## 6. ATOMIC AND NUCLEAR PROPERTIES OF MATERIALS

Table 6.1. Revised May 2002 by D.E. Groom (LBNL). Gases are evaluated at 20°C and 1 atm (in parentheses) or at STP [square brackets]. Densities and refractive indices without parentheses or brackets are for solids or liquids, or are for cryogenic liquids at the indicated boiling point (BP) at 1 atm. Refractive indices are evaluated at the sodium D line. Data for compounds and mixtures are from Refs. 1 and 2. Futher materials and properties are given in Ref. 3 and at http://pdg.lbl.gov/AtomicNuclearProperties.

Material	Z	A	$\langle Z/A \rangle$	collision	interaction length $\lambda_I$	$\frac{dE/dx _{\min}}{\left\{\frac{\text{MeV}}{\text{g/cm}^2}\right\}}$		$X_0$ $X_0$ $X_0$ $X_0$	Density $\{g/cm^3\}$ $(\{g/\ell\}$ for gas)	Liquid boiling point at 1 atm(K)	Refractive index $n$ $((n-1)\times10^6$ for gas)
$H_2$ gas	1	1.00794	0.99212	43.3	50.8	(4.103)	$61.28 \frac{d}{}$	(731000)	(0.0838)[0.0899]		[139.2]
$H_2$ liquid	1	1.00794	0.99212	43.3	50.8	4.034	$61.28 \ ^d$	866	0.0708	20.39	1.112
$D_2$	1	2.0140	0.49652	45.7	54.7	,	122.4	724	0.169[0.179]	23.65	1.128 [138]
He	2	4.002602	0.49968	49.9	65.1	(1.937)	94.32	756	0.1249[0.1786]	4.224	1.024 [34.9]
Li	3	6.941	0.43221	54.6	73.4	1.639	82.76	155	0.534		
Be	4	9.012182	0.44384	55.8	75.2	1.594	65.19	35.28	1.848		_
C	6	12.011	0.49954	60.2	86.3	1.745	42.70	18.8	2.265 <sup>e</sup>	00	
$N_2$	7	14.00674	0.49976	61.4	87.8	(1.825)	37.99	47.1	0.8073[1.250]	77.36	1.205 [298]
$O_2$	8	15.9994	0.50002	63.2	91.0	(1.801)	34.24	30.0	1.141[1.428]	90.18	1.22 [296]
$F_2$ Ne	9	18.9984032	0.47372	65.5 66.1	95.3 96.6	(1.675)	32.93	21.85	1.507[1.696]	85.24	[195]
Al	10 13	$20.1797 \\ 26.981539$	0.49555 $0.48181$	$66.1 \\ 70.6$	106.4	(1.724) $1.615$	28.94 $24.01$	24.0 8.9	1.204[0.9005] 2.70	27.09	1.092 [67.1]
Si	14	28.0855	0.49848	70.6	106.4	1.664	24.01 $21.82$	9.36	2.33		3.95
Ar	18	39.948	0.49848 $0.45059$	76.4	100.0 $117.2$	(1.519)	19.55	9.50	1.396[1.782]	87.28	3.93 1.233 [283]
Ti	22	47.867	0.45948	79.9	124.9	1.476	16.17	3.56	4.54	01.20	
Fe	26	55.845	0.46556	82.8	131.9	1.451	13.84	1.76	7.87		_
re Cu	29	63.546	0.46536 $0.45636$	85.6	134.9	1.401 $1.403$	13.84 $12.86$	1.43	8.96		_
Ge	32	72.61	0.44071	88.3	140.5	1.371	12.25	2.30	5.323		_
Sn	50	118.710	0.42120	100.2	163	1.264	8.82	1.21	7.31		_
Xe	54	131.29	0.41130	102.8	169	(1.255)	8.48	2.87	2.953[5.858]	165.1	[701]
W	74	183.84	0.40250	110.3	185	1.145	6.76	0.35	19.3		
Pt	78	195.08	0.39984	113.3	189.7	1.129	6.54	0.305	21.45		_
Pb	82	207.2	0.39575	116.2	194	1.123	6.37	0.56	11.35		_
U	92	238.0289	0.38651	117.0	199	1.082	6.00	$\approx 0.32$	$\approx 18.95$		_
Air, (20°C, 1	atm.), [S	STP]	0.49919	62.0	90.0	(1.815)	36.66	[30420]	(1.205)[1.2931]	78.8	(273) [293]
$H_2O$	// [	,	0.55509	60.1	83.6	1.991	36.08	36.1	1.00	373.15	1.33
$CO_2$ gas			0.49989	62.4	89.7	(1.819)	36.2	[18310]	[1.977]		[410]
$CO_2$ solid (dr			0.49989	62.4	89.7	1.787	36.2	23.2	1.563	${\rm sublimes}$	_
Shielding cond	$\operatorname{crete} f$		0.50274	67.4	99.9	1.711	26.7	10.7	2.5		_
