

List of frequently used symbols and abbreviations; conversion factors

1. Symbols

| Symbol | Unit | Property |
|---------------------------------|---|---|
| a | | activity |
| a_{O_2} | | oxygen activity |
| a, b, c | Å, nm | lattice parameters |
| A | | absorbance |
| b | | electron-hole mobility ratio |
| B | T, G | magnetic induction |
| B_i | | internal magnetic induction |
| B | $\text{cm}^2 \text{K}^{-1} \text{s}^{-1}$ | Nernst coefficient |
| $B_0, (B_S, B_T)$ | Pa, bar, N m^{-2} | bulk modulus (adiabatic, isothermal) |
| c_{lm} | Pa, dyn cm^{-2} | elastic moduli (stiffnesses) |
| C | $\text{J mol}^{-1} \text{K}^{-1}$, $\text{cal mol}^{-1} \text{K}^{-1}$, | heat capacity |
| C_p | | heat capacity at constant pressure |
| C_m | | magnetic heat capacity |
| c | $\text{J g}^{-1} \text{K}^{-1}$ | specific heat capacity |
| C_m | $\text{emu K mol}^{-1} =$ $\text{cm}^3 \text{K mol}^{-1}$ | Curie constant per mole |
| d | g cm^{-3} | density |
| d_X | | X-ray density |
| d | μm , Å, nm | thickness, (interatomic) distance, bond length |
| d_{opt} | | optical density ($\log I_0/I$) |
| D | $\text{cm}^2 \text{s}^{-1}$ | diffusion coefficient |
| $D_{n(p)}$ | | diffusion coefficient for electrons (holes) |
| Dq | eV | crystal field splitting parameter |
| e | | polarization vector |
| e | C | elementary charge |
| $e^{*(T)}$ | e | (transverse) effective ionic charge |
| E | Pa, bar | Young's modulus |
| $E_{[\text{hkl}]}$ | | Young's modulus measured in [hkl] direction |
| E | V cm^{-1} , V m^{-1} | electric field strength |
| E | eV, Ry, J | energy |
| $E_{a(d)}$ | | ionization energies of acceptors (donors), energy of acceptor (donor) state |
| E_A | | activation energy (mostly of conductivity) |
| E_b | | binding energy |
| E_{mc}^b | | binding energy of bound multiexcitons |
| E_C | | carbon donor level |
| $E_{C(V)}$ | | band edge of conduction (valence) band |
| E_{crit} | | critical point energy |
| E_F | | Fermi energy |
| E_g | | energy gap, band gap |
| $E_{g,\text{th}}$ | | energy gap extrapolated to 0 K (thermal energy gap) |
| $E_{g,\text{dir}} (\text{ind})$ | | direct (indirect) energy gap |
| E_{kin} | | kinetic (photoelectron) energy |

