

Quantum-chemical modeling of the interaction of 1,2-di-phenylcyclopropyl with *N*-benzylideneaniline

© Victoria V. Esenbaeva,¹ Alexander N. Vasyanin,²
Sergey N. Shurov,²⁺ and Lidia P. Yunnikova^{1*+}

¹ Perm State Agricultural Academy n.a. Academician D.N. Pryanishnikov. Petropavlovskaya St., 23.
Perm, 614000. Perm Region. Russia. Phone: +7 (342) 212-95-68. E-mail: yunnikova@yahoo.com.

² Perm State National Research University. Bukerev St., 15. Perm, 614990. Perm Region. Russia.
Phone: +7 (342) 239-64-35. E-mail: seshurov@yandex.ru

*Supervising author; +Corresponding author

Keywords: *N*-benzylideneaniline, 1,2-diphenylcyclopropyl, 1*a*,2,7*b*-triphenyl-1*a*,2,3,7*b*-tetrahydro-1*H*-cyclopropa[*c*]quinoline, quantum-chemical calculations.

Abstract

Mechanism has been suggested for the formation of 1*a*,2,7*b*-triphenyl-1*a*,2,3,7*b*-tetrahydro-1*H*-cyclopropa[*c*]quinoline in the reaction of 1,2-diphenylcyclopropyl with *N*-benzylideneaniline in the presence of zinc chloride.