$SiO_2$ (fused qu			0.49926	66.5	97.4	1.699	27.05	12.3	$2.20^{\ g}$		1.458
Dimethyl ethe	$er, (CH_3)$	) <sub>2</sub> O	0.54778	59.4	82.9		38.89			248.7	
$Methane,\ CH_{2}$	4		0.62333	54.8	73.4	(2.417)	46.22	[64850]	0.4224[0.717]	111.7	$[444]_{}$
Ethane, $C_2H_6$			0.59861	55.8	75.7	(2.304)	45.47	[34035]	$0.509(1.356)^{1}$		$(1.038)^{h}$
Propane, $C_3H$	-		0.58962	56.2	76.5	(2.262)	45.20		(1.879)	231.1	_
Isobutane, (C.			0.58496	56.4	77.0	(2.239)	45.07	[16930]	[2.67]	261.42	[1900]
Octane, liquid			0.57778	56.7	77.7	2.123	44.86	63.8	0.703	398.8	1.397
Paraffin wax,		$l_2)_{n\approx 23}CH_3$	0.57275	56.9	78.2	2.087	44.71	48.1	0.93		_
Nylon, type $6$			0.54790	58.5	81.5	1.974	41.84	36.7	1.14		_
Polycarbonate			0.52697	59.5	83.9	1.886	41.46	34.6	1.20		_
Polyethylene t		ate (Mylar) <sup>k</sup>	0.52037	60.2	85.7	1.848	39.95	28.7	1.39		_
Polyethylene <sup>l</sup>			0.57034	57.0	78.4	2.076	44.64	$\approx 47.9$	0.92 – 0.95		_
Polyimide film		on) <sup>m</sup>	0.51264	60.3	85.8	1.820	40.56	28.6	1.42		
Lucite, Plexig		0	0.53937	59.3	83.0	1.929	40.49	≈34.4	1.16-1.20		≈1.49
Polystyrene, s			0.53768	58.5	81.9	1.936	43.72	42.4	1.032		1.581
Polytetrafluor			0.47992	64.2	93.0	1.671	34.84	15.8	2.20		
Polyvinyltolul			0.54155	58.3	81.5	1.956	43.83	42.5	1.032		
Aluminum oxi	\ _	-,	0.49038	67.0	98.9	1.647	27.94	7.04	3.97		1.761
Barium fluorio			0.42207	92.0	145	1.303	9.91	2.05	4.89		1.56
Bismuth germanate (BGO) <sup>r</sup>		0.42065	98.2	157	1.251	7.97	1.12	7.1		2.15	
Cesium iodide			0.41569	102	167	1.243	8.39	1.85	4.53		1.80
Lithium fluori			0.46262	62.2	88.2	1.614	39.25	14.91	2.632		1.392
Sodium fluorio		)	0.47632	66.9 04.6	98.3 151	1.69	29.87	11.68	2.558		1.336
Sodium iodide	\ /		0.42697	94.6	151	1.305	9.49	2.59	3.67		1.775
Silica Aerogel			0.50093	66.3	96.9	1.740	27.25	$136@\rho=0.$			$1.0+0.21\rho$
NEMA G10 p	late '			62.6	90.2	1.87	33.0	19.4	1.7		

Material	Dielectric	Young's	Coeff. of	Specific	Electrical	Thermal conductivity	
	constant $(\kappa = \epsilon/\epsilon_0)$	modulus	thermal	heat	resistivity		
	() is $(\kappa - 1) \times 10^6$	$[10^6 \text{ psi}]$	expansion	[cal/g-°C]	$[\mu\Omega cm(@^{\circ}C)]$	[cal/cm-°C-sec]	
	for gas	,	$[10^{-6} \text{cm/cm-}^{\circ}\text{C}]$	. , , , ,			
$H_2$	(253.9)	_	_	_	_	_	
He	(64)		_	_	_	_	
Li	_	_	56	0.86	$8.55(0^{\circ})$	0.17	
Be	_	37	12.4	0.436	$5.885(0^{\circ})$	0.38	
C	_	0.7	0.6-4.3	0.165	1375(0°)	0.057	
$N_2$	(548.5)		_	_		_	
$O_2$	(495)		_	_	_	_	
Ne	(127)	_	_	_	_	_	
Al	<u> </u>	10	23.9	0.215	$2.65(20^{\circ})$	0.53	
Si	11.9	16	2.8 – 7.3	0.162		0.20	
Ar	(517)	_	_			_	
Ti	_	16.8	8.5	0.126	$50(0^{\circ})$	_	
Fe	_	28.5	11.7	0.11	9.71(20°)	0.18	
Cu	_	16	16.5	0.092	$1.67(20^{\circ})$	0.94	
Ge	16.0	_	5.75	0.073	_	0.14	
$\operatorname{Sn}$	_	6	20	0.052	$11.5(20^{\circ})$	0.16	
Xe	_	_	_			_	
W	_	50	4.4	0.032	$5.5(20^{\circ})$	0.48	
Pt	_	21	8.9	0.032	$9.83(0^{\circ})$	0.17	
Pb	_	2.6	29.3	0.038	$20.65(20^{\circ})$	0.083	
U	_	_	36.1	0.028	$29(20^{\circ})$	0.064	

- R.M. Sternheimer, M.J. Berger, and S.M. Seltzer, Atomic Data and Nuclear Data Tables 30, 261–271 (1984).
