

TABLE OF PERIODIC PROPERTIES OF THE ELEMENTS

Percent Ionic Character of a Single Chemical Bond

Difference in electronegativity	0.1	0.2	0.3	0.4	0.5	0.6	0.7	0.8	0.9	1.0	1.1	1.2	1.3	1.4	1.5	1.6	1.7	1.8	1.9	2.0	2.1	2.2	2.3	2.4	2.5	2.6	2.7	2.8	2.9	3.0	3.1	3.2
Percent ionic character %	0.5	1	2	4	6	9	12	15	19	22	26	30	34	39	43	47	51	55	59	63	67	70	74	76	79	82	84	86	88	89	91	92

GROUP 1/IA

H	0.32	2.20
	0.79	0.4581
	14.10	0.0585
	13.998	—
	14.304	0.1815

2/IIA

Li	1.23	0.98	0.90	1.57
	2.05	147.1	1.40	297
	13.10	3.0	5.0	11.71
	5.392	11.7	9.323	25
	3.582	84.7	1.825	200

Na	1.54	0.93	1.36	1.31
	2.23	98.01	1.72	127.6
	23.70	2.601	14.0	8.95
	5.139	20.1	17.646	22.4
	1.228	141	1.02	156

DATA CONCERNING THE MORE STABLE ELEMENTARY (SUBATOMIC) PARTICLES

	Neutron	Proton	Electron*	Neutrino†	Photon
Symbol	n	p	e (-)	ν	γ
Rest mass (kg)	1.67495x10 ⁻²⁷	1.67265x10 ⁻²⁷	9.1095x10 ⁻³¹	~0	0
Relative atomic mass (12C=12)	1.008665	1.007276	5.48580x10 ⁻⁴	~0	0
Charge (C)	0	1.60219x10 ⁻¹⁹	-1.60219x10 ⁻¹⁹	0	0
Radius (m)	8x10 ⁻¹⁶	8x10 ⁻¹⁶	<1x10 ⁻¹⁶	~0	0
Spin quantum number	1/2	1/2	1/2	1/2	1
Magnetic Moment†	-1.913 μ _N	2.793 μ _N	1.001 μ _B	0	0

* The positron (e⁺) has properties similar to those of the (negative) electron or beta particle except that its charge has opposite sign (+). The antineutrino (ν̄) has properties similar to those of the neutrino except that its spin (or rotation) is opposite in relation to its direction of propagation.
An antineutrino accompanies release of an electron in radioactive β (particle) decay, whereas a neutrino accompanies the release of a positron in β⁺ decay.
† μ_B=Bohr magneton and μ_N=Nuclear magneton.

