

AlN - Aluminium Nitride

Electrical properties

Basic Parameter

Wurtzite crystal structure

Breakdown field	$1.2 \div 1.8 \times 10^6 \text{ V cm}^{-1}$	300 K
Mobility electrons	$300 \text{ cm}^2 \text{ V}^{-1} \text{ s}^{-1}$	300 K
Mobility holes	$14 \text{ cm}^2 \text{ V}^{-1} \text{ s}^{-1}$	300 K
Diffusion coefficient electrons	$7 \text{ cm}^2 \text{ s}^{-1}$	300 K
Diffusion coefficient holes	$0.3 \text{ cm}^2 \text{ s}^{-1}$	300 K
Electron thermal velocity	$1.85 \times 10^5 \text{ m s}^{-1}$	300 K
Hole thermal velocity	$0.41 \times 10^5 \text{ m s}^{-1}$	300 K

Conductivity σ	$10^{-3} \div 10^{-5} \Omega^{-1} \text{ cm}^{-1}$	290 K ; doped (Al_2O_3) single <i>p</i> -type crystals (blue)	Edwards et al. (1965)
	$10^{-11} \div 10^{-15} \Omega^{-1} \text{ cm}^{-1}$	300 K ; undoped single crystals (colorless or pale yellow) see also Conductivity vs. reciprocal temperature .	
Electron drift mobility μ_n	$\approx 300 \text{ cm}^2 \text{ V}^{-1} \text{ s}^{-1}$	300 K ; calculated	Chin et al. (1994)
Phonon-limited electron drift mobility μ_n	$\approx 2000 \text{ cm}^2 \text{ V}^{-1} \text{ s}^{-1}$	77 K ; calculated for very weak doped AlN	
Mobility holes μ_p	$14 \text{ cm}^2 \text{ V}^{-1} \text{ s}^{-1}$	290 K ; doped single crystal, the authors point out that this result must be viewed with some caution	Edwards et al. (1965)