- S.M. Seltzer and M.J. Berger, Int. J. Appl. Radiat. 33, 1189–1218 (1982).
- D.E. Groom, N.V. Mokhov, and S.I. Striganov, "Muon stopping-power and range tables," Atomic Data and Nuclear Data Tables 78, 183–356 (2001).
- 4. S.M. Seltzer and M.J. Berger, Int. J. Appl. Radiat. 35, 665 (1984) & http://physics.nist.gov/PhysRefData/Star/Text/contents.html.
- a.  $\sigma_T$ ,  $\lambda_T$  and  $\lambda_I$  are energy dependent. Values quoted apply to high energy range, where energy dependence is weak. Mean free path between collisions  $(\lambda_I)$  or inelastic interactions  $(\lambda_I)$ , calculated from  $\lambda^{-1} = N_A \sum w_j \, \sigma_j \, / A_j$ , where N is Avogadro's number and  $w_j$  is the weight fraction of the jth element in the element, compound, or mixture.  $\sigma_{\text{total}}$  at 80–240 GeV for neutrons ( $\approx \sigma$  for protons) from Murthy et al., Nucl. Phys. **B92**, 269 (1975). This scales approximately as  $A^{0.77}$ .  $\sigma_{\text{inelastic}} = \sigma_{\text{total}} \sigma_{\text{elastic}} \sigma_{\text{quasielastic}}$ ; for neutrons at 60–375 GeV from Roberts et al., Nucl. Phys. **B159**, 56 (1979). For protons and other particles, see Carroll et al., Phys. Lett. **80B**, 319 (1979); note that  $\sigma_I(p) \approx \sigma_I(n)$ .  $\sigma_I$  scales approximately as  $A^{0.71}$ .
- b. For minimum-ionizing muons (results are very slightly different for other particles). Minimum dE/dx from Ref. 3, using density effect correction coefficients from Ref. 1. For electrons and positrons see Ref. 4. Ionization energy loss is discussed in Sec. 26.
- c. From Y.S. Tsai, Rev. Mod. Phys. **46**, 815 (1974);  $X_0$  data for all elements up to uranium are given. Corrections for molecular binding applied for  $H_2$  and  $D_2$ . For atomic H,  $X_0 = 63.05$  g/cm<sup>2</sup>.
- d. For molecular hydrogen (deuterium). For atomic H,  $X_0 = 63.047 \text{ g cm}^{-2}$ .
- e. For pure graphite; industrial graphite density may vary 2.1–2.3 g/cm<sup>3</sup>.
- f. Standard shielding blocks, typical composition  $O_2$  52%, Si 32.5%, Ca 6%, Na 1.5%, Fe 2%, Al 4%, plus reinforcing iron bars. The attenuation length,  $\ell=115\pm5$  g/cm<sup>2</sup>, is also valid for earth (typical  $\rho=2.15$ ), from CERN–LRL–RHEL Shielding exp., UCRL–17841 (1968).
- g. For typical fused quartz. The specific gravity of crystalline quartz is 2.64.
- h. Solid ethane density at  $-60^{\circ}$ C; gaseous refractive index at  $0^{\circ}$ C, 546 mm pressure.
- i. Nylon, Type 6,  $(NH(CH_2)_5CO)_n$
- j. Polycarbonate (Lexan),  $(C_{16}H_{14}O_3)_n$
- k. Polyethylene terephthlate, monomer, C<sub>5</sub>H<sub>4</sub>O<sub>2</sub>
- l. Polyethylene, monomer  $CH_2 = CH_2$
- m. Polymide film (Kapton),  $(C_{22}H_{10}N_2O_5)_n$
- n. Polymethylmethacralate, monomer  $CH_2 = C(CH_3)CO_2CH_3$
- o. Polystyrene, monomer  $\mathrm{C_6H_5CH}{=}\mathrm{CH_2}$
- p. Teflon, monomer  $CF_2 = CF_2$
- q. Polyvinyltolulene, monomer 2-CH<sub>3</sub>C<sub>6</sub>H<sub>4</sub>CH=CH<sub>2</sub>
- r. Bismuth germanate (BGO),  $(Bi_2O_3)_2(GeO_2)_3$
- s. 97% SiO<sub>2</sub> + 3% H<sub>2</sub>O by weight; see A. R. Buzykaev et al., Nucl. Instrum. Methods A433, 396 (1999). Aerogel in the density range  $0.04-0.06 \text{ g/cm}^3$  has been used in Čerenkov counters, but aerogel with higher and lower densities has been produced.  $\rho = \text{density in g/cm}^3$ .
- t. G10-plate, typically  $60\%~\mathrm{SiO_2}$  and 40% epoxy.