18/VIII

He	0.93	—
	0.49	0.084
	31.80	0.021
	24.587	—
	5.193	0.152

3/IIIA 4/IVA 5/VA 6/VIA 7/VIIA 8 VIIIA 9 10 11/IB 12/IIB

K	Ca	Sc	Ti	V	Cr	Mn	Fe	Co	Ni	Cu	Zn	Ga	Ge	As	Se	Br	Kr	
2.03	0.82	1.74	1.00	1.44	1.36	1.32	1.54	1.22	1.63	1.18	1.66	1.17	1.55	1.17	1.83	1.16	1.88	
2.77	76.9	2.23	154.67	2.09	304.80	2.00	425.2	1.92	446.7	1.85	399.5	1.79	219.74	1.72	349.5	1.67	373.3	
45.30	2.33	29.90	6.53	15.0	16.11	10.69	18.6	8.35	22.9	7.23	20	7.39	14.64	7.1	13.8	6.70	16.2	
4.341	16.4	6.113	31.3	5.56	1.5	6.83	2.6	6.75	4.0	4.734	0.5	7.902	11.2	7.88	17.9	7.640	14.6	
0.757	102.5	0.647	200	0.568	15.8	0.523	21.9	0.489	30.7	0.449	93.7	0.479	7.82	0.449	80.2	0.421	100	
Rb	Sr	Y	Zr	Nb	Mo	Tc	Ru	Rh	Pd	Ag	Cd	In	Sn	Sb	Te	I	Xe	
2.16	0.82	1.91	0.95	1.62	1.22	2.16	1.34	1.6	1.30	2.26	1.95	2.2	1.25	2.28	1.28	2.28	1.28	2.28
2.90	87.62	2.45	136.9	2.27	393.20	2.16	590.5	2.08	690.1	1.95	502	1.89	567.77	1.83	495.39	1.79	393.3	
55.9	2.34	33.7	8.2	19.8	17.15	14.1	21	10.80	26.9	9.4	36	8.5	23	8.3	25.52	8.30	21.76	
4.177	47.8	5.695	5.0	6.22	1.8	6.63	2.3	6.76	6.6	7.092	17.3	7.28	0.001	7.36	14.9	7.46	23	
0.363	58.2	0.301	35.3	0.30	17.2	0.278	22.7	0.265	53.7	0.265	138	0.243	50.6	0.238	117	0.243	150	
Cs	Ba	La	Hf	Ta	W	Re	Os	Ir	Pt	Au	Hg	Tl	Pb	Bi	Po	At	Rn	
2.35	0.79	1.98	0.89	1.25	1.10	1.44	1.3	1.34	1.5	1.30	2.36	1.28	1.9	1.26	2.2	1.27	2.20	
3.34	67.740	2.78	140.2	2.74	399.57	2.02	422.58	1.97	707.10	1.89	652.58	1.83	510.45	1.79	324.43	1.76	59.30	
70	2.09	39.0	8.01	22.50	11.30	13.60	21.76	9.53	35.4	8.85	33.05	8.43	29.29	8.54	26.36	9.10	19.66	
3.894	5.3	5.212	2.8	5.58	1.9	6.825	3.4	7.55	8.1	7.864	18.2	8.438	12.3	8.967	21.3	8.959	9.4	
0.24	35.9	0.204	18.4	0.195	13.5	0.144	23.0	0.140	57.5	0.132	174	0.137	47.9	0.130	87.6	0.131	147	
Fr	Ra	Ac	Rf	Db	Sg	Bh	Hs	Mt	Ds	Rg	Uub	Uut	Uuq	Uup	Uuh	Uu	Uu	
0.7	—	0.9	—	—	—	—	—	—	—	—	—	—	—	—	—	—	—	
2.7	64*	136.82	—	—	—	—	—	—	—	—	—	—	—	—	—	—	—	
4.073	2.1*	45.2	8.37	—	—	—	—	—	—	—	—	—	—	—	—	—	—	
—	—	5.279	1.0	5.17	—	—	—	—	—	—	—	—	—	—	—	—	—	
—	—	0.094	18.6	0.120	12	—	—	—	—	—	—	—	—	—	—	—	—	

* ESTIMATED VALUES

13/IIIB 14/IVB 15/VB 16/VIB 17/VII B

Ce	Pr	Nd	Pm	Sm	Eu	Gd	Tb	Dy	Ho	Er	Tm	Yb	Lu
1.65	1.12	1.65	1.13	1.64	1.14	1.63	1.13	1.62	1.17	1.85	1.2	1.61	1.20
2.70	313.80	2.97	332.63	2.94	293.68	2.92	—	2.59	191.63	2.56	175.73	2.54	311.71
5.538	1.4	5.473	1.5	5.525	1.6	5.582	2	5.644	1.1	5.67	1.1	6.15	0.8
0.19	11.4	0.193	12.5	0.190	16.5	—	17.9*	0.197	13.3	0.182	13.9	0.236	10.6
Th	Pa	U	Np	Pu	Am	Cm	Bk	Cf	Es	Fm	Md	No	Lr
1.65	1.3	—	1.5	1.42	1.38	—	1.36	1.08	1.28	—	1.3*	—	1.3*
—	543.92	—	—	—	422.68	—	—	—	—	—	—	—	—
19.90	15.65	15.0	—	12.50	15.48	—	—	12.32	—	—	—	—	—
6.31	7.1	5.89	5.6	6.194	3.6	6.266	0.8	6.026	0.7	5.974	0.7	5.991	—
0.113	54	—	47	0.124	27.6	—	6.3	0.13*	6.74	—	—	—	—

- NOTES: (1) For representative oxides (higher valence) of group. Oxide is acidic if color is red, basic if color is blue and amphoteric if both colors are shown. Intensity of color indicates relative strength.
 (2) Cubic, face centered; Cubic, body centered; Cubic; hexagonal; rhombohedral; tetragonal; orthorhombic; monoclinic.
 (3) At 300 K (27°C) (6) Generally at 293 K (20°C)
 (4) At boiling point for polycrystalline material
 (5) At melting point (7) Quantum mechanical value for free atom
 (8) From density at 300 K (27°C) for liquid and solid elements; values for gaseous elements refer to liquid state at boiling point

The A & B subgroup designations, are those recommended by the International Union of Pure and Applied Chemistry.



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