Supplementary Material

Of

Coordination of Au and Cu in Peridotite Melts Studied by First Principle Molecular Dynamic Simulations

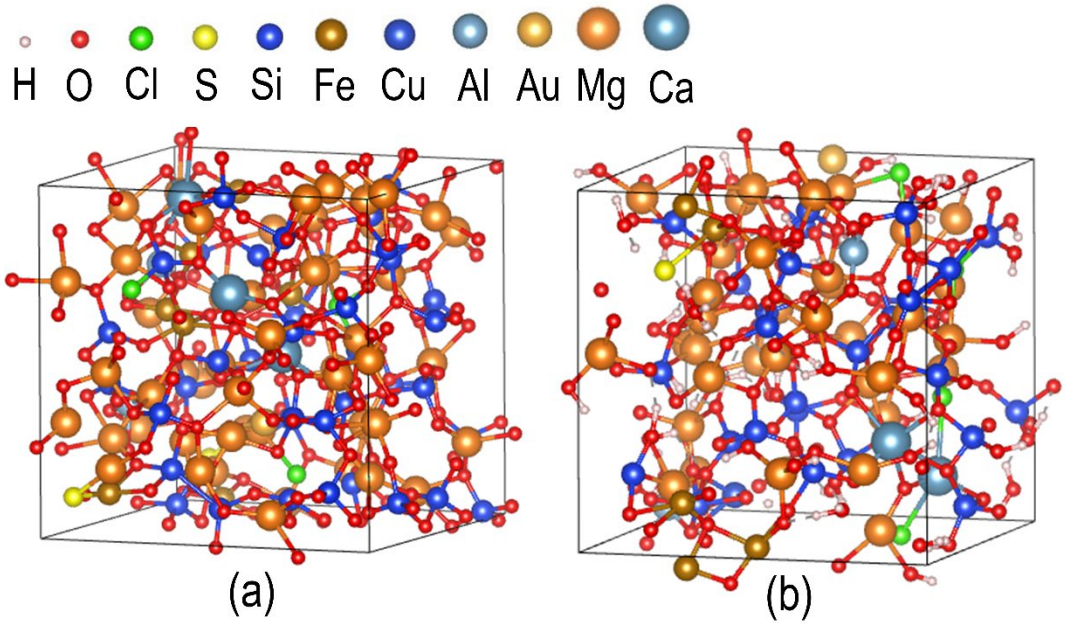


Figure S1. The simulated structures of (a) anhydrous and (b) hydrous peridotite melt system.