψ Dra in the TESS passband is dominated by a sinusoidal single-wave component, but the contribution of higher order harmonics is also significant.

The periodogram of the residual light curve obtained by subtracting the model of rotational variability from the observed data is shown in Fig. 2 (middle panel). There are clear residual signals associated with the rotational frequency and its harmonics. In particular, there are two frequency peaks separated by \( \pm 0.0080(15) \) d\(^{-1}\) from the rotational frequency. This frequency splitting is consistent with the orbital frequency 0.007818(4) d\(^{-1}\) derived in Section 1. This suggests that the photometric rotational variability of ψ Dra is modulated by the orbital motion within the Aab binary system. In addition, Fig. 2 reveals the presence of a signal with a period of 6.649(80) d, which likely corresponds to the rotation of either component Ab or B.

To gain further insight into long-term changes of the photometric rotational modulation of ψ Dra we constructed average binned light curves for each of the 24 continuous data segments. The time evolution of these average light curves is shown in Fig. 3 (left-hand panel). No significant changes on the time span of 1 yr covered by TESS photometry are immediately obvious. However, when we subtract the median light curve and examine evolution of the residuals (Fig. 3, right-hand panel), a clear repetitive S-shaped pattern emerges. This cyclic structure in the residuals is particularly clear for the first 200 d of TESS observations and appears to be consistent with \( P_{\text{orb}} \approx 130 \) d derived earlier.
variation shown by the dashed line in Fig. 4 (lower panel). Both the phase and shape of this curve are in good agreement with the experimental time-delay measurements, confirming that the light traveltime effect is the most likely explanation of the long-term cyclic changes of the photometric rotational variability of ϕ Dra. We also attempted to derive the orbital parameters \( K, e, \) and \( \omega \) directly from the photometric time-delay measurements. This approach yielded \( K = 42.1 \pm 6.7 \, \text{km s}^{-1}, e = 0.674 \pm 0.079; \) and \( \omega = 133.0^\circ \pm 3.4^\circ \). The resulting orbital semi-amplitude is significantly higher than the one inferred from hydrogen lines in Section 3.1, confirming that the latter radial velocity indicator is biased. On the other hand, the photometric radial velocity semi-amplitude is consistent within \( 1.2\sigma \) with the estimate obtained from modelling Fe lines in Section 3.4, but is considerably less precise compared to the latter.

### 3.3 Multiline profile analysis

#### 3.3.1 Model atmosphere and mean abundances

Two previous spectroscopic studies derived atmospheric parameters and analysed the chemical composition of ϕ Dra. Kuchnerg (1998) adopted \( T_{\text{eff}} = 12,500 \, \text{K}, \log g = 4.0, \) and obtained surface abundance distributions of He, Si, Ti, Cr, and Fe with the Doppler imaging method. This combination of atmospheric parameters and chemical abundance maps allowed Prváč et al. (2015) to reproduce photometric variability of ϕ Dra from UV to near-infrared. Sikora et al. (2019a) determined \( T_{\text{eff}} = 11630 \, \text{K} \) and \( \log g = 4.11 \) from spectroscopic and photometric observations but estimated \( \log g = 3.85 \) from the comparison with stellar evolutionary models. They also provided abundances of Si, Ti, Cr, and Fe. This abundance analysis was based on the same set of Narval observations as analysed in our study, but appears to have been carried out using a single high-quality observation since the authors do not mention correcting for the orbital radial velocity variation, which would have been necessary to obtain a phase-averaged spectrum.

