

*Poster paper*

**INVESTIGATION OF ROTATIONAL VELOCITY  
OF E-PERSEI (EPSILON-PERSEI)**

N. GAVRILOVIĆ<sup>1</sup>, S. JANKOV<sup>1</sup>, P. MATHIAS<sup>2</sup> and P. De CAT<sup>3</sup>

<sup>1</sup>*Astronomical Observatory, Volgina 7, 11160 Belgrade, Serbia*

<sup>2</sup>*Observatoire de la Cote d'Azur, Département Fresnel, U.M.R. 6528, B.P. 4229,  
06304 Nice Cedex 04, France*

<sup>3</sup>*Instituut voor Sterrenkunde, Katholieke Universiteit Leuven,  
Celestijnenlaan 200 B, 3001 Heverlee, Belgium*

*E-mail: ngavrilovic@aob.bg.ac.yu*

We present the analysis of spectral line profiles of the Si III triplet at 455,3 nm, 456,8 nm and 457,4 nm of a variable star  $\epsilon$ -Persei, and we investigate the  $v\sin(i)$  value of the star using Fourier transform technique. Since the star is a strong non-radial pulsator the spectra averaged over several pulsational cycles have been used.

The derived average value using all lines is  $v\sin(i)=134$  km/s.

*Poster paper*

**ELECTRIC DIPOLE TRANSITION PROBABILITIES  
IN Al IV AND Al V IONS**

RAFIK HAMDI and NÉBIL BEN NESSIB

*Groupe de Recherche en Physique Atomique et Astrophysique,  
Faculté des Sciences de Bizerte, 7021 Zarzouna, Tunisia*

*E-mail: hamdi\_rafik12003@yahoo.fr*

Electric dipole transition probabilities in triply and four times ionized aluminium have been calculated in intermediate coupling.

The present calculations were carried out with the general purpose atomic-structure program SUPERSTRUCTURE (Eissner et al.1974), as modified by Nussbaumer and Storey (1978). The wavefunctions are of the type  $\psi = \sum_i \phi_i C_i$ , where the basis functions  $\phi_i$  are constructed using one-electron orbitals  $\psi$ . The latter are calculated with a scaled Thomas-Fermi statistical model potential (Eissner and Nussbaumer 1969) or obtained from the Coulomb potential (Nussbaumer and Storey 1978).

The relativistic corrections to the non-relativistic Hamiltonian are taken into account through the Breit-Pauli approximation.

We have also introduced a semi-empirical correction (TEC) for the calculation of the energy-levels

The adopted atomic model for Al IV includes 12 configurations corresponding to 103 fine structure levels. For Al V the model includes 25 configurations corresponding to 434 fine structure levels.