

PERIODIC TABLE OF THE ELEMENTS

Table of Selected Radioactive Isotopes

Selected Radioactive Isotopes

Naturally occurring radioactive isotopes are designated by a mass number in parentheses. Radioactive isotopes that are not naturally occurring are designated by a mass number in parentheses and a half-life in parentheses. Half-lives follow in parentheses, where s, min, h, d, and y stand respectively for seconds, minutes, hours, days, and years. The table contains information on the decay mode, the decay energy, and the half-life of the isotopes. Isotopes that have been produced in laboratories but are not naturally occurring have been produced in laboratories known to be radioactive. Isotopes that have been produced in laboratories but are not naturally occurring and whose half-lives exceeding 10¹⁰ y have not been included. Symbols designating the principal mode (or modes) of decay are as follows: α - alpha particle emission; β^- - beta minus (electron) emission; β^+ - beta plus (positron) emission; γ - gamma ray emission; ϵ - orbital electron capture; EC - electron capture; SF - spontaneous fission.

GROUP
1/A

1 1.00794
1.00814
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2 4.002602
4.0015061
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3 6.941
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4 9.012182
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5 10.811
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6 12.0107
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7 13.00335
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8 14.00307
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9 15.994915
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11 19.99244
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12 24.30509
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13 26.9815386
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14 27.976927
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15 28.9764151
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16 30.973762
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18 35.453
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19 39.0983
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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	5	10	15	20	25	30	35	40	45	50	55	60	65	70	75	80	85	90	95	100	100	100	100	100	100	100	100	100	100	100	100	100

GROUP 1/IIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
H	1.008	2.20	13.6	1s ¹	Orthorhombic	Gas
Li	6.941	0.98	5.39	1s ² 2s ¹	BCC	Solid
Na	22.990	0.93	5.14	1s ² 2s ² 3s ¹	BCC	Solid
K	39.098	0.82	4.19	1s ² 2s ² 3s ² 4s ¹	BCC	Solid
Rb	85.468	0.79	4.18	1s ² 2s ² 3s ² 4s ² 5s ¹	BCC	Solid
Cs	132.905	0.77	4.19	1s ² 2s ² 3s ² 4s ² 5s ² 6s ¹	BCC	Solid

DATA CONCERNING THE MORE STABLE ELEMENTARY (SUBATOMIC) PARTICLES

Symbol	Neutron	Proton	Electron	Neutrino	Photon
Rest mass (kg)	1.67493x10 ⁻²⁷	1.67262x10 ⁻²⁷	9.10938x10 ⁻³¹	~0	0
Relative atomic mass (m ₀)	1.008665	1.007276	5.4858x10 ⁻⁴	~0	0
Charge (e)	0	1	-1	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 position (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 helicity) 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.

T_{1/2} - Bohr magneton and μ_N - Nuclear magneton.

18/III

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
He	4.0026	2.20	23.72	1s ²	Face-centered cubic	Gas
Ne	20.1797	2.20	21.51	1s ² 2s ² 2p ⁶	Face-centered cubic	Gas
Ar	39.948	2.20	15.21	1s ² 2s ² 2p ⁶ 3s ² 3p ⁶	Face-centered cubic	Gas
Kr	83.801	2.20	14.01	1s ² 2s ² 2p ⁶ 3s ² 3p ⁶ 4s ² 4p ⁶	Face-centered cubic	Gas
Xe	131.29	2.20	12.13	1s ² 2s ² 2p ⁶ 3s ² 3p ⁶ 4s ² 4p ⁶ 5s ² 5p ⁶	Face-centered cubic	Gas

2/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Be	9.0122	1.57	9.00	1s ² 2s ²	Hexagonal	Solid
Mg	24.304	1.31	7.38	1s ² 2s ² 3s ²	Hexagonal	Solid
Ca	40.078	1.00	6.11	1s ² 2s ² 3s ² 4s ²	Face-centered cubic	Solid
Sc	44.956	1.36	6.30	1s ² 2s ² 3s ² 4s ² 3d ¹	Hexagonal	Solid
Ti	47.88	1.54	6.82	1s ² 2s ² 3s ² 4s ² 3d ²	Hexagonal	Solid

* The position (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 helicity) 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.