Taking into account results of the previous studies, we adopted \( T_{\text{eff}} = 12,000 \, \text{K}, \log g = 4.0, \) and calculated a stellar model atmosphere with these parameters, using the LLMODELS code (Shulyak et al. 2014). These atmospheric parameters together with the abundances reported by Sikora et al. (2019a) were employed to obtain a preliminary line list from the VALD3 database (Ryabchikova et al. 2015). We constructed a mean spectrum of ϕ Dra by removing the orbital radial velocity offsets predicted by our spectroscopic orbital solution and averaging all 22 observations. The resulting mean spectrum was compared with the theoretical synthetic spectrum calculated with the SYNTH3 code (Kochukhov 2007) for the parameters listed above and \( v_\text{t} \sin i = 81.9 \, \text{km s}^{-1} \) (Sikora et al. 2019a). Abundances of several elements with a large number of absorption features were then adjusted to reduce the discrepancy between the average observed spectrum and calculations. This analysis yielded the following set of mean abundances: \( \log (N(\text{He})/N(\text{tot})) = -2.40, \) \( \log (N(\text{Si})/N(\text{tot})) = -3.52, \) \( \log (N(\text{Ti})/N(\text{tot})) = -6.90, \) \( \log (N(\text{Cr})/N(\text{tot})) = -5.58, \) and \( \log (N(\text{Fe})/N(\text{tot})) = -4.09 \) with an uncertainty of about 0.2 dex. A new atomic line list was extracted from VALD3 using this revised abundance table.

#### 3.3.2 Least-squares deconvolved profiles

Due to a large rotational line broadening and moderate magnetic field strength, circular polarization spectra of ϕ Dra do not readily show Zeeman signatures except in the few strongest lines. In this situation, it is convenient to apply the least-squares deconvolution...
For the absorption example, was due in well of the continuum, initially, spectra of many individual lines into high-quality mean line profiles. In this work, we used the iLSD code (Kochukhov et al. 2010) and the VALD3 line list tailored for ϕ Dra for the LSD profile calculation. Initially, we considered all metal lines deeper than 5 per cent of the continuum, excluding wavelength regions contaminated by telluric absorption and affected by broad wings of the hydrogen Balmer lines. In practice, the wavelength range was restricted to 4000–7850 Å. These selection criteria yielded 1260 metal lines and resulted in Stokes V LSD profiles with a typical uncertainty of 6.3 × 10⁻⁵ per 2 km s⁻¹ velocity bin. This set of Stokes V spectra and the corresponding Stokes I profiles are illustrated in Fig. 5. The LSD line mask normalization adopted for this calculation (see Kochukhov et al. 2010 for details) was λ₀ = 5000 Å and z₀ = 1.2.

As evident from Fig. 5, variable Stokes V signal with an amplitude of about ±0.05 per cent of the unpolarized continuum is detected well above the noise level in all observations. No signatures are seen in the LSD profiles computed from the diagnostic null spectra. The Stokes I profiles exhibit a strong rotational modulation, primarily due to spots of Fe and Cr that dominate (941 out of 1260 lines) the metal line mask. In addition to this prominent variability, we detected weak variable absorption outside the ±v, sin i velocity range, which was correlated with the orbital rather than rotational phase. For example, the group of four profiles with a conspicuous red-wing absorption component corresponds to the observations obtained in the 0.074–0.097 orbital phase interval when the primary (component Aa) had a negative velocity with respect to the centre of mass and the secondary (component Ab) had a positive velocity offset (see Fig. 1). Conversely, another group of four observations, obtained in the 0.601–0.624 orbital phase interval, shows depression on the blue side of the line when the secondary was predicted to have a negative velocity offset relative to the primary. Therefore, distortions seen in the outer wings of the Stokes I profiles are consistent with contribution of the secondary star. This indicates that ϕ Dra is an SB2 system. Unfortunately, our observations have a poor coverage of the 127.9-d binary orbit, precluding us from pursuing further binary-star analysis such as deriving radial velocities of the secondary and spectral disentangling.

