

Reanalysis of the odd configuration system of scandium atom

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We report the fine- and hyperfine structure analysis of the odd configurations system of scandium atom in a large multiconfiguration basis. The configurations selected for the basis have to satisfy the following conditions: they should contribute to eigenvector composition of the levels with experimentally determined hyperfine structure constants and they should extend in the energy region up to 65000 cm^{-1} . We have tried to take into account first- and second-order effects in the fine- and hyperfine structure as comprehensively as possible with all experimental energy level and hyperfine structure constants available at present [1-7].

The good agreement between calculated and experimental energy levels and hyperfine structure constants was achieved. For all the levels belonging to the selected configuration basis the predicted hfs A constants have been calculated.

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