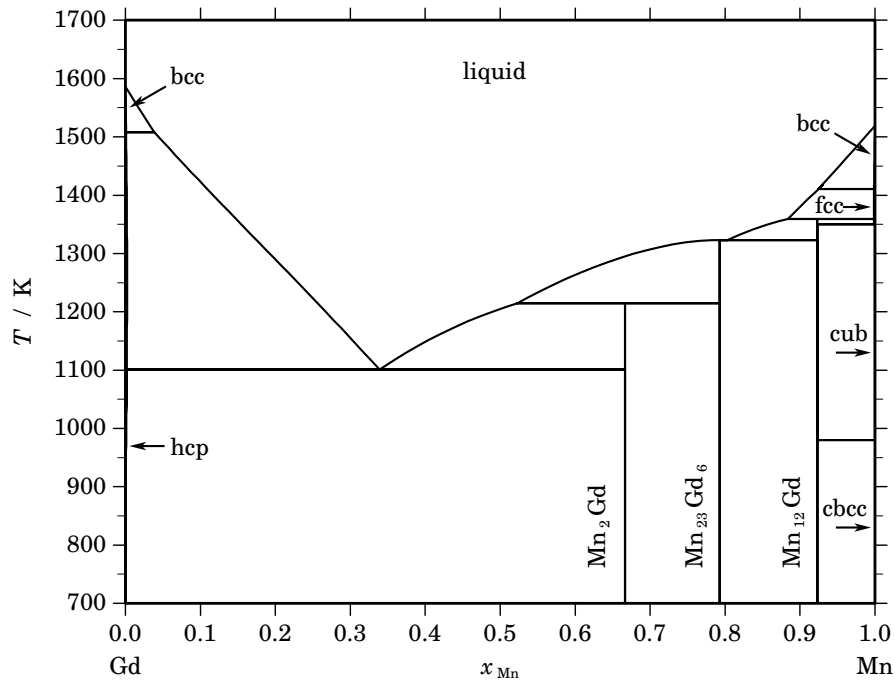


Gd – Mn (Gadolinium – Manganese)**Fig. 1.** Calculated phase diagram for the system Gd-Mn.

Manganese and rare earth metals are added to magnesium alloys in order to improve their creep resistance and strength. A review and a thermodynamic optimisation of the Gd-Mn system has been prepared by [2001Grö]. The phase diagram of the system Gd-Mn has been determined by Kirchmayr and Lugscheider [1967Kir]. The partial enthalpies of the components in molten Gd-Mn alloys have been measured calorimetrically by Nikolaenko and Nosova [1989Nik]. However, their data have a large range of scatter and they are not used in the optimisation [2001Grö]. Enthalpies of formation of the intermetallic compounds have been estimated from the corresponding values of the Mn-Y system which been assessed in the same publication [2001Grö].

Table I. Phases, structures and models.

Phase	Strukturbericht	Prototype	Pearson symbol	Space group	SGTE name	Model
liquid					LIQUID	(Gd,Mn) ₁
bcc	A2	W	<i>cI2</i>	<i>Im</i> $\bar{3}m$	BCC_A2	(Gd,Mn) ₁
hcp	A3	Mg	<i>hP2</i>	<i>P6</i> ₃ / <i>mmc</i>	HCP_A3	(Gd,Mn) ₁
Mn ₂ Gd	C15	Cu ₂ Mg	<i>cF24</i>	<i>Fd</i> $\bar{3}m$	M2R	Mn ₂ Gd ₁
Mn ₂₃ Gd ₆	D8 _a	Mn ₂₃ Th ₆	<i>cF116</i>	<i>Fm</i> $\bar{3}m$	M23R6	Mn ₂₃ Gd ₆
Mn ₁₂ Gd	D2 _b	Mn ₁₂ Th	<i>tI26</i>	<i>I4</i> / <i>mmm</i>	M12R	Mn ₁₂ Gd ₁
fcc	A1	Cu	<i>cF4</i>	<i>Fm</i> $\bar{3}m$	FCC_A1	(Gd,Mn) ₁
cbcc	A12	α Mn	<i>cI58</i>	<i>I</i> $\bar{4}3m$	CBCC_A12	Mn ₁
cub	A13	β Mn	<i>cP20</i>	<i>P4</i> ₁ 32	CUB_A13	Mn ₁

Table II. Invariant reactions.