The LSD profiles calculated with the complete metal line mask allow one to achieve the highest precision in detecting Zeeman signatures. However, such profiles are unsuitable for detailed quantitative modelling of magnetic field and starspot geometries because the line strength modulation influencing both I and V profiles is a result of compound effect of several chemical elements with different surface distributions. The modelling task can be accomplished only by considering LSD profiles of individual chemical elements. To this end, we derived three additional sets of LSD profiles for Fe, Si, and Cr. The corresponding LSD masks included 828, 103, and 113 lines, respectively, with the remaining metal lines treated as a background with the multiprofile LSD approach (Kochukhov et al. 2010). The Fe and Si profiles obtained using this method are modelled in Section 3.4 to derive both the magnetic field topology and abundance distributions of these elements. Chromium lines, being less numerous than those of Fe and, on average, weaker than those of Si, yield rather noisy Stokes V LSD profiles. For this reason, we use only the Stokes I profiles of this element for abundance spot mapping.

### 3.3.3 Longitudinal magnetic field

The mean longitudinal magnetic field, ⟨B₀⟩, was determined from the first moment of the Stokes V LSD profile normalized by the equivalent width of Stokes I (Wade et al. 2000; Kochukhov et al. 2010). Numerical integration required for the evaluation of line profile moments was carried out in the ±90 km s⁻¹ velocity interval after removing the orbital radial velocity variation. The resulting ⟨B₀⟩ measurements obtained from the LSD profiles derived using all metal lines are provided in Table 1. Fig. 6 illustrates these measurements as a function of rotational phase together with ⟨B₀⟩ measurements reported in earlier studies. According to more precise Narval measurements, ϕ Dra exhibits a smooth, approximately sinusoidal variability of the mean longitudinal magnetic field between about −290 and +440 G. Our measurements agree within uncertainties with the results derived by Sikora et al. (2019b) from the same
data, so we do not pursue further quantitative analysis of the \( (B_p) \) phase curve of \( \varphi \) Dra. At the same time, it may be noted that all investigations, including ours, are slightly underestimating the field strength of \( \varphi \) Dra because the spectral contribution of the secondary is not removed from Stokes \( I \) prior to \( (B_p) \) calculation.

### 3.4 Zeeman Doppler imaging

The magnetic field geometry and chemical abundance maps of \( \varphi \) Dra were reconstructed with the help of the Zeeman Doppler imaging (ZDI) code INVERS-LSD, introduced by Kochukhov et al. (2014) and subsequently used by Kochukhov et al. (2017), Kochukhov et al. (2019), Rosén, Kochukhov & Wade (2015), and Oksala et al. (2018). This tomographic mapping procedure allows one to interpret rotational modulation of the LSD Stokes parameter profiles of magnetic stars based on detailed polarized radiative transfer calculations in a realistic model atmosphere. This is accomplished by calculating a grid of local Stokes spectra, including all relevant absorption lines for the entire observed wavelength interval, and a range of magnetic field strengths, field vector inclinations, chemical abundances, and limb angles, and then applying the LSD procedure to these calculations using the same line mask and relative weights as in the treatment of observations. The resulting grid of theoretical local Stokes parameter profiles is tabulated once and then used by the inversion module, which interpolates within this local profile table during calculation of the disc-integrated spectra and their derivatives. This approach allows us to separate time-consuming detailed polarized spectrum synthesis, usually carried out at a supercomputing facility, from the actual ZDI inversions for different sets of nuisance parameters, which require relatively modest computational resources.

In this study, we reconstructed the magnetic field geometry from the Stokes \( I \) and \( V \) profiles of Fe and Si. Chemical abundance maps were derived simultaneously with magnetic mapping for both elements. In addition, we obtained a non-uniform distribution of Cr from its Stokes \( I \) LSD profiles. For each of the three chemical elements, relevant abundance ranges were covered with a step of 0.25 dex during the calculation of local LSD profiles. A step of 200 G was employed for the magnetic field strength, while both the angle between the field vector and the line of sight and limb angle parameter spaces were sampled with 15 values each. The same LLMODELS atmosphere with \( T_{\text{eff}} = 12\,000 \text{K} \) and \( g = 4.0 \), the mean abundances discussed above, and a complete VALD3 line list comprising about 5000 transitions in the 4000–7850 Å wavelength interval were adopted for the spectrum synthesis calculations.

