

B-Li
Cs-Li

77Cal: B. Callmer, "An Accurate Refinement of the β -Rhombohedral Boron Structure," *Acta Crystallogr. B*, **33**, 1951-1954 (1977). (Equi Diagram, Crys Structure; Experimental)

77Sor: V.P. Sorokin, P.I. Gavrilov, and E.V. Levakov, "Preparation of Lithium Monoboride," *Zh. Neorg. Khim.*, **22**(3), 595-597 (1977) in Russian; TR: *Russ. J. Inorg. Chem.*, **22**(3), 329-330 (1977). (Equi Diagram, Crys Structure; Experimental)

78Mit: M.A. Mitchell and R.A. Sutula, "The Density, Electrical Resistivity and Hall Coefficient of Li-B Alloys," *J. Less-Common Met.*, **57**(2), 161-175 (1978). (Equi Diagram, Crys Structure; Experimental)

***78Wan:** F.E. Wang, M.A. Mitchell, R.A. Sutula, and J.R. Holden, "Crystal Structure of a New Compound Li_5B_4 ," *J. Less-Common Met.*, **61**(2), 237-251 (1978). (Equi Diagram, Crys Structure; Experimental)

***79Dal:** S. Dallek, D.W. Ernst, and B.F. Larrick, "Thermal Analysis of Lithium-Boron Alloys," *J. Electrochem. Soc.*, **126**(5), 866-870 (1979). (Equi Diagram, Crys Structure, Thermo; Experimental)

79Dev: L.E. DeVries, L.D. Jackson, and S.D. James, "Structure and Anodic Discharge Behavior of Lithium-Boron Alloys in the LiCl-KCl Eutectic Melt (II)," *J. Electrochem. Soc.*, **126**(6), 993-996 (1979). (Equi Diagram; Experimental)

79Vla: M. Vlasse, R. Naslain, J.S. Kasper, and K. Ploog, "The Crystal Structure of Tetragonal Boron," *J. Less-Common Met.*, **67**, 1-6 (1979). (Equi Diagram, Crys Structure; Experimental)

84Mal: G. Mair, "The Lithium-Boron System," dissertation, Univ. Stuttgart, 99 p (1984) in German. (Equi Diagram; Experimental)

* Indicates key paper.

B-Li evaluation contributed by H. Okamoto, ASM INTERNATIONAL, Metals Park, OH 44073. This work was supported by ASM INTERNATIONAL. Literature searched through early 1987. Dr. Okamoto is ASM/NIST Data Program Category Editor for miscellaneous binary alloys.

The Cs-Li (Cesium-Lithium) System

By C.W. Bale
Ecole Polytechnique de Montréal

Equilibrium Diagram

The assessed Cs-Li phase diagram (Fig. 1) is a simple system that exhibits virtually complete immiscibility in both the liquid and solid states. At higher temperatures, the diagram is dominated by the vapor phase protruding from the Cs side.

The diagram is similar to the K-Li, Li-Na, and Li-Rb systems. The extent of immiscibility is expected to be the most pronounced in this system, because Li and Cs have the smallest and largest alkali-metal atomic diameters, respectively.

As reported by [Shunk] and [Hansen], the Cs-Li system has been investigated by [39Boh] (thermal analyses) and by [62Cue] (solubility measurements). More recently, density measurements of Cs in Li have been reported by [70Nov]. The crystal structures of the pure elements are listed in Table 1.

Table 2 lists data of [62Cue] for the solubility of Cs in Li corresponding to two experimental runs. Applying a regression analysis to the data, the following expression for the solubility of Cs in Li as a function of temperature, in K, can be calculated:

$$\log(\text{at.}\% \text{ Cs}) = 4.43 - 6594/T \quad (\text{Eq } 1)$$

A comparison of the measured and smoothed data is also given in Table 2.

[70Nov] reported density measurements in the range 25 to 1100 °C and calculated a solubility of 7.5 wt.% Cs (42.2 at.% Cs) in Li at 1100 °C. It is unlikely that solubility is extensive at this temperature. At 1100 °C, the vapor pressure of Cs is approximately 11.3 atm, and at 1 atm a Cs-Li liquid alloy with 42.2 at.% Cs would be unstable. [70Nov] did not report the total pressure.

In Fig. 1, the Li(L₂) portion of the phase diagram up to 1100 °C was taken from Eq 1. Transition temperatures were taken from [83Cha], [King1], and [King2]. The

Table 1 Cs-Li Crystal Structure and Lattice Parameter Data

Phase	Composition, at.% Li	Pearson symbol	Space group	Strukturbericht		Lattice parameters, nm		Comment
				designation	Prototype	a	c	
(Cs).....	0.00	cI2	Im $\bar{3}m$	A2	W	0.6141	...	25 °C
(α Li).....	100.00	hP2	P6 ₃ /mmc	A3	Mg	0.3111	0.5093	< 75 K
(β Li).....	100.00	cI2	Im $\bar{3}m$	A2	W	0.35093	...	25 °C

