

## Erratum to: Phase Stability of Low-Density, Multiprincipal Component Alloys Containing Aluminum, Magnesium, and Lithium

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Due to calculation error, Table II, Table V, Fig. 5, and Fig. 6 in the paper were not presented correctly. The updated versions of Table II, Table V, Fig. 5, and Fig. 6 are shown below.

**Table II. Theoretical ( $\rho_{\text{theor}}$ ) and measured densities ( $\rho_{\text{exp}}$ ), and chemical compositions (at.%) of phases identified in the microstructures of AlLiMgZnSn, AlLi<sub>0.5</sub>MgZn<sub>0.5</sub>Sn<sub>0.2</sub>, AlLi<sub>0.5</sub>MgZn<sub>0.5</sub>Cu<sub>0.2</sub>, AlLi<sub>0.5</sub>MgCu<sub>0.5</sub>Sn<sub>0.2</sub>, Al<sub>80</sub>Li<sub>5</sub>Mg<sub>5</sub>Zn<sub>5</sub>Sn<sub>5</sub>, and Al<sub>80</sub>Li<sub>5</sub>Mg<sub>5</sub>Zn<sub>5</sub>Cu<sub>5</sub> alloys**

<b>Alloys</b>	$\rho_{\text{theor}}$ (g/cm <sup>3</sup> )	$\rho_{\text{exp}}$ (g/cm <sup>3</sup> )	<b>Regions</b>	<b>Measured chemical composition (at.%)</b>				
				<b>Al</b>	<b>Mg</b>	<b>Zn</b>	<b>Sn</b>	<b>Cu</b>
AlLiMgZnSn	3.88	4.23	A	0	62.8	1.79	35.41	—
			B	54.65	0	44.59	0.76	—
			C	89.44	0	10.56	0	—
			D	0	0	11.31	88.69	—
AlLi <sub>0.5</sub> MgZn <sub>0.5</sub> Sn <sub>0.2</sub>	2.90	3.22	A	36.25	35.51	28.23	0	—
			B	0	58.93	3.02	38.04	—
			C	83.45	12.75	3.79	0	—
AlLi <sub>0.5</sub> MgZn <sub>0.5</sub> Cu <sub>0.2</sub>	2.75	3.73	A	39.26	37.79	16.34	—	6.61
			B	32.87	32.46	15.70	—	18.97
			C	70.36	20.12	8.77	—	0.74
AlLi <sub>0.5</sub> MgCu <sub>0.5</sub> Sn <sub>0.2</sub>	2.96	3.69	A	49.05	34.06	—	0	16.88
			B	36.76	35.09	—	0	28.15
			C	4.30	22.12	—	73.58	0
Al <sub>80</sub> Li <sub>5</sub> Mg <sub>5</sub> Zn <sub>5</sub> Sn <sub>5</sub>	3.05	3.05	A	87.83	1.47	3.89	6.81	—
			B	8.20	2.47	3.95	85.38	—
Al <sub>80</sub> Li <sub>5</sub> Mg <sub>5</sub> Zn <sub>5</sub> Cu <sub>5</sub>	2.91	3.08	A	94.54	1.88	2.38	—	1.20
			B	72.23	0	1.87	—	25.91

**Table V. Complete list of investigated alloy compositions, and calculated values for  $\Omega$ ,  $\delta$ ,  $\Delta\chi$  and VEC**