The global magnetic field geometry of \( \varphi \) Dra was parametrized using a spherical harmonic expansion (Donati et al. 2006; Kochukhov et al. 2014), which included both poloidal (potential) and toroidal (non-potential) terms. Three independent sets of the spherical harmonic coefficients, \( \alpha \), \( \beta \), and \( \gamma \), were employed to describe the radial poloidal, horizontal poloidal, and horizontal toroidal field components, respectively. The expansion was truncated at the maximum angular degree, \( \ell = 10 \), which is sufficient to describe the global magnetic fields of Ap stars. A harmonic penalty function described by Morin et al. (2008) and Kochukhov et al. (2014) was applied to prevent the inversion procedure from introducing unnecessarily complex magnetic field distributions. Chemical abundance maps were prescribed using an 1876-element surface grid and regularized with the Tikhonov method (Piskunov & Kochukhov 2002). Both the abundance and magnetic field regularization parameters were adjusted following the procedure outlined by Kochukhov (2017) to achieve a balance between reaching a satisfactory fit to observations and avoiding spurious small-scale surface features.

### 3.4.1 Refinement of the stellar and orbital parameters

As discussed above, the blending of hydrogen lines of the Ap primary by the secondary star in the inner spectroscopic binary introduces a systematic bias in the radial velocity measurements, yielding an underestimated orbital velocity amplitude. Metal lines are intrinsically narrower, allowing one to better isolate the spectral signature of the primary and potentially obtain a more accurate orbital amplitude. However, metal lines exhibit strong rotational modulation due to chemical spots on the primary. Thus, in practice, a refinement of the orbital parameters using metal lines is only possible if one can reproduce Stokes \( I \) profile shapes in every observation. This necessitates modelling of abundance spots with a DI code. To this end, we carried out chemical spot inversions using the Fe Stokes \( I \) LSD profiles for a two-dimensional parameter grid spanning a range of \( K \) and \( y \) values while keeping all other orbital elements fixed. A well-defined \( y \) minimum was found at \( K = 33.87 \pm 0.25 \text{ km s}^{-1} \) and \( y = -16.44 \pm 0.08 \text{ km s}^{-1} \). This centre-of-mass velocity differs little from the orbital solution based on the hydrogen-line radial velocity measurements. On the other hand, the orbital semi-amplitude found using Fe lines is 26 per cent higher than the one inferred from H lines. This orbital parameter derived from Fe lines is more accurate than the orbital amplitude obtained from H lines since spectroscopic contribution of the primary can be better isolated using intrinsically narrower metal lines. Consequently, we adopt the orbital parameters determined in this section for all subsequent ZDI calculations.

Aiming to improve the \( v_\text{r}\sin i \) estimate, we carried out ZDI inversions for a range of projected rotational velocities and compared the resulting fit quality. This analysis was performed for both Si and Fe LSD profiles, yielding \( v_\text{r}\sin i = 82.0 \pm 0.5 \) and \( 81.0 \pm 0.5 \text{ km s}^{-1} \) for these two elements, respectively. We adopted \( v_\text{r}\sin i = 81.5 \pm 0.5 \text{ km s}^{-1} \) as a compromise value. This determination of the projected rotational velocity of \( \varphi \) Dra is consistent with \( v_\text{r}\sin i = 81.9 \pm 1.6 \text{ km s}^{-1} \) found by Sikora et al. (2019a) from the conventional spectrum synthesis assuming uniform chemical abundance distributions.