| Symbol | Unit | Property |
|------------------------|---|---|
| E_p | | primary electron energy (in ELS) |
| $E_{pl(pc)}, E_{peak}$ | | photoluminescence (photoconductivity) peak energy |
| E_r | | relative energy |
| E_{tot} | | total energy |
| E_t | | trapping energy of charge carrier |
| E_{thr} | | photoelectric threshold energy |
| $E_{0,1,2,...}$ | | energies of critical points in optical spectra (see also E_{crit}) |
| $E(\Gamma_6)_{..}$ | | energy of band edge of type $\Gamma_6 \dots$ |
| ΔE | | energy loss |
| f | Hz | frequency |
| f | | oscillator strength |
| f | kg | frictional force |
| g | | spectroscopic splitting factor |
| $g_{ }$ | | longitudinal g -factor |
| g_{\perp} | | transverse g -factor |
| g_c, g_n | | g -factor of conduction electrons |
| g_v, g_p | | g -factor of holes |
| $g(E)$ | various units | density of states (also $N(E)$ or DOS is used) |
| $g_c(v)$ | | density of conduction (valence) band states |
| g_v | J | Gibbs energy of formation of a neutral cation vacancy |
| G | Pa, bar | shear modulus |
| $G_{[hkl]}$ | | shear modulus measured in $[hkl]$ direction |
| G | J mol ⁻¹ , cal mol ⁻¹ | free energy |
| ΔG_f^0 | J mol ⁻¹ | standard free energy of formation |
| ΔG_{tr} | J mol ⁻¹ | Gibbs energy of phase transition |
| h | | Planck constant |
| h | Å | height (of icosahedron) |
| $h\nu$ | eV | photon energy |
| H | Pa, N m ⁻² , kg mm ⁻² | hardness, microhardness |
| $H_{B,V,K}$ | | hardness (Brinell, Vickers, Knoop) |
| H | Oe, A m ⁻¹ | magnetic field (strength), also given as $\mu_0 H$ in tesla (T)) |
| H_A | | applied magnetic field |
| H_{cr} | | critical magnetic field |
| H_{ext} | | external magnetic field |
| $\Delta H_{(p-p)}$ | Oe, A m ⁻¹ | (peak to peak) linewidth of resonance spectrum |
| H_T | J mol ⁻¹ | enthalpy at temperature T |
| $\Delta H_f^{(0)}$ | J mol ⁻¹ , cal mol ⁻¹ , J g-atom ⁻¹ | (standard) heat of formation |
| ΔH_{tr} | J mol ⁻¹ | heat of transformation |
| i | A cm ⁻² | current density |
| I | A | electric current |
| I_{ph} | | photocurrent |
| I | various units | intensity |
| I_{lum} | | luminescence intensity |
| I_R | | Raman intensity |
| J | | total orbital angular momentum quantum number |
| J | eV | exchange interaction energy (J/k_B in K) |
| k | | extinction coefficient (absorption index) |

| Symbol | Unit | Property |
|----------------------------------|--|--|
| k | $\text{\AA}^{-1}, \text{nm}^{-1}$ | wavevector of electrons |
| k | $\text{dyn } \text{\AA}^{-1}$ | force constant |
| $k_{\text{BB}}, k_{\text{B-B}}$ | | force constant between B atoms |
| k, k_{B} | J K^{-1} | Boltzmann constant |
| K | cm^{-1} | absorption coefficient (sometimes α is used) |
| K, K_{B} | | Knight shift |
| K_{C} | $\text{N m}^{-3/2}$ | fracture toughness |
| L, l | cm | length |
| $\Delta l/l$ | | linear thermal elongation |
| m_0 | g | electron mass |
| $m_{\text{n(p)}}$ | m_0 | effective mass of electrons (holes) |
| $m_{\omega_{\text{c}}}$ | m_0 | effective cyclotron resonance mass |
| $m_{\omega_{\text{p}}}$ | m_0 | effective plasma frequency mass |
| m^* | m_0 | effective mass |
| m^{**} | m_0 | polaronic mass |
| $m_{\perp, \parallel}$ | m_0 | transverse, longitudinal effective mass |
| $M_{(\text{s})}$ | G | (saturation) magnetization per unit volume |
| n | | (linear) refractive index |
| $n_{\text{a,b,c}}$ | | refractive index in a, b, c direction |
| Δn | | birefringence |
| n, n_{e} | $\text{g}^{-1}, \text{atom}^{-1}, \text{unit cell}^{-1}, \text{cm}^{-3}$ | electron concentration (n also carrier concentration in general) |
| $n_{\text{a(d)}}$ | | acceptor (donor) concentration |
| $n_{\text{eff}}, N_{\text{eff}}$ | | effective number of electrons per FU, per atom (contributing to opt. properties) |
| n_{H} | | Hall density |
| n_{i} | | intrinsic carrier concentration |
| n_{S} | | spin density |
| n_{t} | | density of trapping levels |
| N | | count rate, number of electrons |
| $N(E)$ | various units | density of states (see also $g(E)$) |
| p | Pa, bar | hydrostatic pressure |
| p_{tr} | | (phase) transition pressure |
| p | μ_{B} | magnetic moment |
| p_{eff} | | effective (paramagnetic) moment |
| $p_{\text{FU}}, p_{\text{m}}$ | | magnetic moment per formula unit |
| $p_{\text{M}}, p_{\text{A}}$ | | magnetic moment per ion M, per atom A |
| p_{RE} | | magnetic moment per rare earth atom |
| p_{s} | | saturation magnetic moment |
| p | cm^{-3} | hole concentration |
| P | C m^{-2} | spontaneous polarization |
| q | e | Mulliken effective charge |
| q | $\text{\AA}^{-1}, \text{nm}^{-1}$ | wavevector of phonons |
| Q | \AA^{-1} | wavevector (frequently reciprocal vector component for neutron scattering) |
| Q^{-1} | | internal friction |
| $r_{(\text{M})}$ | $\text{cm}, \text{nm}, \text{\AA}$ | radius (of atom M) |
| R | | reflectivity, reflectance |
| R | Ω | resistance |
| R | $\text{J K}^{-1} \text{mol}^{-1}, \text{cal mol}^{-1} \text{K}^{-1}$ | gas constant |
| R_{H} | $\text{m}^3 \text{C}^{-1}, \text{cm}^3 \text{C}^{-1}$ | normal Hall coefficient |

