caesium chloride

Other names:	Cesium chloride
Inchi:	InChI=1S/CIH.Cs/h1H;/q;+1/p-1
InchiKey:	AIYUHDOJVYHVIT-UHFFFAOYSA-M
Formula:	CICs
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 ± 0.10	eV	NIST Webbook
ie	7.84 ± 0.05	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
tf	918.00	К	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(Pvp) = 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

Indices for the Ternary Systems Cherenachemicstry H200 (RPS) Na, calcus, m resign of the second of the se in mixtures of water and methanol by

Water Activity and Solubility

+ Dusone user instanted standard the approximation of the second secon Solutions by Experiment and Molecular Bhase Diagrams and Physicochemical Properties for the Ternary System (Cecino duramines of a empleted of of a gage us description of a empleted of of a gage us description of a empleted of of a gage us description of a empleted of of a gage us description of a empleted of of a gage us description of a empleted of the a description of a empleted of the a state of the state of the a terms of the a terms of a description of the a terms of the a terms of a description of the a terms of the a terms of a description of the a terms of the a terms of a description of the a terms of the a terms of the a description of the a terms of the a terms of the a description of the a terms of the a terms of the a description of the a terms of the a terms of the a description of the a terms of the a terms of the a description of the a terms of the a terms of the a description of the a terms of the a terms of the a description of the a terms of the a terms of the a description of the a terms of the a terms of the a description of the a terms of the a terms of the a description of the a terms of the a terms of the a terms of the a description of the a terms of the a terms of the a terms of the a description of the a terms of the a terms of the a terms of the a description of the a terms of te H2O systems:

Ultrasound velocity in dissolving alkali https://www.doi.org/10.1016/j.jct.2010.10.021 halide melts: Thermodynamic Characterization of the https://www.doi.org/10.1021/je8002223 Thermodynamic Characterization of the Congruently Melting Cs3CeCl6 Consecutivitionia of CsCl-polyethylene dlycol (PEG)-H2O at 298.15 K: Effect of infernation and accountering control and the poly of the second se https://www.doi.org/10.1016/j.tca.2006.05.020 https://www.doi.org/10.1016/j.fluid.2017.10.034 https://www.doi.org/10.1016/j.jct.2016.03.027 **Baseq angli i brit i monometric in ternary systems at By any any internation in ternary systems at By any any internation in ternary systems at By any any internation i brit internation i brit** http://webbook.nist.gov/cgi/cbook.cgi?ID=C7647178&Units=SI https://www.doi.org/10.1021/acs.jced.7b00459 Water Activity and Solubility https://www.doi.org/10.1021/acs.jced.7b00455 Measurements and Model Simulation of Adiabatic Monophysic Score Action line for the Source Verpoints and Model Simulation of Adiabatic Monophysic Score Action line for the Source Verpoints and Verpoints and Source Action and Action a https://www.doi.org/10.1016/j.tca.2008.11.003 https://www.doi.org/10.1016/j.tca.2004.11.004 https://www.doi.org/10.1021/acs.jced.7b00023 https://www.doi.org/10.1016/j.jct.2005.05.005 https://www.doi.org/10.1021/je400959n https://www.doi.org/10.1021/acs.jced.7b00690 https://www.sciencedirect.com/book/9780128029992/the-yaws-handbook-of-vapor-pressure https://www.doi.org/10.1016/j.jct.2010.01.017 https://www.doi.org/10.1016/j.jct.2016.09.031

https://www.doi.org/10.1021/acs.jced.6b00952

Legend

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

Latest version available from:

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

Generated by Cheméo on 2024-04-28 08:25:57.95459971 +0000 UTC m=+16582006.875177026.

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