### 3.4.2 Magnetic field topology

Results of the tomographic reconstruction of the magnetic field structure of \( \varphi \) Dra using the Si and Fe LSD profiles are illustrated in Figs 7–11. The first of these figures compares the observed and model Stokes profiles. Figs 8 and 9 present spherical maps of the field strength, the radial field component, and the field vector orientation obtained from modelling Si and Fe lines, respectively. In both cases, we derived a distorted dipolar topology characterized by a large obliquity relative to the stellar rotational axis and a significant asymmetry between the positive and negative magnetic poles. The former is represented by an arc-like structure in which the field strength reaches 1.32–1.48 kG. The latter is much more diffuse, with a field strength below 1 kG.

Fig. 10 provides a closer look at the inter-agreement of the magnetic field maps derived from the mean profiles of the two chemical elements. This figure shows that the results of independent magnetic inversions are broadly consistent with each other despite a noticeably different pattern of the Stokes \( I \) and \( V \) profile variability.
Field of spherical and magnetic is applied as discrepant (8.4). Likewise, corresponding radial, meridional, and azimuthal field components of the two maps are consistent to within 113, 112, and 104 G, respectively. This corresponds to 7.4–8.1 per cent of the peak surface field strength. Likewise, the field modulus distributions agree to within 118 G (8.4 per cent of the peak field) and the local field inclination is discrepant by 11.6° on average. These figures can be considered as an estimate of the internal precision of the magnetic inversions applied in this study.

The harmonic content of the reconstructed magnetic field maps is assessed in Fig. 11. We find that most (75–77 per cent) of the magnetic field energy is concentrated in the dipolar component and that the field structure is predominately poloidal (all toroidal spherical harmonic modes combined contribute only 3.5–5.6 per cent of the total field energy) and does not contain a significant small-scale field (all modes with \( \ell > 5 \) combined contribute less than 2 per cent of the field energy).

To conclude the ZDI analysis, we assessed the feasibility of reproducing spectropolarimetric observations of \( \varphi \) Dra with a canonical oblique dipolar topology (e.g. Sikora et al. 2019b) or with a superposition of dipolar and general non-axisymmetric quadrupolar field geometries (Baguñol et al. 2002). In these tests, conducted using the Fe Stokes / and V LSD profiles, the magnetic field geometry was reconstructed by setting \( \ell_{\text{max}} = 1 \) or 2 and adopting \( \beta_{lm} = \alpha_{lm} = \gamma_{lm} = 0 \) spherical harmonic expansion (see Kochukhov et al. 2014 and Kochukhov & Wade 2016 for detailed discussions of different harmonic parametrizations). No magnetic field regularization was applied and the Fe abundance distribution was reconstructed simultaneously with the magnetic field map. With these restrictions, INVERSLSD was not able to obtain a satisfactory fit to the Fe Stokes V profiles, with the final standard deviation remaining 55–68 per cent higher than for the inversion based on the general high-\( \ell \) harmonic parametrization normally adopted in ZDI. This confirms that the global distortion of the dipolar field as well as higher-order harmonic perturbations are essential for reproducing.

Figure 7. Comparison of the observed (symbols) and model (solid lines) Stokes IV Si, Fe, and Cr LSD profiles. The spectra are offset vertically, with the rotational phase indicated next to each profile. The Stokes V spectra are magnified by a factor of 12 relative to Stokes I.
3.4.3 Chemical spot distributions

Surface abundance distributions of Si and Fe were reconstructed simultaneously with the magnetic field maps from the Stokes $I$ and $V$ profiles of these elements. The resulting chemical spot maps are presented in Fig. 12. This figure also shows the surface distribution of Cr recovered from the mean intensity profiles of this element. The corresponding observed and computed Stokes $I$ spectra are shown in the right-most column of Fig. 7. This DI reconstruction of the Cr spot geometry was carried out using the magnetic field map obtained by averaging the field distributions inferred in the ZDI analyses of Si and Fe.

Figure 8. Magnetic field topology of $\varphi$ Dra derived from the Si LSD profiles. The spherical plots show maps of the total magnetic field strength (a), the radial field component (b), and the field vector orientation (c) at five rotational phases, which are indicated next to each column. The contours over spherical maps in panels (a) and (b) are plotted with a step of 0.2 kG. The stellar equator and rotational pole are indicated with double lines. The side colour bars give the field strength in kG.