| Symbol | Unit | Property |
|-----------------------|--|--|
| R_s | $\text{m}^3 \text{C}^{-1}, \text{cm}^3 \text{C}^{-1}$ | anomalous Hall coefficient |
| S | | spin quantum number |
| S | $\text{J mol}^{-1} \text{K}^{-1}, \text{cal mol}^{-1} \text{K}^{-1}$ | entropy |
| ΔS_f^0 | $\text{J mol}^{-1} \text{K}^{-1}$ | entropy of formation |
| ΔS_{tr} | $\text{J mol}^{-1} \text{K}^{-1}$ | entropy of phase transition |
| S | $\text{mV K}^{-1}, \mu\text{V K}^{-1}$ | Seebeck coefficient (thermoelectric power) |
| $S(Q, E)$ | $\text{barn eV}^{-1} \text{sr}^{-1} \text{atom}^{-1}$ | INS response function |
| t | s, min, h, d, yr | time |
| t_i | | transport numbers (t_+ , t_- : for cations, anions) |
| T | | transmission, transmittance |
| T | K, °C | temperature |
| T_c | | superconductor transition temperature |
| T_C | | Curie temperature |
| T_{dec}, T_{decomp} | | decomposition temperature |
| T_g | | glass transition temperature |
| T_m | | melting temperature |
| T_N | | Néel temperature |
| T_{perit} | | peritectic (decomposition) temperature |
| T_{prep} | | preparation temperature |
| T_Q | | antiferroquadrupolar temperature |
| T_s | | substrate temperature, spin flip temperature |
| T_{tr} | | crystallographic transition temperature |
| T_v | | Verwey temperature |
| T_α | | critical temperature for cluster formation |
| T_0 | | characteristic temperature in Mott's law of variable range hopping |
| u | | position parameter |
| U | V | voltage (also V is used) |
| U_H | | Hall voltage |
| U_i | | induced voltage |
| U | eV | Hubbard self-energy, electrostatic correlation energy |
| U_{ij} | \AA^2 | temperature factors |
| U_{iso} | | isotropic temperature factor |
| U_{equiv} | | equivalent temperature factor |
| v | cm s^{-1} | velocity, sound velocity |
| v_b | | bulk velocity |
| $v_{l(L)}, v_{t(T)}$ | | longitudinal, transverse sound velocity |
| v_s | | shear sound velocity |
| v | | valence |
| V | $\text{\AA}^3, \text{nm}^3$ | volume |
| V_m | $\text{cm}^3 (\text{mol}^{-1})$ | molar volume |
| W | eV | width of valence or conduction bands |
| x, y, z | | fractional coordinates of atoms in the unit cell |
| x | | concentration |
| X_{ik} | bar | stress tensor (6×6) (in the literature often labeled T_{ij}) |
| X_k | bar | stress vector (6-component) |
| Y | | yield of electrons, photoemissive yield |
| z | K^{-1} | figure of merit ($z = S^2 \sigma / \kappa$) |
| Z | | atomic number, coordination number |

