

APPENDIX AII

FUNDAMENTAL PHYSICAL CONSTANTS

Quantity	Symbol	Value	Unit
Electronic charge (magnitude)	q	1.60×10^{-19}	coulomb (C)
Electronic mass (magnitude)	m	9.11×10^{-31}	gram (g)
Unit nucleonic mass	m_{nucl}	1.66×10^{-24}	gram (g)
Avogadro's number	N_A	6.02×10^{23}	atoms/g-atom (mole)
Boltzmann's constant	k	1.38×10^{-13}	joule/kelvin ($J/^{\circ}K$)
Thermal energy (300°K)	kT	0.026	electron-volt (eV)
Thermal voltage (300°K)	kT/q	0.026	volt (V)
Speed of light (vacuum)	c	3.0×10^{10}	centimeter/second (cm/sec)
Planck's constant	h	6.63×10^{-34}	joule-second (J-sec)
Frequency (cycles/sec)	Hz	1.0	hertz (Hz)
Permittivity of free space	ϵ_0	8.85×10^{-14}	farad/centimeter (F/cm)
Relative dielectric constants:			
Germanium	$(\epsilon_r)_{Ge}$	16.0	—
Silicon	$(\epsilon_r)_{Si}$	11.9	—
Gallium arsenide	$(\epsilon_r)_{GaAs}$	13.1	—
Silicon dioxide	$(\epsilon_r)_{SiO_2}$	3.9	—
Silicon nitride	$(\epsilon_r)_{Si_3N_4}$	7.5	—
Sapphire	$(\epsilon_r)_{Sapp}$	9.2	—
Thermal conductivities (300°K)	κ_{Ge}	0.6	watts/cm ⁻² C (W/cm ⁻² C)
	κ_{Si}	1.5	watts/cm ⁻² C (W/cm ⁻² C)
	κ_{GaAs}	0.46	watts/cm ⁻² C (W/cm ⁻² C)
Lattice constants (300°K)	a_{Ge}	5.646	angstrom (Å)
	a_{Si}	5.430	angstrom (Å)
	a_{GaAs}	5.653	angstrom (Å)
Energy gaps (300°K)	$E_{\epsilon_{Ge}}$	0.66	electron-volt (eV)
	$E_{\epsilon_{Si}}$	1.12	electron-volt (eV)
	$E_{\epsilon_{GaAs}}$	1.42	electron-volt (eV)
Intrinsic concentrations (300°K)	n_{Ge}	2.4×10^{13}	centimeter ⁻³ (cm ⁻³)
	n_{Si}	1.5×10^{10}	centimeter ⁻³ (cm ⁻³)
	n_{GaAs}	1.8×10^6	centimeter ⁻³ (cm ⁻³)
Micrometer (micron)	μm	1.0×10^{-4}	centimeter (cm)
Angstrom unit	Å	1.0×10^{-8}	centimeter (cm)
Electron-volt	eV	1.6×10^{-19}	joules (J)

Appendix B

International System of Units

Quantity	Unit	Symbol	Dimensions
Length	meter	m	
Mass	kilogram	kg	
Time	second	s	
Temperature	kelvin	K	
Current	ampere	A	
Frequency	hertz	Hz	1/s
Force	newton	N	kg·m/s ²
Pressure	pascal	Pa	N/m ²
Energy	joule	J	N·m
Power	watt	W	J/s
Electric charge	coulomb	C	A·s
Potential	volt	V	J/C
Conductance	siemens	S	A/V
Resistance	ohm	Ω	V/A
Capacitance	farad	F	C/V
Magnetic flux	weber	Wb	V·s
Magnetic induction	tesla	T	Wb/m ²
Inductance	henry	H	Wb/A

Appendix C

Unit Prefixes^a

Multiple	Prefix	Symbol	Multiple	Prefix	Symbol
10^{18}	exa	E	10^{-1}	deci	d
10^{19}	peta	P	10^{-2}	centi	c
10^{20}	tera	T	10^{-3}	milli	m
10^{21}	giga	G	10^{-6}	micro	μ
10^{22}	mega	M	10^{-9}	nano	n
10^3	kilo	k	10^{-12}	pico	p
10^2	hecto	h	10^{-15}	femto	f
10	deka	da	10^{-18}	atto	a

^aAdopted by International Committee on Weights and Measures. (Compound prefixes should not be used; e.g., not $\mu\mu$ but p.)

