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Me-Me'-X系三元化合物形成规律的人工神经网络

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摘要: 用 r_{Me} , $r_{Me'}$, r_X , x_{Me} 和 $x_{Me'}$ 等原子参数作为人工神经网络的输入, 203个可信的Me-Me'-X体系的三元化合物的形成情况作为输出, 研究了Me-Me'-X系三元化合物的形成规律, 这里 r 是离子半径, x 是电负性, Me代表一价金属, Me'代表二价金属, X代表卤族元素。利用所得规律预报了M-Eu-I (M=Li, Na, K, Rb和Cs)系三元碘化物的形成情况, 应用差热分析和粉末X射线法测定了它们的相图, 预报结果和实验测定结果的对比是令人满意的。

关键字: 三元化合物 相图 计算机预报 人工神经网络

ARTIFICIAL NEURAL NETWORK APPLIED TO REGULARITY OF FORMATION OF TERNARY COMPOUNDS IN Me-Me'-X SYSTEMS

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Abstract: Using the atomic parameters of r_{Me} , $r_{Me'}$, r_X , x_{Me} and $x_{Me'}$ as the inputs, the regularity of formation of ternary complex halides of Me-Me'-X halides systems has been investigated by artificial neural networks (ANNs). Where, r is the radius of the ion, x is the electronegativity of element, Me is the mono-valent metal, Me' is di-valent metal, X represents the F, Cl, Br or I. The regularity was found by training the ANNs with 203 known samples (such as Ag-Ca-Cl system and K-Mg-Cl system etc.). The formation of ternary complex iodides in M-Eu-I systems (where M represents Li, Na, K, Rb or Cs) was predicted by this trained ANNs. The predicted results are completely in agreement with the experimental facts.

Key words: ternary compound phase diagram computer prediction artificial neural network

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