

## Group 2 / IIA; Alkaline Earth Metals Group

Calcium (Ca) and magnesium (Mg) are ubiquitous; the rest is rarer; Radium is radioactive and thus only found in radioactive decay chains. Altogether the earth alkali metals account for about 4 % of the crust of the earth. Like the [alkali metals](#) they are very reactive and thus never found in the elemental state.

All are silvery-shining metals that oxidize immediately in air. Stable oxide layers, however, protect some of beryllium (Be) and magnesium (Mg) so they can be used in their metallic state and are stable in water (which dissolves the others)

Magnesium (Mg), calcium (Ca) are technically important (not to mention the [biological](#) importance of calcium). The rest also finds special uses.

*Table of Basic Data*

Name (German)	Beryllium <i>Beryllium</i>	Magnesium <i>Magnesium</i>	Calcium <i>Calcium</i>	Strontium <i>Strontium</i>	Barium <i>Barium</i>	Radium <i>Radium</i>
Atomic number	4	12	20	38	56	88
Atomic mass [u]	9,01	24,31	40,08	87,62	137,33	226,03
Melting point [K]	1551	921,95	1112	1042	998	973
Melting point [°C]	1278	648,95	839	769	725	700
Melting point [°F]	2332	1200	1542	1416	1337	1292
Boiling point [K]	3243	1380	1760	1657	1913	1413
Density [g/cm <sup>3</sup> ]	1,85	1,74	1,54	2,63	3,65	5,50
Ionization energy [eV]	9,32	7,65	6,11	5,70	5,21	5,28
Electro-negativity	1,5	1,2	1,0	1,0	1,0	1,0
Atomic radius [pm]	113,3	160	197,3	215,1	217,3	223
Ionic radius [pm]	34	78	106	127	143	152
Oxidation numbers	2	2	2	2	2	2
lattice typ Transformation temp. [°C]	hcp	hex	hcp 440 fcc	bcc 540 hcp 234 fcc	bcc	?
Lattice constant [Å] (a or c)	2,29	3,20	5,56 ?	6,09 ?	5,02	?
Young's - Modul us [GPa]	310	44,3	19,6	15,7	12,7	?

Therm. expansion coefficient $\alpha$ [ $10^{-6}\text{K}^{-1}$ ]	11,5	26	22	?	19	?
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- In case of doubt all numbers are for room temperatures
- fcc = [face centered cubic](#); lattice const. = a
- bcc = [body centered cubic](#)
- sc = [simple cubic](#)
- hp = simple [hexagonal](#)
- hcp = [hexagonal close packed](#); lattice constants a and c.
- op = [simple orthorhombic](#), [monoclinic](#), [triclinic](#)
- tp = [simple tetragonal](#)
- dia = [diamond structure](#)
- r = [trigonal](#) or rhomboedral trigonal