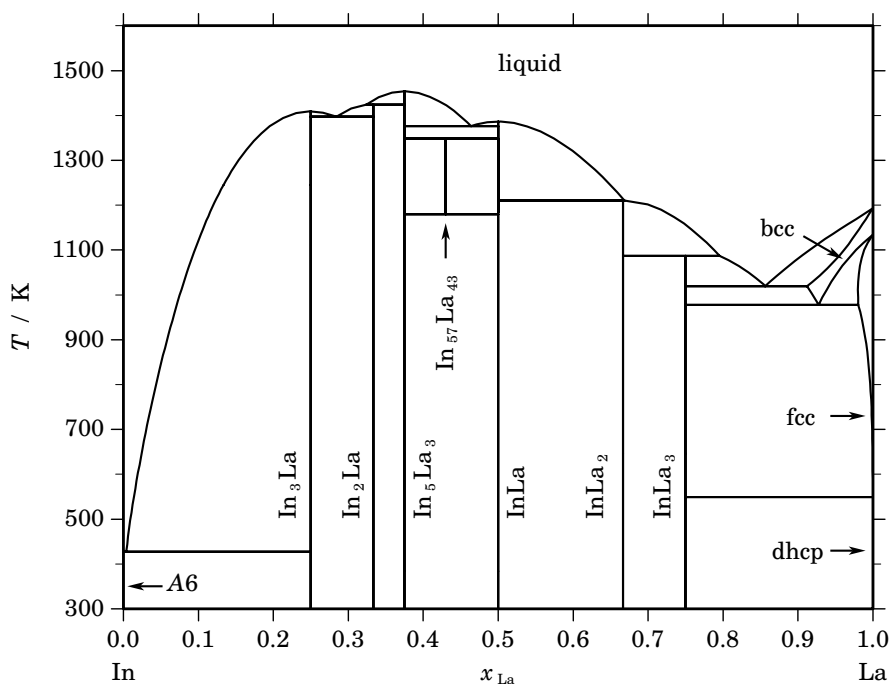


**In – La** (Indium – Lanthanum)**Fig. 1.** Calculated phase diagram for the system In-La.

Most of the interest in the In-La system has probably been attracted by the superconducting compound  $\text{InLa}_3$  which has one of the highest transition temperatures among the superconducting intermetallics of lanthanum. Variations in the reported transition temperatures between 8 and 10 K have been attributed to impurities and prompted for investigations on the solubility range and the phase diagram. The literature on the In-La system has been reviewed by [1992Pal] and a thermodynamic dataset has been optimised in [2002Wei] which is presented here. The phase diagram for the In-La system has been essentially worked out by [1974McM] and in addition the liquidus of In-rich melts with up to 10 at.% La has been investigated by [1971Deg]. Seven intermetallic compounds have been established and only for  $\text{In}_5\text{La}_3$  a small range of solid solubility has been indicated but not quantified. Therefore, the assessment [2002Wei] describes all of them as stoichiometric. Enthalpies of formation have been measured by [1977Bor] across the whole composition range. For  $\text{In}_3\text{La}$  several other measurements of the enthalpy of formation have been reported and also EMF measurements of the Gibbs energy of formation [1971Deg].

**Table I.** Phases, structures and models.

Phase	Struktur- bericht	Prototype	Pearson symbol	Space group	SGTE name	Model
liquid					LIQUID	(In,La) <sub>1</sub>
A6	A6	In	<i>tI2</i>	<i>I4/mmm</i>	TETRAGONAL_A6	In <sub>1</sub>
In <sub>3</sub> La	<i>L1</i> <sub>2</sub>	AuCu <sub>3</sub>	<i>cP4</i>	<i>Pm3̄m</i>	IN3LA	In <sub>3</sub> La <sub>1</sub>
In <sub>2</sub> La	...	CeCu <sub>2</sub>	<i>oI12</i>	<i>Imma</i>	IN2LA	In <sub>2</sub> La <sub>1</sub>
In <sub>5</sub> La <sub>3</sub>	...	Pd <sub>5</sub> Pu <sub>3</sub>	<i>oC32</i>	<i>Cmcm</i>	IN5LA3	In <sub>5</sub> La <sub>3</sub>
In <sub>57</sub> La <sub>43</sub>	...	...	...	...	IN57LA43	In <sub>57</sub> La <sub>43</sub>
InLa	<i>B2</i>	CsCl	<i>cP2</i>	<i>Pm3̄m</i>	INLA	In <sub>1</sub> La <sub>1</sub>
InLa <sub>2</sub>	<i>B8</i> <sub>2</sub>	InNi <sub>2</sub>	<i>hP6</i>	<i>P6<sub>3</sub>/mmc</i>	INLA2	In <sub>1</sub> La <sub>2</sub>
InLa <sub>3</sub>	<i>L1</i> <sub>2</sub>	AuCu <sub>3</sub>	<i>cP4</i>	<i>Pm3̄m</i>	INLA3	In <sub>1</sub> La <sub>3</sub>
bcc	<i>A2</i>	W	<i>cI2</i>	<i>Im3̄m</i>	BCC_A2	(In,La) <sub>1</sub>
fcc	<i>A1</i>	Cu	<i>cF4</i>	<i>Fm3̄3m</i>	FCC_A1	(In,La) <sub>1</sub>
dhcp	<i>A3'</i>	αLa	<i>hP4</i>	<i>P6<sub>3</sub>/mmc</i>	DHCP	(In,La) <sub>1</sub>

