

Metal reduction by hydrogen from the B_2O_3 -CaO-Ni(Zn,Pb,Cu)O melts thermodynamic modeling

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Abstract

Thermodynamic modeling is used to describe the metal reduction processes by hydrogen from oxide melt in the B_2O_3 -CaO- MeO (Me – Ni, Zn, Pb, Cu) system. Open systems approximation with periodic removal of metal particles and gases from the working melt composition is used in the method. By this work we present the thermodynamic modeling results of metal reduction processes (Ni, Cu, Pb, Zn) by Hydrogen. The reducible metals oxides content in the all melts was 3 mass %, and the mass ratio of B_2O_3 /CaO was taken as 3 to be close to eutectic composition. The calculations made it possible to determine such parameters as oxide melt compositions and elements reduction degree depending on the induced gas quantity. of the Nickel, Copper, Lead and Zinc reduction process simulation from B_2O_3 -CaO-MeO melts proved the reduction process by Hydrogen is similar to that which was earlier established when Carbon monoxide was used as the reducing agent. When Copper is reduced from CuO, the process occurs with intermediate Cu_2O oxide formation ($CuO \rightarrow Cu_2O \rightarrow Cu$). The Nickel ($NiO \rightarrow Ni$), Lead ($PbO \rightarrow Pb_s + Pb_g$) and Zinc ($ZnO \rightarrow Zn_g$) recovery have been realized by one stage. The non-ferrous metals change content in the oxide melt and the degrees of its reduction depending on temperature and reducing agent quantity introduced are described by the second-order polynomial functional equations. Comparison of the Carbon monoxide and Hydrogen used for Nickel, Copper, Lead, and Zinc reducing to 90% metallization degree proved much less Hydrogen consumption.

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