Appendix D

Greek Alphabet

Letter	Lowercase	Uppercase	Letter	Lowercase	Uppercase
Alpha	α	A	Nu	ν	N
Beta	β	B	Xi	ξ	Ξ
Gamma	γ	Γ	Omicron	\circ	O
Delta	δ	Δ	Pi	π	Π
Epsilon	ϵ	E	Rho	ρ	\Rho
Zeta	ζ	Z	Sigma	σ	Σ
Eta	η	H	Tau	τ	T
Theta	θ	Θ	Upsilon	υ	Υ
Iota	ι	I	Phi	ϕ	Φ
Kappa	κ	K	Chi	χ	X
Lambda	λ	Λ	Psi	ψ	Ψ
Mu	μ	M	Omega	ω	Ω

Appendix E
Physical Constants

Quantity	Symbol	Value
Angstrom unit	\AA	$1 \text{\AA} = 10^{-8} \mu\text{m} = 10^{-8} \text{ cm}$
Avogadro constant	N_{AVO}	$6.02204 \times 10^{23} \text{ mol}^{-1}$
Bohr radius	a_B	0.52917\AA
Boltzmann constant	k	$1.38066 \times 10^{-23} \text{ J/K} (R/N_{AVO})$
Elementary charge	q	$1.60218 \times 10^{-19} \text{ C}$
Electron rest mass	m_0	$9.1095 \times 10^{-30} \text{ kg}$
Electron volt	eV	$1 \text{ eV} = 1.60218 \times 10^{-19} \text{ J}$ $= 23.053 \text{ kcal/mol}$
Gas constant	R	$1.98719 \text{ cal mol}^{-1} \text{ K}^{-1}$
Permeability in vacuum	μ_0	$1.25663 \times 10^{-8} \text{ H/cm} (4\pi \times 10^{-9})$
Permittivity in vacuum	ϵ_0	$8.85418 \times 10^{-14} \text{ F/cm} (1/\mu_0 c^2)$
Planck constant	h	$6.62617 \times 10^{-34} \text{ J-s}$
Reduced Planck constant	\hbar	$1.05458 \times 10^{-34} \text{ J-s} (h/2\pi)$
Proton rest mass	M_p	$1.67264 \times 10^{-27} \text{ kg}$
Speed of light in vacuum	c	$2.99792 \times 10^{10} \text{ cm/s}$
Standard atmosphere		$1.01325 \times 10^5 \text{ N/m}^2$
Thermal voltage at 300 K	kT/q	0.0259 V
Wavelength of 1-eV quantum	λ	$1.23977 \mu\text{m}$

Appendix F
Lattice Constants

	Element or Compound	Name	Crystal ^a Structure	Lattice Constant at 300 K (Å)
Element	C	Carbon (diamond)	D	3.56683
	Ge	Germanium	D	5.64613
	Si	Silicon	D	5.43095
	Sn	Grey Tin	D	6.48920
IV-IV	SiC	Silicon carbide	W	$a = 3.086, c = 15.117$
III-V	AlAs	Aluminum arsenide	Z	5.6605
	AlP	Aluminum phosphide	Z	5.4510
	AlSb	Aluminum antimonide	Z	6.1355
	BN	Boron nitride	Z	3.6150
	BP	Boron phosphide	Z	4.5380
	GaAs	Gallium arsenide	Z	5.6533
	GaN	Gallium nitride	W	$a = 3.189, c = 5.185$
	GaP	Gallium phosphide	Z	5.4512
	GaSb	Gallium antimonide	Z	6.0959
	InAs	Indium arsenide	Z	6.0584
	InP	Indium phosphide	Z	5.8686
	InSb	Indium antimonide	Z	6.4794
II-VI	CdS	Cadmium sulfide	Z	5.8320
	CdS	Cadmium sulfide	W	$a = 4.16, c = 6.756$
	CdSe	Cadmium selenide	Z	6.050
	CdTe	Cadmium telluride	Z	6.482
	ZnO	Zinc oxide	R	4.580
	ZnS	Zinc sulfide	Z	5.420
	ZnS	Zinc sulfide	W	$a = 3.82, c = 6.26$
IV-VI	PbS	Lead sulfide	R	5.9362
	PbTe	Lead telluride	R	6.4620

^aD = Diamond, W = Wurtzite, Z = Zincblende, R = Rock salt.

Appendix G

Properties of Important Semiconductors

Semiconductor	Bandgap (eV)		Mobility at 300 K (cm ² /V-s) ^a			Effective Mass m^*/m_0			
	300 K	0 K	Elec.	Holes	Band ^b	Elec.	Holes	ϵ_s/ϵ_0	
Element	C	5.47	5.48	1800	1200	I	0.2	0.25	5.7
	Ge	0.66	0.74	3900	1900	I	1.64 ^c	0.04 ^c	16.0
	Si	1.12	1.17	1500	450	I	0.082 ^d	0.28 ^f	
	Sn		0.082	1400	1200	D	0.98 ^e	0.16 ^e	11.9
IV-IV	α -SiC	2.996	3.03	400	50	I	0.60	1.00	10.0
III-V	AlSb	1.58	1.68	200	420	I	0.12	0.98	14.4
	BN	~7.5				I			7.1
	BP	2.0							
	GaN	3.36	3.50	380			0.19	0.60	12.2
	GaSb	0.72	0.81	5000	850	D	0.042	0.40	15.7
	GaAs	1.42	1.52	8500	400	D	0.067	0.082	13.1
	GaP	2.26	2.34	110	75	I	0.82	0.60	11.1
	InSb	0.17	0.23	80000	1250	D	0.0145	0.40	17.7
	InAs	0.36	0.42	33000	460	D	0.023	0.40	14.6
II-VI	InP	1.35	1.42	4600	150	D	0.077	0.64	12.4
	CdS	2.42	2.56	340	50	D	0.21	0.80	5.4
	CdSe	1.70	1.85	800		D	0.13	0.45	10.0
	CdTe	1.56		1050	100	D			10.2
	ZnO	3.35	3.42	200	180	D	0.27		9.0
	ZnS	3.68	3.84	165	5	D	0.40		5.2
IV-VI	PbS	0.41	0.286	600	700	I	0.25	0.25	17.0
	PbTe	0.31	0.19	6000	4000	I	0.17	0.20	30.0