T_{1/2} - Bohr magneton and μ_N - Nuclear magneton.

3/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Al	26.9815	1.61	5.78	1s ² 2s ² 3s ² 3p ¹	Face-centered cubic	Solid
Ga	69.723	1.61	5.79	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹	Orthorhombic	Solid
In	114.818	1.61	5.79	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹ 5s ² 5p ¹	Orthorhombic	Solid
Tl	204.38	1.61	5.79	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹ 5s ² 5p ¹ 6s ² 6p ¹	Orthorhombic	Solid

4/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Si	28.0855	1.90	8.45	1s ² 2s ² 3s ² 3p ²	Diamond	Solid
Ge	72.630	1.90	8.45	1s ² 2s ² 3s ² 3p ² 4s ² 4p ²	Diamond	Solid
Sn	118.710	1.90	8.45	1s ² 2s ² 3s ² 3p ² 4s ² 4p ² 5s ² 5p ²	Diamond	Solid
Pb	207.2	1.90	8.45	1s ² 2s ² 3s ² 3p ² 4s ² 4p ² 5s ² 5p ² 6s ² 6p ²	Diamond	Solid

5/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
As	74.9216	2.20	9.41	1s ² 2s ² 3s ² 3p ³	Orthorhombic	Solid
Sb	121.757	2.20	9.41	1s ² 2s ² 3s ² 3p ³ 4s ² 4p ³	Orthorhombic	Solid
Bi	208.980	2.20	9.41	1s ² 2s ² 3s ² 3p ³ 4s ² 4p ³ 5s ² 5p ³	Orthorhombic	Solid
Po	209	2.20	9.41	1s ² 2s ² 3s ² 3p ³ 4s ² 4p ³ 5s ² 5p ³ 6s ² 6p ³	Orthorhombic	Solid

6/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Se	78.96	2.55	10.36	1s ² 2s ² 3s ² 3p ⁴	Orthorhombic	Solid
Te	127.6	2.55	10.36	1s ² 2s ² 3s ² 3p ⁴ 4s ² 4p ⁴	Orthorhombic	Solid
Po	209	2.55	10.36	1s ² 2s ² 3s ² 3p ⁴ 4s ² 4p ⁴ 5s ² 5p ⁴	Orthorhombic	Solid

7/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Br	79.904	2.96	11.81	1s ² 2s ² 3s ² 3p ⁵	Orthorhombic	Solid
I	126.905	2.66	10.45	1s ² 2s ² 3s ² 3p ⁵ 4s ² 4p ⁵	Orthorhombic	Solid
At	210	2.66	10.45	1s ² 2s ² 3s ² 3p ⁵ 4s ² 4p ⁵ 5s ² 5p ⁵	Orthorhombic	Solid

8/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Kr	83.801	2.20	14.01	1s ² 2s ² 3s ² 3p ⁶	Face-centered cubic	Gas
Xe	131.29	2.20	12.13	1s ² 2s ² 3s ² 3p ⁶ 4s ² 4p ⁶	Face-centered cubic	Gas
Rn	222	2.20	10.72	1s ² 2s ² 3s ² 3p ⁶ 4s ² 4p ⁶ 5s ² 5p ⁶ 6s ² 6p ⁶	Face-centered cubic	Gas

9/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Ag	107.868	1.93	7.57	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹ 5s ² 5p ¹	Face-centered cubic	Solid
Cu	63.546	1.90	7.73	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹	Face-centered cubic	Solid
Ni	58.693	1.91	7.64	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹ 3d ⁸	Face-centered cubic	Solid

10/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Zn	65.38	1.65	9.00	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹	Hexagonal	Solid
Cd	112.411	1.69	8.99	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹ 5s ² 5p ¹	Hexagonal	Solid
Hg	200.59	1.69	8.99	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹ 5s ² 5p ¹ 6s ² 6p ¹	Hexagonal	Solid

11/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Al	26.9815	1.61	5.78	1s ² 2s ² 3s ² 3p ¹	Face-centered cubic	Solid
Ga	69.723	1.61	5.79	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹	Orthorhombic	Solid
In	114.818	1.61	5.79	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹ 5s ² 5p ¹	Orthorhombic	Solid

