

# caesium chloride

Other names:	Cesium chloride
Inchi:	InChI=1S/ClH.Cs/h1H;/q;+1/p-1
InchiKey:	AIYUHDOJVVYHVIT-UHFFFAOYSA-M
Formula:	ClCs
SMILES:	[Cl-].[Cs+]
Mol. weight [g/mol]:	168.36
CAS:	7647-17-8

## Physical Properties

Property code	Value	Unit	Source
ea	0.46 ± 0.10	eV	NIST Webbook
hfus	3.80	kJ/mol	Thermodynamic Characterization of the Congruently Melting Cs3CeCl6 Compound
ie	7.40	eV	NIST Webbook
ie	8.80	eV	NIST Webbook
ie	8.30 ± 0.10	eV	NIST Webbook
ie	8.30 ± 0.30	eV	NIST Webbook
ie	8.70 ± 0.10	eV	NIST Webbook
ie	8.50	eV	NIST Webbook
ie	8.30 ± 0.10	eV	NIST Webbook
ie	7.84 ± 0.05	eV	NIST Webbook
tf	918.00	K	Adiabatic compressibility along the two-phase saturation line for the molten (LiF + CsCl) system

## Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
rhos	3074.50	kg/m3	843.00	Density of Crystalline Alkali Chlorides and Their Eutectic Mixtures Near the Melting Point

rhos	3067.60	kg/m3	853.00	Density of Crystalline Alkali Chlorides and Their Eutectic Mixtures Near the Melting Point
rhos	3060.20	kg/m3	863.00	Density of Crystalline Alkali Chlorides and Their Eutectic Mixtures Near the Melting Point
rhos	3052.40	kg/m3	873.00	Density of Crystalline Alkali Chlorides and Their Eutectic Mixtures Near the Melting Point
rhos	3046.00	kg/m3	883.00	Density of Crystalline Alkali Chlorides and Their Eutectic Mixtures Near the Melting Point
rhos	3040.90	kg/m3	893.00	Density of Crystalline Alkali Chlorides and Their Eutectic Mixtures Near the Melting Point
rhos	3039.50	kg/m3	903.00	Density of Crystalline Alkali Chlorides and Their Eutectic Mixtures Near the Melting Point
rhos	3041.80	kg/m3	913.00	Density of Crystalline Alkali Chlorides and Their Eutectic Mixtures Near the Melting Point

## Correlations

Information	Value
Property code	pvap
Equation	$\ln(P_{vp}) = A + B/(T + C)$
Coeff. A	1.58087e+01
Coeff. B	-1.66564e+04
Coeff. C	-8.16900e+01
Temperature range (K), min.	1017.15
Temperature range (K), max.	1570.15

## Sources

Equilibrium Phase Behavior of Water + Propan-1-ol or Propan-2-ol + Cesium Chloride at Compressibilities and 318.15 Two-phase saturation line for the Electrical Conductivity of Electrolytes Found in Natural Waters from (5 to 90) Water Activity and Solubility Measurements and Model Simulation of Aqueous Cesium Chloride Systems of H<sub>2</sub>O and PEG 600 H<sub>2</sub>O Mixtures at 298.15 K of Crystalline Alkali Chlorides and Their Eutectic Mixtures Near the Melting Chemistry of dicesium calcium tetraborate octahydrate: The Yaws Handbook of Vapor Pressure: Partial molar volume of NaCl and CsCl in mixtures of water and methanol by experimental and molecular simulation of the Congruently Melting Cs<sub>3</sub>CeCl<sub>6</sub> Compound specific heat capacity of water and aqueous cesium chloride solutions Measurements and Correlations of the Speed of Sound in Equilibrium of RbCl/CsCl + Dimethyl Sulfoxide (DMSO) and Ternary Thermodynamic Properties of 5, and 60 wt% (DMSO) and activity coefficients in ternary systems at 298.15 K Measurements and Correlation of Solubilities and Solution Thermodynamic Data in dissolving alkali halides formamide + MCl (M = Na, K, Rb and Cs) + Water Systems in the Formamide/N-methylformamide/K<sup>+</sup>/N-methylformamide/depresion of aqueous formamide systems solubility and correlations: Square Molar Volumes and Compressed Electrolyte Solutions of Densities of Methanolic Alkali Halide Salt Solutions by Experiment and Molecular Simulation: Temperature and Density of Water at Super-temperatures, Heat Capacity of H<sub>2</sub>O and H<sub>2</sub>O + NaCl at T from 298.15 and 1100 K and NaCl in H<sub>2</sub>O: Modeling of the NaCl-H<sub>2</sub>O System: Chloride Ion Activity Coefficients of Chloride Ions in Aqueous Solutions of NaCl and H<sub>2</sub>O: Binary Glycol-water (H<sub>2</sub>O)-Ethylene Glycol and Propylene Glycol Ternary Systems: Vapor Pressure Depression Ternary Systems of H<sub>2</sub>O and Alkali Halide Salt Solutions by Experiment and Molecular Simulation: Electrical Conductivity of LiCl-KCl-CsCl Melts: Solubilities, densities and refractive indices for the ternary systems Ethanol-glycols and Phisococams, K, Rb, Cs (to the ternary System (Cocodimethyl Oxidies of (298.15, RbCl 302.29) and (318.15) CsCl + H<sub>2</sub>O) ternary systems from thermodynamic compressibilities of Molten MCl + NdCl<sub>3</sub> Mixtures (M = Rb, Cs and Cs) metric interaction parameters for the Chemodynamics and Phase Equilibrium of the Systems CsCl-MgCl<sub>2</sub> Solubilities, Densities, and Refractive Indices for the Ternary Systems Phase equilibria of 20 systems of RbCl + CsCl 208.29 and 318.15 K: RbCl + HCl Phase equilibria of H<sub>2</sub>O + polyethers glycol of (RbCl)-H<sub>2</sub>O and CsCl-H<sub>2</sub>O: Effect of Alkali Halide Ion on Phase Equilibria: Experimental and Calculated metal acetates:

