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Structure and properties of Al^+ KYLE ROLLIN, MICHAEL BROMLEY¹, Department of Physics, San Diego State University, San Diego, CA, JIM MITROY, School of Engineering, Charles Darwin University, Darwin, NT, Australia — The properties of a number of states of single charged Al^+ ion are determined from a large basis configuration interaction calculation. The main focus is on the polarizabilities of the low-lying states (the $3s^2$ 1Se , 3s3d $^{1,3}De$, 3s3p $^{1,3}Po$ and 3s4s $^{1,3}Se$ states) and the dispersion interactions of those states with the Al^+ ground state, the hydrogen atom and the rare gases.

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