| Symbol | Unit | Property |
|--|--|---|
| α (α_F) | | Fröhlich polaron coupling constant |
| α | cm^{-1} | absorption coefficient |
| α | K^{-1} | linear thermal expansion coefficient |
| $\alpha_{a,b,c}$ | | lin. thermal expansion coefficient in a , b , c direction |
| α, β, γ | deg | unit cell angles |
| β | K^{-1} | volume thermal expansion coefficient |
| γ, γ_G | | Grüneisen constant |
| γ | | microbrittleness |
| γ | cm^{-1} | damping constant |
| γ | $\text{Hz T}^{-1}, \text{Hz Oe}^{-1}$ | nuclear gyromagnetic ratio |
| γ | $\text{J mol}^{-1} \text{K}^{-2}$ | coefficient of electronic heat capacity |
| γ | Hz | hopping rate |
| Γ | $\text{Oe}, \text{\AA}^{-1}$ | linewidth (e.g. of nuclear Bragg reflection, in ESR) |
| Γ | cm^{-1} | linewidth (e.g. of phonon wavenumber) |
| δ | ppm | chemical shift |
| δ | | Anderson-Grüneisen constant |
| $\tan \delta$ | | dielectric loss tangent (ϵ_2/ϵ_1) |
| Δ_{cf} | eV | CEF splitting energy |
| Δ_{el} | eV | electronic stabilization energy |
| Δ_{ex} | eV | exchange splitting energy |
| ϵ_0 | F cm^{-1} | permittivity of free space |
| $\epsilon = \epsilon_1 - i \epsilon_2$ | | dielectric constant |
| ϵ_∞ | | high frequency limit of dielectric constant |
| ϵ_0 | | static dielectric constant |
| $\epsilon(0), \epsilon_0$ | | low frequency limit of dielectric constant |
| ϵ_1, ϵ_2 | | real, imaginary part of dielectric constants |
| ϵ_L | | lattice part of dielectric constant |
| ϵ^e | | contribution of free electron to ϵ |
| $\text{Im } \epsilon^{-1}$ | | energy loss function |
| $\Delta \epsilon$ | | oscillator strength |
| ζ | | reduced wavevector coordinate |
| θ | deg | diffraction angle |
| Θ, Θ_p | K | paramagnetic Curie temperature, characteristic temperature |
| Θ_a | K | asymptotic Curie temperature |
| Θ_D | K | Debye temperature |
| Θ_E | K | Einstein temperature |
| κ | $\text{cm}^2 \text{dyn}^{-1}, \text{m}^2 \text{N}^{-1}, \text{Pa}^{-1}$ | compressibility (= 1/bulk modulus) |
| κ_L | | linear compressibility |
| κ_V | | volume compressibility |
| κ | $\text{W cm}^{-1} \text{K}^{-1}, \text{W m}^{-1} \text{K}^{-1}, \text{W mol}^{-1} \text{K}^{-1}$ | thermal conductivity |
| κ_e | | charge carrier contribution to κ |
| κ_L | | lattice contribution to κ |
| λ | nm, μm | wavelength |
| λ_L | | wavelength of laser light |
| Λ | cm | phonon mean free path |