Reaction	Type	T / K	Compositions / x_{Mn}			$\Delta_r H / (\text{J/mol})$
$\text{bcc} \rightleftharpoons \text{hcp} + \text{liquid}$	metatectic	1507.9	0.001	0.001	0.039	–3502
$\text{bcc} \rightleftharpoons \text{liquid} + \text{fcc}$	metatectic	1410.8	0.999	0.925	0.999	–1905
$\text{liquid} + \text{fcc} \rightleftharpoons \text{Mn}_{12}\text{Gd}$	peritectic	1359.4	0.884	0.998	0.923	–15890
$\text{fcc} \rightleftharpoons \text{Mn}_{12}\text{Gd} + \text{cub}$	eutectoid	1350.2	0.999	0.923	1.000	–2330
$\text{liquid} \rightleftharpoons \text{Mn}_{23}\text{Gd}_6$	congruent	1323.2	0.793	0.793		–21254
$\text{liquid} \rightleftharpoons \text{Mn}_{23}\text{Gd}_6 + \text{Mn}_{12}\text{Gd}$	eutectic	1323.0	0.803	0.793	0.923	–21198
$\text{liquid} + \text{Mn}_{23}\text{Gd}_6 \rightleftharpoons \text{Mn}_2\text{Gd}$	peritectic	1214.7	0.523	0.793	0.667	–8918
$\text{liquid} \rightleftharpoons \text{hcp} + \text{Mn}_2\text{Gd}$	eutectic	1100.8	0.339	0.002	0.667	–15053
$\text{cub} \rightleftharpoons \text{Mn}_{12}\text{Gd} + \text{cbcc}$	eutectoid	980.0	1.000	0.923	1.000	–2254

Table IIIa. Integral quantities for the liquid phase at 1600 K.

x_{Mn}	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	–4044	–1159	1.803	281	–0.900	0.000
0.200	–6186	–2089	2.561	471	–1.600	0.000
0.300	–7545	–2779	2.979	581	–2.100	0.000
0.400	–8331	–3218	3.196	622	–2.400	0.000
0.500	–8617	–3396	3.263	604	–2.500	0.000
0.600	–8416	–3303	3.196	537	–2.400	0.000
0.700	–7693	–2927	2.979	433	–2.100	0.000
0.800	–6355	–2258	2.561	302	–1.600	0.000
0.900	–4171	–1286	1.803	154	–0.900	0.000
1.000	0	0	0.000	0	0.000	0.000

Reference states: Gd(liquid), Mn(liquid)

Table IIIb. Partial quantities for Gd in the liquid phase at 1600 K.

