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Prediction of relative density of furoxanes with the use of model «*ANSAB*»

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Abstract

In the work feasibility of using the earlier proposed model «*ANSAB*» for prognosticating the relative density of furoxanes has been shown. Satisfactory agreement of design and experimental values has been obtained. There has been estimated the molecular volume of substances, when the volume of atomic formations, composing molecules, did not have spherical symmetry.