Figure 9. Same as Fig. 8 for the magnetic field topology derived from the Fe LSD profiles.
Significant surface inhomogeneities are found for all three chemical elements. Si shows the smallest (∼1 dex) variation across the northern stellar hemisphere. Fe and Cr exhibit stronger inhomogeneities, with the abundance contrasts exceeding 1.5 dex. The two studied Fe-peak elements possess similar surface distributions over a significant fraction of the stellar surface, but Cr shows larger and higher contrast overabundance spots compared to Fe. At the same time, Fe also exhibits a more complex surface distribution with an additional series of spots (phases 0.6–0.8) not visible in the Cr map.

For all three chemical elements, inversions yield major underabundance or overabundance areas below the stellar rotational equator. This part of the stellar surface is poorly sampled by observations owing to the moderate inclination of ϕ Dra. In this situation, abundance mapping in the sub-equatorial regions is more prone to artefacts, including those possibly related to unaccounted blending by the secondary component. Using the projected surface area summed over all rotational phases as a proxy for relative visibility, one can conclude that the latitudes below −11.8° contribute less than 25 per cent to the disc-integrated profiles compared to the high-latitude zones, which have the best visibility. Therefore, the prominent stellar abundance features reconstructed in this study below the rotational equator are less reliable than the structures at higher latitudes.

4 CONCLUSIONS AND DISCUSSION

This study presented a comprehensive investigation of the rotational variability, binarity, magnetism, and surface structure of the bright Ap star ϕ Dra. We significantly improved the precision of the stellar rotational period, finding $P_{\text{rot}} = 1.71650213(21)$ d, based on the photometric observations collected for this star by the TESS satellite over a period of nearly 1 yr. An in-depth analysis of the rotational photometric phase curve revealed its modulation with the 127.9-d period corresponding to the motion of the Ap star in a spectroscopic binary with a lower mass, fainter companion. We successfully interpreted this long-term light curve modulation as the light traveltime effect (Shibahashi & Kurtz 2012; Murphy et al. 2014) – the first time that this phenomenon has been observed to affect the stellar rotational variability.

Using a set of high-resolution spectra, we obtained new radial velocity measurements for ϕ Dra and revised parameters of its spectroscopic binary orbit, taking into account all available high-precision radial velocity determinations. We also reported the first detection of the spectroscopic contribution of the secondary in metal line profiles, thereby changing the status of ϕ Dra from an SB1 to SB2 system.
We derived mean intensity and polarization profiles of Fe, Cr, and Si with the help of the LSD technique. This multiline analysis is a necessary prerequisite to studying the magnetic field topology of \( \varphi \) Dra since its relatively weak magnetic field and rapid rotation do not allow one to use individual lines for magnetic diagnostic. We were able to derive mean circular polarization signatures with a quality sufficient for detailed modelling from the lines of Fe and Si. A magnetic tomographic mapping technique was then applied to these observations in order to obtain vector magnetic field maps simultaneously with chemical spot distributions. This analysis yielded two independent magnetic field distributions, which agreed at the level of \( \sim 10 \) per cent of the peak field strength. The surface field topology of \( \varphi \) Dra was found to be primarily poloidal and dipolar, with small but statistically significant contributions of higher-order poloidal harmonic components. This magnetic field topology is characterized by a maximum local field modulus of 1.39 kG and a surface-averaged field strength of 0.54 kG, making it one of the lowest-field Ap stars studied with ZDI. In general, the surface magnetic field properties of \( \varphi \) Dra are unremarkable compared to other Ap stars studied with magnetic mapping (see Kochukhov et al. 2019, and references therein).