**Table II.** Invariant reactions.

Reaction	Type	<i>T</i> / K	Compositions / <i>x</i> <sub>La</sub>			$\Delta_r H$ / (J/mol)
liquid $\rightleftharpoons$ In <sub>5</sub> La <sub>3</sub>	congruent	1453.8	0.375	0.375		–21411
liquid + In <sub>5</sub> La <sub>3</sub> $\rightleftharpoons$ In <sub>2</sub> La	peritectic	1424.3	0.324	0.375	0.333	–18701
liquid $\rightleftharpoons$ In <sub>3</sub> La	congruent	1409.0	0.250	0.250		–26980
liquid $\rightleftharpoons$ In <sub>3</sub> La + In <sub>2</sub> La	eutectic	1397.4	0.284	0.250	0.333	–24731
liquid $\rightleftharpoons$ InLa	congruent	1386.6	0.500	0.500		–23309
liquid $\rightleftharpoons$ In <sub>5</sub> La <sub>3</sub> + InLa	eutectic	1376.3	0.464	0.375	0.500	–22165
In <sub>5</sub> La <sub>3</sub> + InLa $\rightleftharpoons$ In <sub>57</sub> La <sub>43</sub>	peritectoid	1348.9	0.375	0.500	0.430	0
InLa + liquid $\rightleftharpoons$ InLa <sub>2</sub>	peritectic	1210.7	0.500	0.668	0.667	–12261
In <sub>57</sub> La <sub>43</sub> $\rightleftharpoons$ In <sub>5</sub> La <sub>3</sub> + InLa	eutectoid	1179.4	0.430	0.375	0.500	0
InLa <sub>2</sub> + liquid $\rightleftharpoons$ InLa <sub>3</sub>	peritectic	1086.7	0.667	0.796	0.750	–5978
liquid $\rightleftharpoons$ InLa <sub>3</sub> + bcc	eutectic	1019.5	0.856	0.750	0.912	–6203
bcc $\rightleftharpoons$ InLa <sub>3</sub> + fcc	eutectoid	977.9	0.927	0.750	0.980	–2560
fcc $\rightleftharpoons$ InLa <sub>3</sub> + dhcp	eutectoid	549.4	1.000	0.750	1.000	–369
liquid $\rightleftharpoons$ A6 + In <sub>3</sub> La	eutectic	427.4	0.004	0.000	0.250	–3377

**Table IIIa.** Integral quantities for the liquid phase at 1500 K.

$x_{\text{La}}$	$\Delta G_{\text{m}}$ [J/mol]	$\Delta H_{\text{m}}$ [J/mol]	$\Delta S_{\text{m}}$ [J/(mol·K)]	$G_{\text{m}}^{\text{E}}$ [J/mol]	$S_{\text{m}}^{\text{E}}$ [J/(mol·K)]	$\Delta C_P$ [J/(mol·K)]
0.000	0	0	0.000	0	0.000	0.000
0.100	–15954	–18215	–1.508	–11899	–4.210	0.000
0.200	–26035	–30946	–3.274	–19794	–7.435	0.000
0.300	–31814	–38732	–4.612	–24195	–9.691	0.000
0.400	–34006	–42112	–5.404	–25612	–10.999	0.000
0.500	–33200	–41623	–5.615	–24555	–11.378	0.000
0.600	–29927	–37804	–5.251	–21533	–10.847	0.000
0.700	–24676	–31194	–4.345	–17057	–9.424	0.000
0.800	–17877	–22331	–2.969	–11636	–7.130	0.000
0.900	–9835	–11753	–1.279	–5780	–3.982	0.000
1.000	0	0	0.000	0	0.000	0.000

Reference states: In(liquid), La(liquid)

