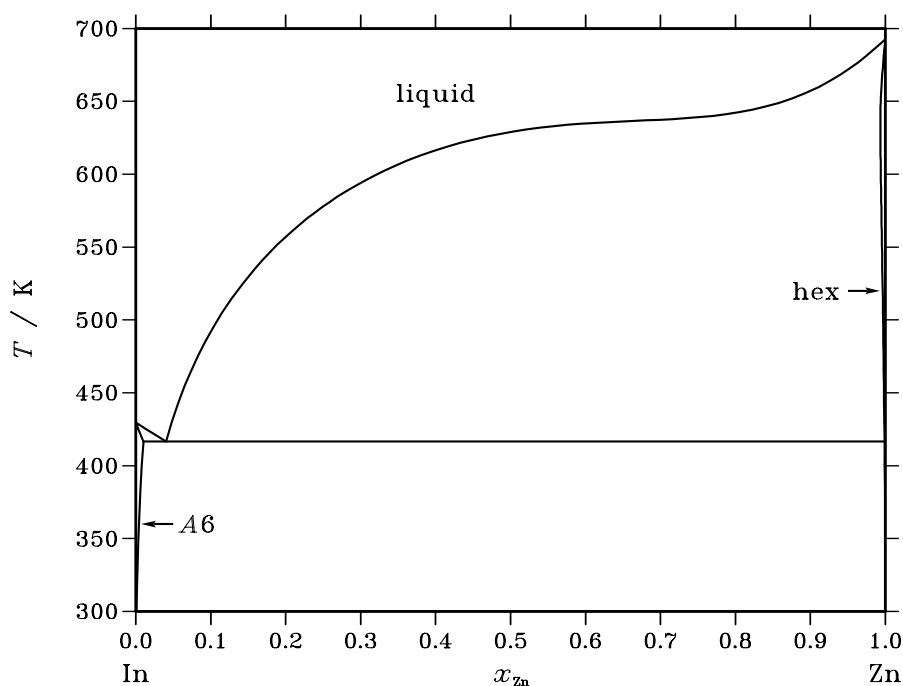


In – Zn (Indium – Zinc)**Fig. 1.** Calculated phase diagram for the system In-Zn.

A good understanding of the thermodynamics and phase equilibria in the In-Zn system is currently very important because of the drive towards the development of Pb-free solder materials for environmental and health reasons. The critically assessed data for this system adopted by SGTE were taken from the assessment of Lee [96Lee]. The experimental information had also been reviewed by Dutkiewicz and Zakulski [84Dut]. The phase diagram for the system is characterised by a simple eutectic close in composition to pure In. The studies of the liquidus surface are in fairly good agreement and shows a tendency towards a metastable miscibility gap in the liquid phase. There seems to be some uncertainty about the exact composition of the eutectic. The value predicted from Lee's assessment is about 4 at.% Zn. There is limited solubility in each of the terminal solid solutions, Zn dissolving in tetragonal In up to about 1.0 at.%. Thermodynamic properties have been determined by calorimetry, EMF studies, vapour pressure measurements and use of an isopiestic technique. The assessed data of Lee provide excellent agreement between calculated and experimental properties.

Table I. Phases, structures and models.

Phase	Strukturbericht	Prototype	Pearson symbol	Space group	SGTE name	Model
liquid					LIQUID	(In,Zn) ₁
A6	A6	In	<i>tI2</i>	<i>I4/mmm</i>	TETRAGONAL_A6	(In,Zn) ₁
hex	A3	Mg	<i>hP2</i>	<i>P6₃/mmc</i>	HCP_ZN	(In,Zn) ₁

Table II. Invariant reactions.

Reaction	Type	<i>T</i> / K	Compositions / <i>x</i> _{Zn}			$\Delta_r H$ / (J/mol)
liquid \rightleftharpoons A6 + hex	eutectic	416.8	0.040	0.010	0.999	–3647

Table IIIa. Integral quantities for the liquid phase at 730 K.

x_{Zn}	ΔG_{m} [J/mol]	ΔH_{m} [J/mol]	ΔS_{m} [J/(mol·K)]	G_{m}^{E} [J/mol]	S_{m}^{E} [J/(mol·K)]	ΔC_P [J/(mol·K)]
0.000	0	0	0.000	0	0.000	0.000
0.100	–1247	926	2.977	726	0.274	0.000
0.200	–1717	1717	4.704	1321	0.543	0.000
0.300	–1923	2359	5.866	1785	0.787	0.000
0.400	–1973	2830	6.579	2112	0.984	0.000
0.500	–1919	3100	6.876	2288	1.112	0.000
0.600	–1790	3136	6.748	2295	1.152	0.000
0.700	–1603	2895	6.161	2105	1.082	0.000
0.800	–1351	2329	5.041	1686	0.881	0.000
0.900	–973	1385	3.230	1000	0.527	0.000
1.000	0	0	0.000	0	0.000	0.000