| Symbol | Unit | Property |
|----------------------------------|--|--|
| μ | $\text{cm}^2 \text{V}^{-1} \text{s}^{-1}$ | mobility of charge carriers |
| $\mu_{a,b,c}$ | | mobility in a , b , c direction |
| μ_{dr} | | drift mobility |
| μ_{H} | | Hall mobility |
| $\mu_{\text{n}}, \mu_{\text{p}}$ | | electron (hole) mobility |
| μ_{B} | J T^{-1} | Bohr magneton |
| ν | Hz | frequency |
| ν_{p} | | plasma frequency |
| ν_{R} | | Raman frequency |
| $h\nu$ | eV | photon energy |
| ν | | Poisson's ratio |
| $\bar{\nu}$, ν/c | cm^{-1} | wavenumber |
| π_{ik} | $\text{cm}^2 \text{dyn}^{-1}$ | piezoresistance coefficient |
| Π | eV | Peltier energy |
| ρ | $\Omega \text{ cm}$, $\Omega \text{ m}$ | resistivity |
| $\rho_{a,b,c}$ | | resistivity in a , b , c direction |
| $\Delta\rho/\rho_0$ | | magnetoresistance |
| σ | Pa | microstrength |
| σ | $\text{emu g}^{-1} = \text{G cm}^3 \text{g}^{-1}$, $\text{A m}^2 \text{kg}^{-1}$ | magnetic moment per unit mass = specific magnetization |
| σ_{m} | $\text{emu mol}^{-1} =$ $\text{G cm}^3 \text{mol}^{-1}$ | magnetic moment per mole = molar magnetization |
| σ | $\Omega^{-1} \text{cm}^{-1}$ | electrical conductivity |
| $\sigma_{a,b,c}$ | | conductivity in a , b , c direction |
| σ_{d} | | dark conductivity |
| σ_{i} | | intrinsic conductivity |
| $\sigma_{\text{n(p)}}$ | | conductivity of electrons (holes) |
| σ_{ph} | | photoconductivity |
| σ_{opt} | $\Omega^{-1} \text{cm}^{-1}, \text{s}^{-1}$ | optical conductivity |
| τ | s | relaxation time, rise or decay time, lifetime of carriers |
| Φ | eV | work function |
| Φ_{e} | | work function of electrons |
| Φ | eV | inner lattice potential |
| χ | | magnetic susceptibility |
| χ_{g} | $\text{emu g}^{-1} = \text{cm}^3 \text{g}^{-1}$ | magnetic susceptibility per gram |
| χ_{m} | $\text{emu mol}^{-1} = \text{cm}^3 \text{mol}^{-1}$ | magnetic susceptibility per mole |
| χ_{v} | | magnetic volume susceptibility |
| ω | s^{-1} | angular (circular) frequency |
| ω_{p} | | plasma resonance frequency |
| $\hbar\omega$ | eV | photon energy |

2. Abbreviations

| | |
|-----------|---|
| A | actinoid element |
| a | acceptor |
| a | amorphous |
| ac | alternating current |
| AES | atomic emission spectroscopy |
| AF, AFM | antiferromagnetic |
| AFQ | antiferroquadrupolar |
| APB | antiphase boundary |
| APW | augmented plane wave (method) |
| arb | arbitrary |
| av | average |
| bcc | body centered cubic |
| bct | body centered tetragonal |
| BZ | Brillouin zone |
| c | mostly as subscript: cubic, critical or conduction band |
| calc | calculated |
| CB | conduction band |
| CCDW | commensurate charge density wave |
| cr, crit | mostly as subscript: critical |
| CS | crystallographic shear (plane) |
| cub | cubic |
| CVD | chemical vapour deposition |
| d, D | donor |
| dc | direct current |
| dhcp | double hexagonal close-packed |
| dHvA | de Haas van Alphen oscillations (method) |
| DFT | density functional theory |
| dir | direct |
| DOS | density of states |
| DTA | differential thermal analysis |
| e | electron |
| EDC | electron (intensity) distribution curve |
| eff | effective |
| ELNES | energy loss near edge structure |
| ELS, EELS | electron energy loss spectrum |
| EMF | electromotive force |
| EPMA | electron probe microanalysis |
| EPR, epr | electron paramagnetic resonance |
| ESCA | electron spectroscopy for chemical analysis |
| ESR, esr | electron spin resonance |
| EXELFS | electron-loss fine structure |
| exp | experimental |
| fcc | face centered cubic |
| F | ferromagnetic |
| FIR, fir | far infrared |
| FLAPW | full potential linearized augmented plane wave (method) |
| FT | Fourier transform |
| FWHM | full width at half maximum |
| h | hole |
| hcp | hexagonal close-packed |
| hex | hexagonal |