<https://www.doi.org/10.1021/je050265z>

<https://www.doi.org/10.1016/j.jct.2019.07.003>

<https://www.doi.org/10.1021/je101012n>

<https://www.doi.org/10.1021/acs.jced.7b00459>

<https://www.doi.org/10.1021/acs.jced.7b00024>

<https://www.doi.org/10.1021/ie901030f>

<https://www.doi.gov/10.1016/j.tca.2006.05.020>

<https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure>

<https://www.doi.org/10.1016/j.fluid.2017.10.034>

<https://www.doi.org/10.1021/je8002223>

<https://www.doi.org/10.1016/j.jct.2005.10.010>

<https://www.doi.org/10.1021/je4007986>

<https://www.doj.org/10.1016/j.tca.2008.11.003>

<https://www.doi.org/10.1016/j.ijct.2016.03.027>

<https://www.doi.org/10.1021/acs.iced.5b01043>

<https://www.doi.org/10.1016/j.ijct.2010.10.021>

<https://www.doi.org/10.1016/j.ijct.2014.05.013>

<https://www.doi.org/10.1016/j.ijct.2005.05.005>

<https://www.doi.org/10.1021/acs.jced.7b00690>

<https://www.doi.org/10.1021/je5009944>

<https://www.doi.org/10.1016/j.ijct.2016.09.031>

<https://www.doi.org/10.1016/j.ijct.2004.01.004>

<https://www.doi.org/10.1021/je900630n>

<https://www.doi.org/10.1021/ie400959n>

<https://www.doi.org/10.1021/je500420a>

<http://webbook.nist.gov/cgi/cbook.cgi?ID=C7647178&Units=SI>

<https://www.doi.org/10.1021/acs.jced.5b00683>

<https://www.doi.org/10.1016/j.ijct.2010.01.017>

<https://www.doi.org/10.1021/acs.iced.7b00033>

<https://www.doi.org/10.1016/j.ijct.2016.08.014>

<https://www.doi.org/10.1021/jo100554a>

<https://www.doi.org/10.1016/j.ijet.2004.04.010>

<https://www.doi.org/10.1031/acs.joc.4c00953>

<https://www.doi.org/10.1031/jc300443t>

<https://www.doi.org/10.1016/j.ijet.2013.09.016>

<https://www.doi.org/10.1016/j.ijet.2019.03.020>

<https://www.doi.org/10.1016/j.tsc.2024.11.004>

# Legend

<b>ea:</b>	Electron affinity
<b>hfus:</b>	Enthalpy of fusion at standard conditions
<b>ie:</b>	Ionization energy
<b>pvap:</b>	Vapor pressure
<b>rhos:</b>	Solid Density
<b>tf:</b>	Normal melting (fusion) point

Latest version available from:

<https://www.chemeo.com/cid/41-878-2/caesium-chloride.pdf>

Generated by Cheméo on 2025-12-05 09:06:38.525419426 +0000 UTC m=+4673796.055460090.

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