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Modeling of the specific electric conductivity of melts of halogenides RbHal (Hal – F, Cl, Br, I, At)

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Abstract

The presented research is devoted to analytical description and modeling of the specific electrical conductivity for the individual rubidium halogenides melts RbHal (Hal – F, Cl, Br, I). Approximate mathematical equations were obtained during the analytical description using the Table Curve 2D, Table Curve 3D software packages. These equations describe the dependence of the specific electrical conductivity κ of the individual rubidium halogenides melts RbHal (Hal – F, Cl, Br, I) with the halogen order number $\kappa = f(Z)$, with the halide-ion ionic radius $\kappa = f(r)$, with ionic potential $\text{Hal}^- \kappa = f(1/r)$. The relationship of $\kappa(\text{RbHal})$ with such physicochemical properties as viscosity and density is considered. Using M.Kh. Karapetyants comparative methods, analytical dependences were obtained and graphical dependencies were plotted in the temperature range $(T_m + n)$ ($n = 5, 10, 50, 75, 100, 150, 200^\circ$ higher the halogenides melting temperature). These dependencies made it possible to calculate and to forecast the RbAt specific electrical conductivity, the numerical values of κ for which are absent in the reference literature. A comparative analysis of the obtained numerical value $\kappa(\text{RbAt})$ with the specific electrical conductivity of lithium, sodium and potassium astatides, calculated earlier. Analytical description and modeling of specific electrical conductivity of RbAt based on the relationship $\kappa(\text{RbAt})$ with viscosity (reduced viscosity) and density (reduced density) of rubidium halide melts were carried out in the coordinates "property – property" and "property 1 – property 2, property 3". Recommendations are given on the choice of an equation for the analytical description of the rubidium halogenides specific electrical conductivity and forecasting $\kappa(\text{RbAt})$.

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