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Ba⁺ vs. Cs: prospects for parity-nonconservation studies M.S. SAFRONOVA, E. ISKRENOVA-TCHOUKOVA, University of Delaware — We conduct a systematic study of Ba⁺ atomic properties using a relativistic all-order method. Energy levels, transition matrix elements, lifetimes, hyperfine constants, and polarizabilities are calculated. This work is motivated by the possibility to study the parity nonconservation (PNC) with a single trapped ion. Comparison of the experimental weak charge of an atom Q_W (a quantity which depends on input from atomic theory) with predictions provides important constraints on possible extensions of the standard model. The experimental value of Q_W is known to 0.5% for Cs and 3% for Tl, the uncertainty in both cases being dominated by atomic theory. Since the Cs PNC studies provided the most accurate determination of Q_W of such kind, we conduct a detailed comparison of accuracy of our Ba⁺ calculations with the corresponding calculations in Cs. We also calculate the ratio of the light shift in the 6s and $5d_{5/2}$ states in Ba⁺ at various wavelengths and consider the prospects of the precise determination of the Ba⁺ electric-dipole matrix elements from the combination of the theoretical and experimental [1] studies of the light shift ratios. [1] J. A. Sherman, T. W. Koerber, A. Markhotok, W. Nagourney, and E. N. Fortson, Phys. Rev. Lett. 94, 243001 (2005)

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