

AB INITIO DETERMINATION OF ATOMIC STRUCTURE AND STARK BROADENING PARAMETERS: Pb IV AND RECENT RESULTS

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In this work, we present a review of our previous ab initio calculations of Stark broadening parameters. Those results are determined using semi-classical perturbation method for Stark width and shift calculations and SUPERSTRUCTURE (SST) code for determination of atomic structure. SST code takes into account configuration interaction and relativistic effects are introduced as perturbation of the non-relativistic Hamiltonian.

Triply ionized lead (Pb IV) is characterized by a strong resonance lines and it is a candidate for spectroscopic detection in hot DA white dwarfs. Pb IV resonance lines are also detected in photospheres of subdwarfs stars.

New results are also presented for some spectral lines of Pb IV of the type $5d^{10}nl$ - $5d^{10}n'l'$. Energy levels and oscillator strengths are calculated using Hartree-Fock Relativistic approach (HFR) and Stark broadening parameters are determined using semi-classical perturbation approach. Stark widths and shifts are presented as a function of temperature for perturber density of 10^{17} cm^{-3} . Such data are important for laboratory and technological plasmas and also for modelling and spectral diagnostics of stellar atmospheres.