^aThe values are for drift mobilities obtained in the purest and most perfect materials available to date.

^bTransverse effective mass.

^cI = indirect.

^dLight-hole effective mass.

^eLongitudinal effective mass.

^fHeavy-hole effective mass.

Properties of Ge, Si, and GaAs at 300 K

Properties	Ge	Si	GaAs
Atoms/cm ³	4.42×10^{22}	5.0×10^{22}	4.42×10^{22}
Atomic weight	72.60	28.09	144.63
Breakdown field(V/cm)	$\sim 10^5$	$\sim 3 \times 10^5$	$\sim 4 \times 10^5$
Crystal structure	Diamond	Diamond	Zincblende
Density (g/cm ³)	5.3267	2.328	5.32
Dielectric constant	16.0	11.9	13.1
Effective density of states in conduction band, N_c (cm ⁻³)	1.04×10^{19}	2.8×10^{19}	4.7×10^{17}
Effective density of states in valence band, N_v (cm ⁻³)	6.0×10^{18}	1.04×10^{19}	7.0×10^{18}
Effective Mass, m^*/m_0			
Electrons	$m_e^* = 1.64$ $m_e^* = 0.082$	$m_e^* = 0.98$ $m_e^* = 0.19$	0.067
Holes	$m_{lh}^* = 0.044$ $m_{hh}^* = 0.28$	$m_{lh}^* = 0.16$ $m_{hh}^* = 0.49$	$m_{lh}^* = 0.082$ $m_{hh}^* = 0.45$
Electron affinity, χ (V)	4.0	4.05	4.07
Energy gap (eV) at 300 K	0.66	1.12	1.424
Intrinsic carrier concentration (cm ⁻³)	2.4×10^{13}	1.45×10^{10}	1.79×10^6
Intrinsic Debye length (μ m)	0.68	24	2250
Intrinsic resistivity (Ω -cm)	47	2.3×10^5	10^8
Lattice constant (\AA)	5.64613	5.43095	5.6533
Linear coefficient of thermal expansion, $\Delta L/L\Delta T$ ($^{\circ}\text{C}^{-1}$)	5.8×10^{-6}	2.6×10^{-6}	6.86×10^{-6}
Melting point ($^{\circ}\text{C}$)	937	1415	1238
Minority carrier lifetime (s)	10^{-3}	2.5×10^{-3}	$\sim 10^{-8}$
Mobility (drift) ($\text{cm}^2/\text{V}\cdot\text{s}$)			
	3900 1900	1500 450	8500 400
Optical-phonon energy (eV)	0.037	0.063	0.035
Phonon mean free path λ_0 (\AA)	105	76 (electron) 55 (hole)	58
Specific heat ($\text{J/g}\cdot{}^{\circ}\text{C}$)	0.31	0.7	0.35
Thermal conductivity at 300 K (W/cm ⁻² · $^{\circ}\text{C}$)	0.6	1.5	0.46
Thermal diffusivity (cm^2/s)	0.36	0.9	0.24
Vapor pressure (Pa)	1 at 1330°C 10^{-6} at 760°C	1 at 1650°C 10^{-6} at 900°C	100 at 1050°C 1 at 900°C

Appendix I

**Properties of SiO₂
and Si₃N₄ at 300 K**

Insulator:	SiO ₂	Si ₃ N ₄
Structure	Amorphous	Amorphous
Melting point (°C)	~1600	—
Density (g/cm ³)	2.2	3.1
Refractive index	1.46	2.05
Dielectric constant	3.9	7.5
Dielectric strength (V/cm)	10 ⁷	10 ⁷
Infrared absorption band (μm)	9.3	11.5–12.0
Energy gap (eV)	9	~5.0
Thermal-expansion coefficient (°C ⁻¹)	5×10 ⁻⁷	—
Thermal conductivity (W/cm-K)	0.014	—
dc resistivity (Ω-cm)		
at 25°C	10 ¹⁴ –10 ¹⁶	~10 ¹⁴
at 500°C	—	~2×10 ¹³
Etch rate in buffered HF ^a (Å/min)	1000	5–10

^aBuffered HF: 34.6% (wt.) NH₄F, 6.8% (wt.) HF, 58.6% H₂O.