| | |
|-------------------------------|---|
| HIP | hot isostatic pressing |
| HT | high temperature $T > 300$ K |
| I | insulator |
| i | intrinsic; sometimes used for interstitial |
| ICDW | incommensurate charge density wave |
| ind | indirect |
| INDO | intermediate neglect of differential overlap |
| ion | ionic, ionization |
| IR, ir | infrared, irradiation |
| KKR | Kohn-Korringa-Rostoker method |
| L, l | mostly as subscript: longitudinal or lattice |
| L, l, liq | liquid |
| LA | longitudinal acoustic |
| LCAO | linear combination of atomic orbitals |
| LMTO | linearized muffin-tin orbital ((linear combination of muffin-tin orbitals (method)) |
| Ln | lanthanide |
| LO | longitudinal optical |
| LSD | local spin density method |
| LT | low temperature, mainly < 10 K |
| LTO | longitudinal and transverse optical |
| m | monoclinic (mostly subscript) |
| M, Me, M' | metal |
| magn | mostly as subscript: magnetic |
| MAS | magic-angle spinning |
| min (max) | minimum (maximum) |
| MO | molecular orbit |
| Mn ⁺ _{Mn} | Mn ion on Mn site, positively charged |
| MSC | magnetically stimulated current |
| MVFF | modified valence force field |
| NMR | nuclear magnetic resonance |
| obs | observed |
| oct | octahedral |
| O _d | oxygen deficiency (concentration of deficient oxygen atoms) |
| opt | optical |
| OPW | orthogonalized plane wave |
| p | hole |
| P | paramagnetic |
| PC | point contact |
| PDOS | partial density of states |
| PE | photoelectron spectrum |
| PEELS | parallel electron energy-loss spectra |
| ph | as subscript: photon, phonon, photo- |
| PM | paramagnetic |
| PVD | physical vapor deposition |
| R, RE, REE | rare earth element |
| R | rutile (mostly subscript) |
| R | as subscript: Raman |
| rel | relative |
| res | resolution |
| RF | radio frequency |
| rh | rhombohedral |
| RPA | random phase approximation |

| | |
|--------------------|--|
| RT | room temperature |
| s | surface |
| SC | semiconductor |
| SCF | self consistent field |
| SICF | single ion in a crystal field |
| SMT | semiconductor-metal transition |
| SXS | soft X-ray spectrum |
| T, TM | transition element |
| t, T | as subscripts or superscripts: transverse |
| TA | transverse acoustic |
| TDOS | total density of states |
| TEM | transmission electron microscopy |
| tetr | mostly as subscripts: tetragonal |
| TG | thermal gravimetry |
| th | thermal, sometimes for theoretical |
| theor | theoretical |
| TIP | temperature independent paramagnetism |
| TO | transverse optical |
| tot | mostly as subscript: total |
| tr | transition (subscript for phase transition parameters) |
| UPS, UPE | UV photoemission spectroscopy |
| uv, UV | ultraviolet |
| v | mostly as subscript: valence band |
| vac | vacuum, sometimes for vacancy |
| VB | valence band, as subscript v is used |
| VCA | virtual crystal approximation |
| V _{Ga...} | vacancy on Ga... site |
| V' _{Mn} | Mn vacancy, negatively charged |
| VT | vapor transport |
| X | anion (e.g. S, Se, Te) |
| XANES | X-ray absorption near edge structural spectra |
| XCS | X-ray chemical shift |
| XPS, XPE | X-ray photoelectron spectroscopy |
| XRD | X-ray diffraction |
| ⊥, | perpendicular, parallel to a crystallographic axis |

