## THE MODIFIED SEMI-EMPIRICAL STARK BROADENING METHOD OF CALCULATIONS: THE EXAMPLE OF ALKALI LIKE IONS

A. ALMODLEJ<sup>1</sup>,N. ALWADIE<sup>1,2</sup>, N. BEN NESSIB<sup>1,3</sup> and M. S. DIMITRIJEVIĆ<sup>4,5</sup>

<sup>1</sup>Department of Physics and Astronomy, College of Sciences, King Saud University, Saudi Arabia E-mail nbennessib@ksu.edu.sa

<sup>2</sup>Department of Physics, College of Sciences, King Khalid University, Saudi Arabia E-mail nalwadie@ksu.edu.sa

<sup>3</sup>GRePAA, INSAT, Centre Urbain Nord, University of Carthage, Tunis, Tunisia

<sup>4</sup>Astronomical Observatory, Volgina 7, 11060 Belgrade 38, Serbia E-mail mdimitrijevic@aob.rs

<sup>5</sup>Sorbonne Université, Observatoire de Paris, Université PSL, CNRS, LERMA, F-92190 Meudon, France

**Abstract.** In this work, the modified semi-empirical (MSE) Stark broadening method of calculations is presented. The method is compared to the semi-empirical (SE), semiclassical perturbation (SCP) and fully relativistic methods and with more simplest calculation formula. As an example of calculation, we calculate the widths of some alkali like ions as Sr II, Y III and Zr IV.

## References

- Cowley, C.R.: An approximate Stark broadening formula for use in spectrum synthesis, 1971, *The Observatory*, **91**, 139.
- Dimitrijević, M. S. and Konjević, N.: Stark widths of doubly-and triply-ionized atom lines, 1980, Journal of Quantitative Spectroscopy and Radiative Transfer, 24, 451.
- Elabidi, H., Ben Nessib, N. and Sahal-Bréchot, S.: Quantum mechanical calculations of the electron-impact broadening of spectral lines for intermediate coupling, 2003, Journal of Physics B: Atomic, Molecular and Optical Physics, 37, 63.
- Griem, H.R.: Semiempirical Formulas for the Electron-Impact Widths and Shifts of Isolated Ion Lines in Plasmas, 1968, Phys. Rev., 165, 258.
- Sahal-Bréchot, S., Dimitrijević, M. S. and Ben Nessib, N.: Widths and shifts of isolated lines of neutral and ionized atoms perturbed by collisions with electrons and ions: An outline of the semiclassical perturbation (SCP) method and of the approximations used for the calculations, 2014, Atoms, 2, 225.