

# silver

|                      |                             |
|----------------------|-----------------------------|
| Other names:         | argentum<br>silver element  |
| Inchi:               | InChI=1S/Ag                 |
| InchiKey:            | BQCADISMDOOEFD-UHFFFAOYSA-N |
| Formula:             | Ag                          |
| SMILES:              | [Ag]                        |
| Mol. weight [g/mol]: | 107.87                      |
| CAS:                 | 7440-22-4                   |

## Physical Properties

| Property code | Value         | Unit    | Source       |
|---------------|---------------|---------|--------------|
| ea            | 1.30 ± 0.20   | eV      | NIST Webbook |
| ea            | 1.30 ± 0.01   | eV      | NIST Webbook |
| ea            | 1.30 ± 0.00   | eV      | NIST Webbook |
| ea            | 1.30 ± 0.03   | eV      | NIST Webbook |
| hf            | 284.90 ± 0.80 | kJ/mol  | NIST Webbook |
| ie            | 7.62 ± 0.07   | eV      | NIST Webbook |
| ie            | 7.57          | eV      | NIST Webbook |
| ie            | 7.50 ± 0.30   | eV      | NIST Webbook |
| ie            | 7.58          | eV      | NIST Webbook |
| ie            | 7.80 ± 0.20   | eV      | NIST Webbook |
| ie            | 7.50          | eV      | NIST Webbook |
| ie            | 7.51 ± 0.07   | eV      | NIST Webbook |
| ie            | 7.58          | eV      | NIST Webbook |
| ie            | 7.53          | eV      | NIST Webbook |
| ie            | 7.58          | eV      | NIST Webbook |
| ie            | 7.58          | eV      | NIST Webbook |
| ie            | 7.57          | eV      | NIST Webbook |
| ie            | 7.60          | eV      | NIST Webbook |
| sgb           | 173.00 ± 0.00 | J/molxK | NIST Webbook |
| ss            | 42.55 ± 0.20  | J/molxK | NIST Webbook |

## Temperature Dependent Properties

| Property code | Value     | Unit | Temperature [K] | Source  |
|---------------|-----------|------|-----------------|---|
| dvisc         | 0.0035200 | Paxs | 1273.00         | A Novel Vibrating Finger Viscometer for High-Temperature Measurements in Liquid Metals and Alloys |
| dvisc         | 0.0029040 | Paxs | 1323.00         | A Novel Vibrating Finger Viscometer for High-Temperature Measurements in Liquid Metals and Alloys |
| dvisc         | 0.0028120 | Paxs | 1373.00         | A Novel Vibrating Finger Viscometer for High-Temperature Measurements in Liquid Metals and Alloys |
| dvisc         | 0.0026490 | Paxs | 1423.00         | A Novel Vibrating Finger Viscometer for High-Temperature Measurements in Liquid Metals and Alloys |
| dvisc         | 0.0025400 | Paxs | 1473.00         | A Novel Vibrating Finger Viscometer for High-Temperature Measurements in Liquid Metals and Alloys |
| dvisc         | 0.0024980 | Paxs | 1523.00         | A Novel Vibrating Finger Viscometer for High-Temperature Measurements in Liquid Metals and Alloys |
| dvisc         | 0.0023990 | Paxs | 1573.00         | A Novel Vibrating Finger Viscometer for High-Temperature Measurements in Liquid Metals and Alloys |
| dvisc         | 0.0023320 | Paxs | 1623.00         | A Novel Vibrating Finger Viscometer for High-Temperature Measurements in Liquid Metals and Alloys |

|       |             |        |         |   |
|-------|-------------|--------|---------|---|
| dvisc | 0.0022140   | Paxs   | 1673.00 | A Novel Vibrating Finger Viscometer for High-Temperature Measurements in Liquid Metals and Alloys                                 |
| econd | 61070000.00 | S/m    | 298.00  | Structural and physical studies of the Ag-rich alloys from Ag-Li system   |
| hvapt | 265.10      | kJ/mol | 1367.00 | Mass spectrometric study of the vaporization behaviour of alpha-Na2NpO4: Thermodynamic investigation of the enthalpy of formation |