3. Conversion tables

A. Conversion factors from the SIU system to the CGS-esu and the CGS-emu systems.

| Quantities | Sym-bols | SIU | CGS-esu (non-rationalized) | CGS-emu |
|--|------------|---|--|--|
| bulk modulus | B | Pa (= N m ⁻²) | 10 dyn cm ⁻² | 10 dyn cm ⁻² |
| magnetic induction | B | T (= Wb m ⁻²) = Vs m ⁻² | 10 ^{-6/3} esu | 10 ⁴ G |
| molar heat capacity at const. pressure | C_p | J K ⁻¹ mol ⁻¹ | 10 ⁷ erg K ⁻¹ mol ⁻¹ (=0.239 cal K ⁻¹ mol ⁻¹) | 10 ⁷ erg K ⁻¹ mol ⁻¹ (=0.239 cal K ⁻¹ mol ⁻¹) |
| elastic moduli (stiffnesses) | c_{lm} | N m ⁻² (= Pa) | 10 dyn cm ⁻² | 10 dyn cm ⁻² |
| density | d | kg m ⁻³ | 10 ⁻³ g cm ⁻³ | 10 ⁻³ g cm ⁻³ |
| piezoelectric strain coefficient | d_{ik} | C N ⁻¹ (= m V ⁻¹) | 3 · 10 ⁴ esu | 10 ⁻⁶ emu |
| strain tensor | e_{ik} | dimensionless | 1 (dimensionless) | 1 (dimensionless) |
| piezoelectric stress coefficients | e_{ik} | C m ⁻² | 3 · 10 ⁵ esu | 10 ⁻⁵ emu |
| Young's modulus | E | N m ⁻² (= Pa) | 10 dyn cm ⁻² | 10 dyn cm ⁻² |
| electric field strength | E | V m ⁻¹ | 10 ^{-4/3} esu | 10 ⁶ emu |
| piezoelectric strain coefficients | g_{ik} | m ² C ⁻¹ | 10 ^{-5/3} esu | 10 ⁵ emu |
| molar free energy change | ΔG | J mol ⁻¹ (= 0.239 cal mol ⁻¹) | 10 ⁷ erg mol ⁻¹ (= 0.239 cal mol ⁻¹) | 10 ⁷ erg mol ⁻¹ |
| piezoelectric stress coefficient | h_{ik} | V m ⁻¹ (= N C ⁻¹) | 10 ^{-4/3} esu | 10 ⁶ emu |
| hardness | H | N m ⁻² (= Pa) | 10 dyn cm ⁻² | 10 dyn cm ⁻² |
| magnetic field strength | H | A m ⁻¹ | 12 π · 10 ⁷ esu | 4 π · 10 ⁻³ Oe |
| molar enthalpy change | ΔH | J mol ⁻¹ (= 0.239 cal mol ⁻¹) | 10 ⁷ erg mol ⁻¹ (= 0.239 cal mol ⁻¹) | 10 ⁷ erg mol ⁻¹ |
| current density | i | A m ⁻² | 3 · 10 ⁵ esu | 10 ⁻⁵ emu |
| elastoresistance coefficients | m_{ik} | dimensionless | 1 (dimensionless) | 1 (dimensionless) |
| pressure | p | Pa (= 10 ⁻⁵ bar) | 10 dyn cm ⁻² (=1.019 · 10 ⁻⁵ kg cm ⁻² =7.5 · 10 ⁻³ Torr) | 10 dyn cm ⁻² (=1.019 · 10 ⁻⁵ kg cm ⁻² =7.5 · 10 ⁻³ Torr) |
| dielectric polarization | P | C m ⁻² | 3 · 10 ⁵ esu | 10 ⁻⁵ emu |
| pyroelectric coefficient | p_i | C m ⁻² K ⁻¹ | 3 · 10 ⁵ esu K ⁻¹ | 10 ⁻⁵ emu K ⁻¹ |
| elasto-optic constant (in cubic crystals) | p_{ij} | dimensionless | 1 (dimensionless) | 1 (dimensionless) |
| piezo-optic constant (in cubic crystals) | q_{ij} | dimensionless | 1 (dimensionless) | 1 (dimensionless) |
| Hall coefficient | R_H | m ³ C ⁻¹ | 10 ^{-3/3} esu | 10 ⁷ emu |
| linear electro-optic constant | r_{ij} | m V ⁻¹ | 3 · 10 ⁴ esu | 10 ⁻⁶ emu |
| elastic compliances | s_{ml} | m ² N ⁻¹ | 10 ⁻¹ cm ² dyn ⁻¹ | 10 ⁻¹ cm ² dyn ⁻¹ |
| molar entropy change | ΔS | J K ⁻¹ mol ⁻¹ | 10 ⁷ erg K ⁻¹ mol ⁻¹ (= 0.239 cal mol ⁻¹) | 10 ⁷ erg K ⁻¹ mol ⁻¹ (= 0.239 cal mol ⁻¹) |
| stress tensor | X_{ij} | N m ⁻² (= Pa) | 10 dyn cm ⁻² | 10 dyn cm ⁻² |
| thermal conductivity | κ | W m ⁻¹ K ⁻¹ (= J m ⁻¹ s ⁻¹ K ⁻¹) | 10 ⁵ erg cm ⁻¹ s ⁻¹ K ⁻¹ | 10 ⁵ erg cm ⁻¹ s ⁻¹ K ⁻¹ |
| dielectric constant | ϵ | dimensionless | 1 (dimensionless) | 1 (dimensionless) |

