

FIG. 1. (Color online) The charge density distributions of monolayer SiC (unit: $|e|/\text{bohr}^3$).

TABLE I. Born effective charges Z^* of Si and C atoms and the dielectric constants (ϵ) of monolayer SiC. They along other directions are zero except xx , yy and zz directions.

Direction	$Z^*(\text{C})$	$Z^*(\text{Si})$	ϵ
xx	-3.674	3.674	2.531
yy	-3.674	3.674	2.531
zz	-0.271	0.271	1.180

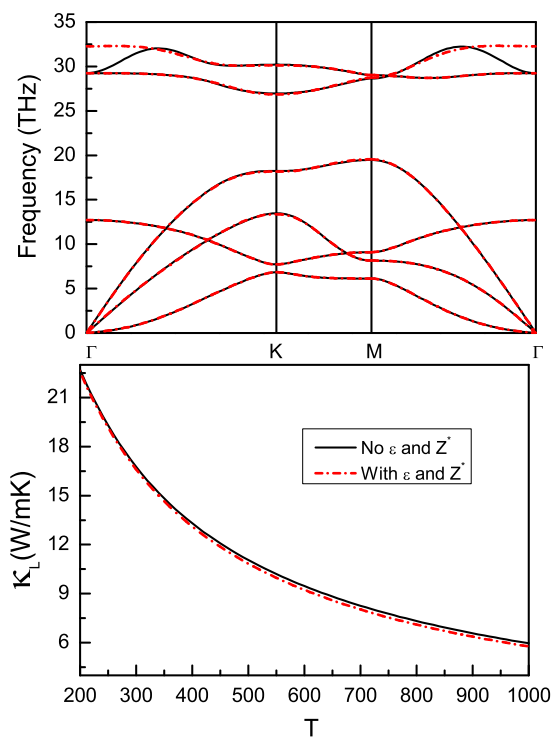


FIG. 2. (Color online) The phonon band structures (Top) and lattice thermal conductivity (Bottom) of monolayer SiC with or without the dielectric constants ϵ and Born effective charges Z^* .