12/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Ge	72.630	1.90	8.45	1s ² 2s ² 3s ² 3p ²	Diamond	Solid
Sn	118.710	1.90	8.45	1s ² 2s ² 3s ² 3p ² 4s ² 4p ²	Diamond	Solid
Pb	207.2	1.90	8.45	1s ² 2s ² 3s ² 3p ² 4s ² 4p ² 5s ² 5p ²	Diamond	Solid

13/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
As	74.9216	2.20	9.41	1s ² 2s ² 3s ² 3p ³	Orthorhombic	Solid
Sb	121.757	2.20	9.41	1s ² 2s ² 3s ² 3p ³ 4s ² 4p ³	Orthorhombic	Solid
Bi	208.980	2.20	9.41	1s ² 2s ² 3s ² 3p ³ 4s ² 4p ³ 5s ² 5p ³	Orthorhombic	Solid

14/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Se	78.96	2.55	10.36	1s ² 2s ² 3s ² 3p ⁴	Orthorhombic	Solid
Te	127.6	2.55	10.36	1s ² 2s ² 3s ² 3p ⁴ 4s ² 4p ⁴	Orthorhombic	Solid
Po	209	2.55	10.36	1s ² 2s ² 3s ² 3p ⁴ 4s ² 4p ⁴ 5s ² 5p ⁴	Orthorhombic	Solid

15/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Br	79.904	2.96	11.81	1s ² 2s ² 3s ² 3p ⁵	Orthorhombic	Solid
I	126.905	2.66	10.45	1s ² 2s ² 3s ² 3p ⁵ 4s ² 4p ⁵	Orthorhombic	Solid
At	210	2.66	10.45	1s ² 2s ² 3s ² 3p ⁵ 4s ² 4p ⁵ 5s ² 5p ⁵	Orthorhombic	Solid

16/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Kr	83.801	2.20	14.01	1s ² 2s ² 3s ² 3p ⁶	Face-centered cubic	Gas
Xe	131.29	2.20	12.13	1s ² 2s ² 3s ² 3p ⁶ 4s ² 4p ⁶	Face-centered cubic	Gas
Rn	222	2.20	10.72	1s ² 2s ² 3s ² 3p ⁶ 4s ² 4p ⁶ 5s ² 5p ⁶ 6s ² 6p ⁶	Face-centered cubic	Gas

17/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Ag	107.868	1.93	7.57	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹ 5s ² 5p ¹	Face-centered cubic	Solid
Cu	63.546	1.90	7.73	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹	Face-centered cubic	Solid
Ni	58.693	1.91	7.64	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹ 3d ⁸	Face-centered cubic	Solid

18/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
He	4.0026	2.20	23.72	1s ²	Face-centered cubic	Gas
Ne	20.1797	2.20	21.51	1s ² 2s ² 2p ⁶	Face-centered cubic	Gas
Ar	39.948	2.20	15.21	1s ² 2s ² 2p ⁶ 3s ² 3p ⁶	Face-centered cubic	Gas

19/IIIA

Symbol	Atomic Weight	Electronegativity	Ionization Potential (eV)	Electron Configuration	Crystal Structure	Phase
Al	26.9815	1.61	5.78	1s ² 2s ² 3s ² 3p ¹	Face-centered cubic	Solid
Ga	69.723	1.61	5.79	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹	Orthorhombic	Solid
In	114.818	1.61	5.79	1s ² 2s ² 3s ² 3p ¹ 4s ² 4p ¹ 5s ² 5p ¹	Orthorhombic	Solid

20/IIIA

122	145	133	134	18	130	216	127	19	126
393.30	218	190.05	20.88	289.1	2.01	590.4	1.95	502	149
17.15	14.1	21	10.80	65.9	9.4	36	8.5	23	8.3
1.8	6.63	2.3	6.78	6.6	7.092	17.3	7.28	0.001	7.38
17.2	0.778	22.7	0.285	53.7	0.25	138	0.24	50.6	0.236
									
10	144	134	155	180	226	19	126	19	126