

THEORETICAL INVESTIGATION OF ENERGY LEVELS FOR BA VI

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This ion was selected for the investigation as it is of great importance to Astrophysics. It is evident that the lines can be observed in hot white dwarf stars [1]. In this work energy levels for Ba VI will be presented. The calculations for Ba VI are performed using general-purpose atomic structure package GRASP2018, based on multiconfiguration Dirac-Hartree-Fock and relativistic configuration interaction (RCI) methods.

As this ion is of Sb isoelectronic sequence, it was calculated using the same principle like in [2]. Active space (AS) method was used to compute radial wave functions layer by layer for four sets of virtual orbitals (AS₁₋₄). Single and double substitutions were done from multi reference set of orbitals up to 11s, 11p, 10d, 8f, 8g, 8h. Substitutions from [Kr]4d¹⁰ core shells were forbidden. After calculating AS₄, RCI method was used to further increase the accuracy of Ba VI energy level calculations. In total, 231 energy levels belonging to the configurations 5s²5p4f², 5s²5p²4f, 5s5p³4f, 5p⁵, 5s²5p²{5d, 6p, 6s}, 5s²5p³, 5s²5p5d², 5s5p³5d, 5s5p³6s and 5s5p⁴ were computed.

50 levels were selected based on the data availability at National Institute of Standards and Technology Atomic Spectra database (NIST) [3] for comparison. The relative difference to NIST data is displayed in Figure 1. The biggest disagreement is observed in 5s²5p²5d and 5s5p⁴ configuration levels, as the energies appear to be reversed in our calculations. The average relative difference compared to NIST [3] for AS₄ is 1.2%. In order to increase the accuracy of energy levels, core, core-valence, core-core correlations need to be investigated. During the conference further data from the calculations will be presented.

Conf.	^M L	J	E _{NIST} cm ⁻¹	AS ₃ %	AS ₄ %	Conf.	^M L	J	E _{NIST} cm ⁻¹	AS ₃ %	AS ₄ %
5s ² 5p ³	⁴ S	3/2	0.0	0.0	0.0	5s ² 5p ² (¹ D)5d	² G	7/2	212208.6	2.6	2.6
5s ² 5p ³	² D	3/2	17260.6	1.2	1.1	5s ² 5p ² (¹ D)5d	² G	9/2	222540	-0.3	-0.3
5s ² 5p ³	² D	5/2	23547.2	1.0	0.9	5s ² 5p ² (³ P)5d	⁴ P	5/2	214870.8	2.3	2.3
5s ² 5p ³	² P	1/2	36155.7	2.2	2.1	5s ² 5p ² (³ P)5d	⁴ P	3/2	217346.5	2.2	2.2
5s ² 5p ³	² P	3/2	49621	0.4	0.4	5s ² 5p ² (³ P)5d	⁴ P	1/2	218693.8	2.2	2.2
5s5p ⁴ (³ P)	⁴ P	5/2	128436.3	-2.2	-2.1	5s ² 5p ² (³ P)5d	² D	3/2	224093.2	0.9	0.8
5s5p ⁴ (³ P)	⁴ P	3/2	139920.1	-2.1	-2.1	5s ² 5p ² (³ P)5d	² D	5/2	228330.9	2.1	2.2
5s5p ⁴ (³ P)	⁴ P	1/2	142851.9	-2.0	-2.0	5s ² 5p ² (¹ D)5d	² P	1/2	234313.6	0.4	0.4
5s5p ⁴ (¹ D)	² D	3/2	158158.4	-0.7	-0.7	5s ² 5p ² (¹ D)5d	² P	3/2	246496.2	2.1	2.1
5s5p ⁴ (¹ D)	² D	5/2	163566.3	-0.8	-0.8	5s ² 5p ² (¹ D)5d	² D	3/2	235818.7	2.3	2.3
5s ² 5p ² (³ P)5d	² P	3/2	176498	-0.3	-0.3	5s ² 5p ² (¹ D)5d	² D	5/2	236640.5	3.1	3.0
5s ² 5p ² (³ P)5d	² P	1/2	229293.1	-0.6	0.6	5s ² 5p ² (³ P)5d	² F	7/2	242157.2	2.3	2.3
5s5p ⁴ (³ P)	² P	1/2	179076.6	4.8	4.8	5s ² 5p ² (³ P)5d	² F	5/2	257298.9	2.2	2.1
5s5p ⁴ (³ P)	² P	3/2	213042.9	3.4	3.4	5s ² 5p ² (¹ S)5d	² D	5/2	243103.4	-0.5	-0.5
5s ² 5p ² (³ P)5d	⁴ F	3/2	180293.1	-0.2	-0.2	5s ² 5p ² (¹ S)5d	² D	3/2	256774.8	1.8	1.8
5s ² 5p ² (³ P)5d	⁴ F	5/2	183552.2	-0.5	-0.5	5s ² 5p ² (³ P)6s	⁴ P	1/2	247725.6	-0.1	-0.1
5s ² 5p ² (³ P)5d	⁴ F	7/2	192561.2	-0.6	-0.6	5s ² 5p ² (³ P)6s	⁴ P	3/2	261563.8	-0.3	-0.3
5s ² 5p ² (³ P)5d	⁴ F	9/2	199080	-0.6	-0.6	5s ² 5p ² (³ P)6s	⁴ P	5/2	267649.7	-0.2	-0.2
5s ² 5p ² (¹ D)5d	² F	5/2	191382.2	-0.6	-0.6	5s ² 5p ² (¹ D)5d	² S	1/2	251008.8	2.6	2.6
5s ² 5p ² (³ P)5d	⁴ D	1/2	193882.1	-0.4	-0.4	5s ² 5p ² (³ P)6s	² P	1/2	265653.5	-0.2	-0.2
5s ² 5p ² (³ P)5d	⁴ D	7/2	196130.2	-0.5	-0.5	5s ² 5p ² (³ P)6s	² P	3/2	271083.2	-0.1	-0.1
5s ² 5p ² (³ P)5d	⁴ D	3/2	197255.4	-0.2	-0.2	5s ² 5p ² (¹ D)6s	² D	5/2	288565.1	-0.2	-0.2
5s ² 5p ² (³ P)5d	⁴ D	5/2	201857	-0.3	-0.3	5s ² 5p ² (¹ D)6s	² D	3/2	290903.6	-0.2	-0.2
5s5p ⁴ (¹ S)	² S	1/2	200937.6	0.0	0.0	5s ² 5p ² (¹ S)6s	² S	1/2	307387	0.1	0.1

Fig. 1. Comparison of energy levels computed in active spaces AS₃ and AS₄ with NIST [3] ASD recommended values for Ba VI.

- [1] Rauch, T. Stellar laboratories III. New Ba V, Va VI, and Ba VII oscillator strengths and the barium abundance in the hot white dwarfs G191-B2B and RE 0503-289, *Astronomy and Astrophysics* 566, A10, 6 (2014)
- [2] Radžiūtė, L., Gaigalas, G. Theoretical investigation of Sb-like sequence: Sb I, Te II, I III, Xe IV, and Cs V, *Atomic Data and Nuclear Data Tables*, Volume 152, 101585 (2023)
- [3] A. Kramida, Yu. Ralchenko, J. Reader, and NIST ASD Team, NIST Atomic Spectra Database (ver. 5.10), [Online]. Available: <https://physics.nist.gov/asd> [2023, October 18]. National Institute of Standards and Technology, Gaithersburg, MD. (2022)