We reconstructed chemical spot distributions of Si and Fe simultaneously with mapping magnetic field using the Stokes I and V LSD profiles of these elements. In addition, the distribution of Cr was recovered from the Stokes I LSD profile of this element keeping the magnetic field map fixed. This set of three chemical maps is small compared to multi-element distributions available for some Ap stars (e.g. Kochukhov et al. 2004; Lüftinger et al. 2010; Nesvacil et al. 2012; Silvester, Kochukhov & Wade 2014). However, an advantage of this study is that a detailed map of magnetic field is recovered.

Figure 11. Relative energies of the poloidal and toroidal harmonic components as a function of the angular degree \( t \) for the magnetic field topologies reconstructed from the Si (top panel) and Fe (bottom panel) LSD profiles.

Figure 12. Surface distributions of Si, Fe, and Cr derived from the LSD profiles of these elements. The fractional element abundances are indicated by the side bars in the \( \log (N_i/N_\text{tot}) \) units, with the minimum and maximum values corresponding to 5 and 95 percentiles, respectively. The contours over spherical maps are plotted every 0.2 dex.
Figure 13. Comparison of the magnetic field topology of ϕ Dra with the surface abundance distributions of Si, Cr, and Fe. The radial field component, the field modulus, and the local field inclination for the average magnetic field map are illustrated from top panel to bottom panel, similar to Fig. 10. Chemical abundance spots are shown with contours plotted every 0.25 dex. The darker and thicker contours correspond to zones with a higher element abundance. Dashed lines trace two small circles on which overabundance spots are found.

self-consistently from the same observational material and that this magnetic field is relatively weak, simplifying both the reconstruction and theoretical interpretation of the chemical abundance maps. In addition, ϕ Dra is one of the fastest rotating Ap stars to which modern ZDI analysis was applied, yielding higher resolution surface maps compared to DI results typically found in the literature (see summary in Kochukhov 2017).

Spherical maps of ϕ Dra (Figs 8, 9, and 12), which are often analysed in the discussion of ZDI and DI results for Ap stars, do not readily reveal a relation between the magnetic field topology and chemical abundance distributions. To investigate this relation in more detail, we plotted abundance maps over the magnetic field geometry using the Hammer–Aitoff map projection (Fig. 13). It is evident that spot maps are morphologically more complex than the magnetic field structure and do not exhibit a simple correlation with the local field inclination as predicted by many theoretical atomic diffusion models (e.g. LeBlanc et al. 2009; Alecian & Stift 2010; Alecian 2015; Kochukhov & Ryabchikova 2018). Nevertheless, one can note that one of the abundance spots in all three studied chemical maps coincides with the same strong-field region where the field vector is directed outwards.

Despite the apparent lack of a simple abundance-magnetic field correlation, we recognized an intriguing regularity of the surface locations of chemical spots relative to the stellar rotational and magnetic field topologies. The spots are found along the two small circles, which have opening angles of ≈ 45°, are located ≈ 180° apart and centred at a latitude of ≈ ±10° at the magnetic equator. These small circles, highlighted by the dashed lines in Fig. 13, appear to account for all the main overabundance features of the three studied elements, including spots below the rotational equator, which we previously deemed to be less reliable. There is a tendency for the overabundance spots to be found along these circles in one of the four quadrants defined by the intersection of the rotational and magnetic equators. At the same time, spots typically avoid the magnetic equator (except for one high-latitude feature in the Fe map) and populate these quadrants differently depending on the element. Nevertheless, a systematic behaviour of spot locations is clearly present, hinting at the importance of stellar rotation alongside the magnetic field for atomic diffusion. Further analysis of the DI results obtained for other Ap stars with the graphical representation similar to Fig. 13 is necessary to assess if the star-spot behaviour uncovered in this study of ϕ Dra is a unique property of this particular star or can be found in other objects.

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DATA AVAILABILITY
The spectropolarimetric data underlying this article are available from the PolarBase archive (http://polarbase.irap.omp.eu). The TESS observations of ϕ Dra can be obtained from the Mikulski Archive for Space Telescopes (MAST, https://mast.stsci.edu).

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