Correlations

| Information                 | Value                         |
|-----------------------------|-------------------------------|
| Property code               | pvap                          |
| Equation                    | $\ln(P_{vp}) = A + B/(T + C)$ |
| Coeff. A                    | 1.66364e+01                   |
| Coeff. B                    | -2.82950e+04                  |
| Coeff. C                    | -8.12200e+01                  |
| Temperature range (K), min. | 1283.00                       |
| Temperature range (K), max. | 2453.15                       |

Sources

Calorimetric studies and thermodynamic properties of liquid Ag-Ga alloys: Enthalpies of mixing of liquid Ag-Ga, Au-Ga and Ag-Au-Ga alloys: Thermodynamic properties of liquid Ag-Li alloys: Thermodynamic properties of liquid silver-gold-gallium alloys determined from isopiestic measurements with the vaporization technique: Vaporization behaviour of alpha-Na2NpO4: Properties of liquid silver-gold alloys determined from enthalpies of mixing of Au-Pb and Ag-Au-Pb liquid alloys at 973K: Mixing enthalpies of liquid Bi-Ni and Ag-Bi-Ni alloys: The Yaws Handbook of Vapor Pressure:

<https://www.doi.org/10.1016/j.jct.2014.05.021>  
<https://www.doi.org/10.1016/j.jct.2016.12.020>  
<https://www.doi.org/10.1016/j.jct.2016.06.013>  
<https://www.doi.org/10.1016/j.tca.2018.02.012>  
<https://www.doi.org/10.1016/j.jct.2012.11.034>  
<https://www.doi.org/10.1016/j.tca.2005.02.005>  
<https://www.doi.org/10.1016/j.tca.2007.07.015>  
<https://www.doi.org/10.1016/j.tca.2017.09.002>  
<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

The measurement of thermal conductivity variation with temperature and integral enthalpies of mixing of Ag-Sb liquid alloys: Thermodynamic behavior of polyvalent solutes in solid silver: System Ag Sb: Experimentally determined thermodynamic properties of Ag-Sb system: Ag-Sb phase diagram at 700 K by a semi-automated viscometer for High-Temperature Measurements in Liquid Metals Laboratory: Studies of the Ag-rich alloys from Ag-Li system: KDB:

NIST Webbook:

Experimental investigation and modelling of phase equilibria for the Ag-Cd-In system: Study of the Ag-Sb-S system by the EMF method: Enthalpy of mixing in the Ag-Cd-In ternary liquid phase: Calorimetric measurements and first principles to study the (Ag-Li) liquid system:

<https://www.doi.org/10.1016/j.tca.2012.12.012>  
<https://www.doi.org/10.1016/j.tca.2011.04.032>  
<https://www.doi.org/10.1016/j.tca.2012.09.037>  
<https://www.doi.org/10.1016/j.jct.2013.11.006>  
<https://www.doi.org/10.1016/j.jct.2013.11.031>  
<https://www.doi.org/10.1007/s10765-016-2104-7>  
<https://www.doi.org/10.1016/j.tca.2019.01.016>  
<https://www.thermo.com/research/kdb/hcprop/showprop.php?cmpid=1944>  
<http://webbook.nist.gov/cgi/cbook.cgi?ID=C7440224&Units=SI>  
<https://www.doi.org/10.1016/j.fluid.2016.02.026>  
<https://www.doi.org/10.1016/j.jct.2016.03.009>  
<https://www.doi.org/10.1016/j.jct.2016.12.005>  
<https://www.doi.org/10.1016/j.jct.2014.10.023>

## Legend

|               |  |
|---------------|--|
| <b>dvisc:</b> | Dynamic viscosity                                |
| <b>ea:</b>    | Electron affinity                                |
| <b>econd:</b> | Electrical conductivity                          |
| <b>hf:</b>    | Enthalpy of formation at standard conditions     |
| <b>hvapt:</b> | Enthalpy of vaporization at a given temperature  |
| <b>ie:</b>    | Ionization energy                                |
| <b>pvap:</b>  | Vapor pressure                                   |
| <b>sgb:</b>   | Molar entropy at standard conditions (1 bar)     |
| <b>ss:</b>    | Solid phase molar entropy at standard conditions |

Latest version available from:

<https://www.cheméo.com/cid/18-740-0/silver.pdf>

Generated by Cheméo on 2025-12-23 02:29:32.638606 +0000 UTC m=+6205170.168646653.

Cheméo (<https://www.cheméo.com>) is the biggest free database of chemical and physical data for the process industry.