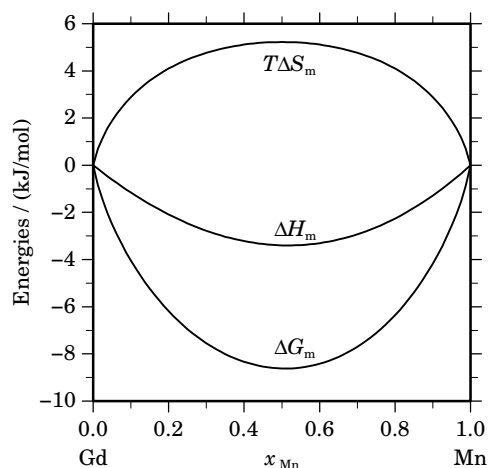
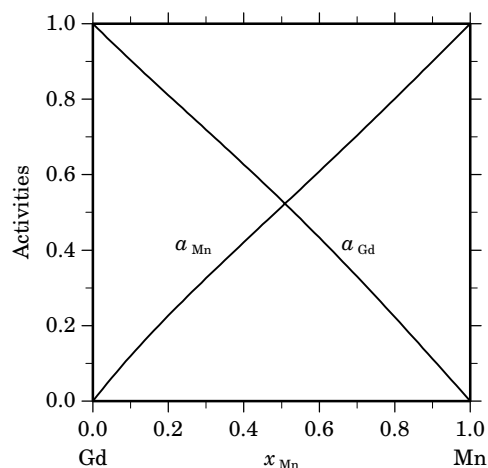
x_{Gd}	ΔG_{Gd} [J/mol]	ΔH_{Gd} [J/mol]	ΔS_{Gd} [J/(mol·K)]	G_{Gd}^{E} [J/mol]	S_{Gd}^{E} [J/(mol·K)]	a_{Gd}	γ_{Gd}
1.000	0	0	0.000	0	0.000	1.000	1.000
0.900	–1355	–113	0.776	47	–0.100	0.903	1.004
0.800	–2794	–466	1.455	174	–0.400	0.811	1.013
0.700	–4385	–1080	2.066	360	–0.900	0.719	1.027
0.600	–6211	–1976	2.647	584	–1.600	0.627	1.045
0.500	–8397	–3176	3.263	825	–2.500	0.532	1.064
0.400	–11130	–4700	4.019	1060	–3.600	0.433	1.083
0.300	–14747	–6570	5.110	1270	–4.900	0.330	1.100
0.200	–19978	–8807	6.982	1433	–6.400	0.223	1.114
0.100	–29105	–11433	11.045	1527	–8.100	0.112	1.122
0.000	– ∞	–14468	∞	1532	–10.000	0.000	1.122

Reference state: Gd(liquid)

Table IIIc. Partial quantities for Mn in the liquid phase at 1600 K.

x_{Mn}	ΔG_{Mn} [J/mol]	ΔH_{Mn} [J/mol]	ΔS_{Mn} [J/(mol·K)]	G_{Mn}^{E} [J/mol]	S_{Mn}^{E} [J/(mol·K)]	a_{Mn}	γ_{Mn}
0.000	$-\infty$	−12702	∞	3298	−10.000	0.000	1.281
0.100	−28247	−10575	11.045	2385	−8.100	0.120	1.196
0.200	−19752	−8581	6.982	1659	−6.400	0.227	1.133
0.300	−14920	−6743	5.110	1097	−4.900	0.326	1.086
0.400	−11511	−5081	4.019	679	−3.600	0.421	1.052
0.500	−8838	−3617	3.263	383	−2.500	0.515	1.029
0.600	−6607	−2371	2.647	189	−1.600	0.609	1.014
0.700	−4671	−1366	2.066	74	−0.900	0.704	1.006
0.800	−2950	−621	1.455	19	−0.400	0.801	1.001
0.900	−1400	−159	0.776	1	−0.100	0.900	1.000
1.000	0	0	0.000	0	0.000	1.000	1.000

Reference state: Mn(liquid)

**Fig. 2.** Integral quantities of the liquid phase at $T=1600$ K.**Fig. 3.** Activities in the liquid phase at $T=1600$ K.**Table IV.** Standard reaction quantities at 298.15 K for the compounds per mole of atoms.

Compound	x_{Mn}	$\Delta_f G^\circ$ / (J/mol)	$\Delta_f H^\circ$ / (J/mol)	$\Delta_f S^\circ$ / (J/(mol·K))	$\Delta_f C_P^\circ$ / (J/(mol·K))
Mn ₂ Gd ₁	0.667	−6111	−6408	−0.996	−4.329
Mn ₂₃ Gd ₆	0.793	−6026	−6371	−1.158	−2.689
Mn ₁₂ Gd ₁	0.923	−3720	−4027	−1.031	−1.003

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 [2001Grö] J. Gröbner, A. Pisch, R. Schmid-Fetzer: J. Alloys Comp. **317-318** (2001) 433–437.