| Quantities | Sym- bols | SIU | CGS-esu (non-rationalized) | CGS-emu |
|---|--------------|------------------------------|---|--|
| Conversion factors (cont.) | | | | |
| piezoresistance tensor coefficients | π_{ik} | $\text{m}^2 \text{N}^{-1}$ | $10^{-1} \text{cm}^2 \text{dyn}^{-1}$ | $10^{-1} \text{cm}^2 \text{dyn}^{-1}$ |
| piezooptic constant (in cubic crystals) | π_{ik} | $\text{m}^2 \text{N}^{-1}$ | $10^{-1} \text{cm}^2 \text{dyn}^{-1}$ | $10^{-1} \text{cm}^2 \text{dyn}^{-1}$ |
| resistivity | ρ | Ωm | $10^{-9}/9 \text{esu}$ | 10^{-11}emu |
| conductivity | σ | $\Omega^{-1} \text{m}^{-1}$ | $9 \cdot 10^9 \text{esu}$ | 10^{-11}emu |
| magnetic volume susceptibility | χ_v | dimensionless | $1/4\pi$ (dimensionless) | $1/4\pi$ (dimensionless) |
| magnetic mass susceptibility | χ_g | $\text{m}^3 \text{kg}^{-1}$ | $10^3/4\pi \text{esu cm}^3 \text{g}^{-1}$ | $10^3/4\pi \text{emu g}^{-1}$ ($= 10^3/4\pi \text{cm}^3 \text{g}^{-1}$) |
| magnetic molar susceptibility | χ_m | $\text{m}^3 \text{mol}^{-1}$ | $10^6/4\pi \text{esu cm}^3 \text{mol}^{-1}$ | $10^6/4\pi \text{emu mol}^{-1}$ ($= 10^6/4\pi \text{cm}^3 \text{mol}^{-1}$) |

Energy conversion

| | |
|---|--|
| Energy: $E = e V = h \nu = h c \bar{\nu}$ | $1 \text{ V As} = 1 \text{ J} = 10^7 \text{ erg} = 2.38845 \cdot 10^{-4} \text{ kcal}$ |
|---|--|

| Energy and equivalent quantities | | E | V | ν | $\bar{\nu}$ |
|-------------------------------------|-----------------------------|--------------------------|--------------------------|-------------------------|--------------------------|
| | | [J] | [V] | [Hz, s ⁻¹] | [cm ⁻¹] |
| 1J | $\underline{\underline{=}}$ | 1 | $6.2415 \cdot 10^{18}$ | $1.50916 \cdot 10^{33}$ | $5.03403 \cdot 10^{22}$ |
| 1V | $\underline{\underline{=}}$ | $1.60219 \cdot 10^{-19}$ | 1 | $2.41797 \cdot 10^{14}$ | $8.06547 \cdot 10^3$ |
| 1s ⁻¹ (= 1 Hz) | $\underline{\underline{=}}$ | $6.62619 \cdot 10^{-34}$ | $4.13550 \cdot 10^{-15}$ | 1 | $3.33564 \cdot 10^{-11}$ |
| 1 cm ⁻¹ | $\underline{\underline{=}}$ | $1.98648 \cdot 10^{-23}$ | $1.23979 \cdot 10^{-4}$ | $2.99792 \cdot 10^{10}$ | 1 |

Error: Experimental errors are frequently given in parentheses referring to the last decimal places. For example, 1.352(12) stands for 1.352 ± 0.012 and 342.5(21) stands for 342.5 ± 2.1 .