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

Table IIIb. Partial quantities for In in the liquid phase at 730 K.

x_{In}	ΔG_{In} [J/mol]	ΔH_{In} [J/mol]	ΔS_{In} [J/(mol·K)]	G_{In}^{E} [J/mol]	S_{In}^{E} [J/(mol·K)]	a_{In}	γ_{In}
1.000	0	0	0.000	0	0.000	1.000	1.000
0.900	–574	65	0.875	66	–0.001	0.910	1.011
0.800	–1094	278	1.879	260	0.023	0.835	1.044
0.700	–1571	678	3.082	593	0.116	0.772	1.103
0.600	–2006	1327	4.566	1094	0.319	0.719	1.198
0.500	–2395	2304	6.437	1812	0.673	0.674	1.348
0.400	–2746	3708	8.841	2815	1.223	0.636	1.590
0.300	–3116	5658	12.019	4192	2.008	0.598	1.995
0.200	–3719	8292	16.454	6049	3.073	0.542	2.709
0.100	–5461	11769	23.603	8515	4.458	0.407	4.067
0.000	– ∞	16266	∞	11736	6.206	0.000	6.914

Reference state: In(liquid)

Table IIIc. Partial quantities for Zn in the liquid phase at 730 K.

x_{Zn}	ΔG_{Zn} [J/mol]	ΔH_{Zn} [J/mol]	ΔS_{Zn} [J/(mol·K)]	G_{Zn}^{E} [J/mol]	S_{Zn}^{E} [J/(mol·K)]	a_{Zn}	γ_{Zn}
0.000	– ∞	9894	∞	7928	2.694	0.000	3.692
0.100	–7312	8672	21.896	6664	2.751	0.300	2.998
0.200	–4207	7477	16.005	5562	2.623	0.500	2.500
0.300	–2743	6282	12.363	4565	2.352	0.636	2.121
0.400	–1924	5084	9.600	3638	1.981	0.728	1.821
0.500	–1443	3897	7.315	2764	1.551	0.788	1.577
0.600	–1153	2754	5.353	1947	1.105	0.827	1.378
0.700	–954	1710	3.651	1210	0.685	0.854	1.221
0.800	–759	838	2.188	596	0.333	0.882	1.103
0.900	–475	231	0.966	165	0.090	0.925	1.028
1.000	0	0	0.000	0	0.000	1.000	1.000

Reference state: Zn(liquid)

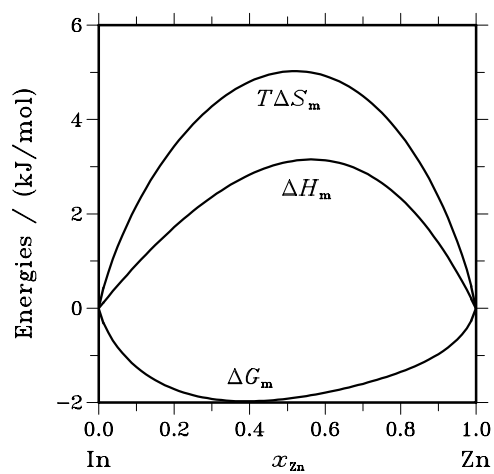


Fig. 2. Integral quantities of the liquid phase at $T=730$ K.

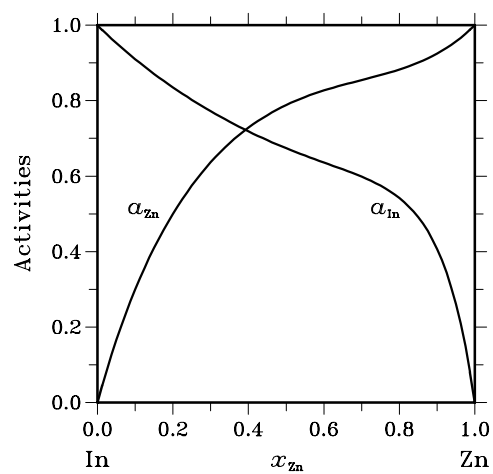


Fig. 3. Activities in the liquid phase at $T=730$ K.

References

- [84Dut] J. Dutkiewicz, W. Zakulski: Bull. Alloy Phase Diagrams **5** (1984) 284–289.
 [96Lee] B.-J. Lee: Calphad **20** (1996) 471–480.