**Table IIIb.** Partial quantities for In in the liquid phase at 1500 K.

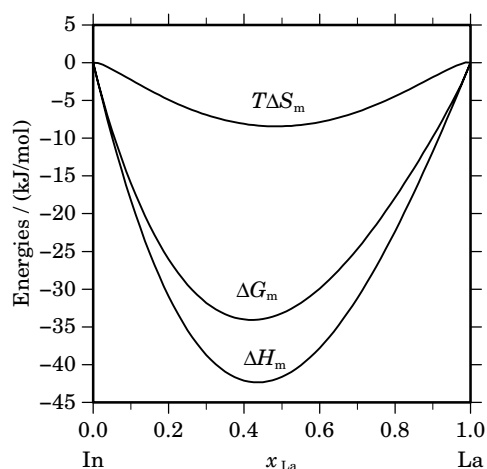
$x_{\text{In}}$	$\Delta G_{\text{In}}$ [J/mol]	$\Delta H_{\text{In}}$ [J/mol]	$\Delta S_{\text{In}}$ [J/(mol·K)]	$G_{\text{In}}^{\text{E}}$ [J/mol]	$S_{\text{In}}^{\text{E}}$ [J/(mol·K)]	$a_{\text{In}}$	$\gamma_{\text{In}}$
1.000	0	0	0.000	0	0.000	1.000	1.000
0.900	–3401	–2832	0.380	–2087	–0.496	0.761	0.846
0.800	–10451	–10608	–0.105	–7668	–1.960	0.433	0.541
0.700	–20172	–22253	–1.388	–15723	–4.353	0.198	0.283
0.600	–31604	–36690	–3.390	–25233	–7.638	0.079	0.132
0.500	–43823	–52841	–6.012	–35178	–11.775	0.030	0.060
0.400	–55965	–69629	–9.109	–44537	–16.728	0.011	0.028
0.300	–67308	–85978	–12.447	–52292	–22.457	0.005	0.015
0.200	–77495	–100810	–15.544	–57422	–28.925	0.002	0.010
0.100	–87625	–113049	–16.950	–58908	–36.094	0.001	0.009
0.000	– $\infty$	–121618	$\infty$	–55729	–43.926	0.000	0.011

Reference state: In(liquid)

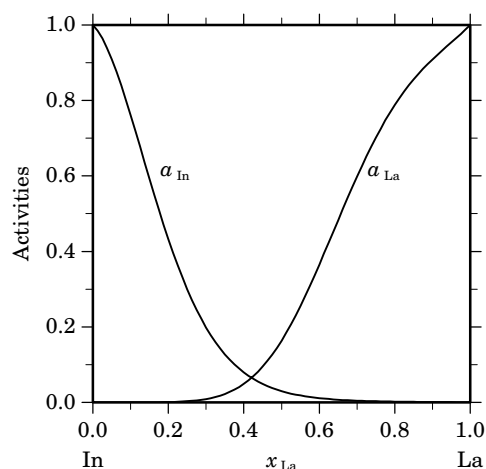
**Table IIIc.** Partial quantities for La in the liquid phase at 1500 K.

$x_{\text{La}}$	$\Delta G_{\text{La}}$ [J/mol]	$\Delta H_{\text{La}}$ [J/mol]	$\Delta S_{\text{La}}$ [J/(mol·K)]	$G_{\text{La}}^{\text{E}}$ [J/mol]	$S_{\text{La}}^{\text{E}}$ [J/(mol·K)]	$a_{\text{La}}$	$\gamma_{\text{La}}$
0.000	– $\infty$	–211362	$\infty$	–140711	–47.101	0.000	0.000
0.100	–128926	–156665	–18.492	–100209	–37.637	0.000	0.000
0.200	–88372	–112297	–15.950	–68300	–29.332	0.001	0.004
0.300	–58979	–77183	–12.136	–43964	–22.146	0.009	0.029
0.400	–37609	–50244	–8.423	–26181	–16.042	0.049	0.123
0.500	–22577	–30405	–5.218	–13932	–10.981	0.164	0.327
0.600	–12568	–16587	–2.679	–6197	–6.927	0.365	0.608
0.700	–6405	–7715	–0.873	–1956	–3.839	0.598	0.855
0.800	–2973	–2711	0.174	–190	–1.681	0.788	0.985
0.900	–1191	–498	0.462	123	–0.414	0.909	1.010
1.000	0	0	0.000	0	0.000	1.000	1.000

Reference state: La(liquid)



**Fig. 2.** Integral quantities of the liquid phase at  $T=1500$  K.



**Fig. 3.** Activities in the liquid phase at  $T=1500$  K.

**Table IV.** Standard reaction quantities at 298.15 K for the compounds per mole of atoms.

Compound	$x_{La}$	$\Delta_f G^\circ$ / (J/mol)	$\Delta_f H^\circ$ / (J/mol)	$\Delta_f S^\circ$ / (J/(mol·K))	$\Delta_f C_P^\circ$ / (J/(mol·K))
In <sub>3</sub> La <sub>1</sub>	0.250	−52657	−57092	−14.876	0.000
In <sub>2</sub> La <sub>1</sub>	0.333	−53128	−56759	−12.181	0.000
In <sub>5</sub> La <sub>3</sub>	0.375	−53323	−56714	−11.374	0.000
In <sub>57</sub> La <sub>43</sub>	0.430	−53434	−57105	−12.314	−0.002
In <sub>1</sub> La <sub>1</sub>	0.500	−53577	−57607	−13.518	0.000
In <sub>1</sub> La <sub>2</sub>	0.667	−35950	−37692	−5.843	0.000
In <sub>1</sub> La <sub>3</sub>	0.750	−27021	−28087	−3.574	0.000

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