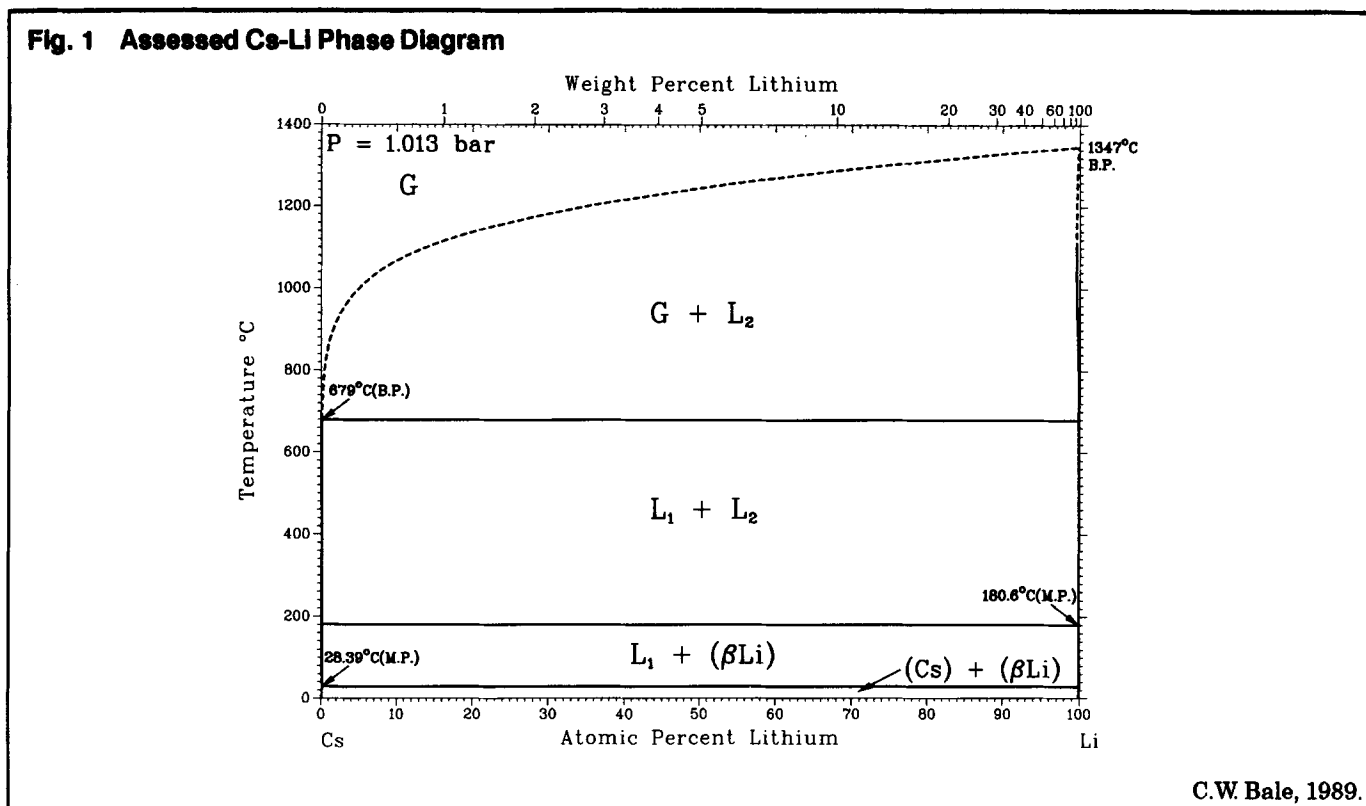


Table 2 Solubility of Cs in Liquid Li

Temperature, °C	Composition, at.% Cs [62Cue](a)	Composition, at.% Cs [62Cue](b)	Eq 1
1093.....	0.336	0.704	0.403
1038.....	0.159	0.200	0.253
982.....	0.174	0.179	0.151
760.....	0.007	0.018	0.011
181.....	10^{-10}

(a) First experimental run. (b) Second experimental run.

vapor phase boundary was calculated by assuming an ideal gas with $P(\text{total}) = P(\text{Li}) + P(\text{Li}_2) + P(\text{Cs}) + P(\text{Cs}_2) = 1 \text{ atm}$ in equilibrium with the liquid where $\alpha_{\text{Li}}(\text{L}) = 1$. Thermodynamic data for the gas phase calculation were taken from the compilation of [73Bar].

Cited References

- 39Boh:** B. Böhm, W. Klemm, "The Behavior of Alkali Metals Toward One Another," *Z. Anorg. Chem.*, **243**, 69 (1939). (Equi Diagram; Experimental)
- 62Cue:** T.R. Cuerou and F. Tepper, ARS Preprint No. 2537-62, (1962). (Equi Diagram; Experimental)
- 70Nov:** I.I. Novikov, V.V. Roshchupkin, Y.S. Trelin, and I.K. Pavlov, "Solubility of Cesium in Lithium in a Fused State," *Teplofiz., Svoistva Zhidk., Mater. Vses. Teplofiz. Konf. Svoistvam Veshchestv Vys. Temp., 3rd*, A.K. Abas-Zade, Ed., 106-110 (1970). (Equi Diagram)
- 73Bar:** I. Barin, O. Knacke, and O. Kubaschewski, *Thermochemical Properties of Inorganic Substances*, Springer-Verlag, Dusseldorf (1973). (Thermo; Compilation)
- 83Cha:** M.W. Chase, "Heats of Transition of the Elements," *Bull. Alloy Phase Diagrams*, **4**(1), 124 (1983). (Equi Diagram; Compilation)

Cs-Li evaluation contributed by C.W. Bale, Centre de Recherche en Calcul Thermochimique, Ecole Polytechnique, Campus de l'Université de Montréal, P.O. Box 6079, Station A, Montréal, Québec, Canada H3C 3A7. This work was partially supported by Department of Energy funds through the Joint Program on Critical Compilation of Physical and Chemical Data coordinated through the Office of Standard Reference Data, National Institute of Standards and Technology. Literature searched through 1988. Professor Bale is the ASM/NIST Data Program Co-Category Editor for binary alkali metal systems.