<b>Alloys</b>	<b>Major phases</b>	$\Delta H_{\text{mix}}$ (kJ/mol)	$\Delta S_{\text{mix}}$ (J/mol K)	$\Omega$	$\delta$	$\Delta\chi$	VEC
<i>Initial compositions</i>							
AlLiMgZnSn	Mg <sub>2</sub> Sn/Li <sub>2</sub> MgSn + Zn + Al + Sn	-6.08	13.38	1.54	5.39	0.33	4.40
AlLi <sub>0.5</sub> MgZn <sub>0.5</sub> Sn <sub>0.2</sub>	Mg <sub>2</sub> Sn/Li <sub>2</sub> MgSn + Mg <sub>32</sub> (AlZn) <sub>49</sub> + Al	-3.89	12.31	2.50	5.66	0.27	3.84
AlLi <sub>0.5</sub> MgZn <sub>0.5</sub> Cu <sub>0.2</sub>	Mg <sub>32</sub> (AlZn) <sub>49</sub> + Unknown	-3.30	12.31	3.15	6.72	0.26	4.28
AlLi <sub>0.5</sub> MgCu <sub>0.5</sub> Sn <sub>0.2</sub>	AlCuMg + Mg <sub>2</sub> Sn/Li <sub>2</sub> MgSn + Sn	-3.65	12.31	3.01	7.60	0.31	3.69
Al <sub>80</sub> Li <sub>5</sub> Mg <sub>5</sub> Zn <sub>5</sub> Sn <sub>5</sub>	Al + Mg <sub>2</sub> Sn/Li <sub>2</sub> MgSn + Sn	-0.53	6.47	10.68	3.61	0.17	3.35
Al <sub>80</sub> Li <sub>5</sub> Mg <sub>5</sub> Zn <sub>5</sub> Cu <sub>5</sub>	Al + Al <sub>2</sub> Cu + AlCu <sub>3</sub>	-1.14	6.47	5.21	4.10	0.17	3.70
<i>Additional compositions investigated for phase content only</i>							
(Al <sub>0.5</sub> Mg <sub>0.5</sub> ) <sub>95</sub> Li <sub>5</sub>	Al <sub>12</sub> Mg <sub>17</sub> + LiMg + Unknown phase	-2.19	7.13	2.95	5.50	0.18	2.43
(Al <sub>0.5</sub> Mg <sub>0.5</sub> ) <sub>90</sub> Li <sub>10</sub>	Al <sub>12</sub> Mg <sub>17</sub> + LiMg + Unknown phase	-2.34	7.89	2.97	5.38	0.20	2.35
(Al <sub>0.5</sub> Mg <sub>0.5</sub> ) <sub>85</sub> Li <sub>15</sub>	Al <sub>12</sub> Mg <sub>17</sub> + LiMg + Unknown phase	-2.47	8.41	2.92	5.26	0.22	2.28
(Al <sub>0.5</sub> Mg <sub>0.5</sub> ) <sub>75</sub> Li <sub>25</sub>	Al <sub>12</sub> Mg <sub>17</sub> + LiMg + Unknown phase	-2.63	9.00	2.77	4.99	0.25	2.13
AlMgLi	Al <sub>12</sub> Mg <sub>17</sub> + LiMg + Unknown phase	-2.67	9.13	2.64	4.74	0.26	2.00
Al <sub>40</sub> Mg <sub>40</sub> Li <sub>10</sub> Cu <sub>10</sub>	Al <sub>12</sub> Mg <sub>17</sub> + Mg <sub>32</sub> Al <sub>47</sub> Cu <sub>7</sub> + AlCuMg	-2.76	9.92	3.32	7.06	0.25	3.20
Al <sub>35</sub> Mg <sub>35</sub> Li <sub>15</sub> Cu <sub>15</sub>	Al <sub>12</sub> Mg <sub>17</sub> + Mg <sub>32</sub> Al <sub>47</sub> Cu <sub>7</sub> + AlCuMg	-3.11	10.84	3.21	7.62	0.28	3.55
Al <sub>40</sub> Mg <sub>40</sub> Li <sub>10</sub> Zn <sub>10</sub>	Mg <sub>32</sub> (AlZn) <sub>49</sub> + HCP	-2.68	9.92	3.17	5.74	0.21	3.30
Al <sub>35</sub> Mg <sub>35</sub> Li <sub>15</sub> Zn <sub>15</sub>	Mg <sub>32</sub> (AlZn) <sub>49</sub> + HCP	-3.08	10.84	2.89	5.79	0.23	3.70
AlLi <sub>0.4</sub> MgZn <sub>0.5</sub> Sn <sub>0.2</sub>	Mg <sub>2</sub> Sn/Li <sub>2</sub> MgSn + Mg <sub>32</sub> (AlZn) <sub>49</sub> + Al	-3.68	12.18	2.65	5.71	0.26	3.94
Al <sub>35</sub> Mg <sub>35</sub> Li <sub>15</sub> Cu <sub>10</sub> La <sub>5</sub>	Al <sub>7</sub> CuLa <sub>2</sub> + AlCuMg + Mg <sub>32</sub> Al <sub>47</sub> Cu <sub>7</sub> + Al <sub>12</sub> Mg <sub>17</sub> + Li-rich phase	-6.07	11.64	1.75	8.73	0.27	3.15
Al <sub>35</sub> Mg <sub>35</sub> Li <sub>15</sub> Cu <sub>10</sub> Ce <sub>5</sub>	Al <sub>3</sub> CeCu + AlCuMg + Mg <sub>32</sub> Al <sub>47</sub> Cu <sub>7</sub> + Al <sub>12</sub> Mg <sub>17</sub> + Li-rich phase	-6.04	11.64	1.75	8.23	0.27	3.15
Al <sub>35</sub> Mg <sub>35</sub> Li <sub>15</sub> Zn <sub>10</sub> La <sub>5</sub>	Al <sub>2</sub> LaZn <sub>2</sub> + Al <sub>12</sub> Mg <sub>17</sub> + Mg-rich phase	-6.25	11.64	1.58	7.59	0.24	3.25
Al <sub>35</sub> Mg <sub>35</sub> Li <sub>15</sub> Zn <sub>10</sub> Ce <sub>5</sub>	Al <sub>4</sub> Ce + Mg <sub>32</sub> (AlZn) <sub>49</sub> + Mg-rich phase	-6.22	11.64	1.57	7.03	0.23	3.25

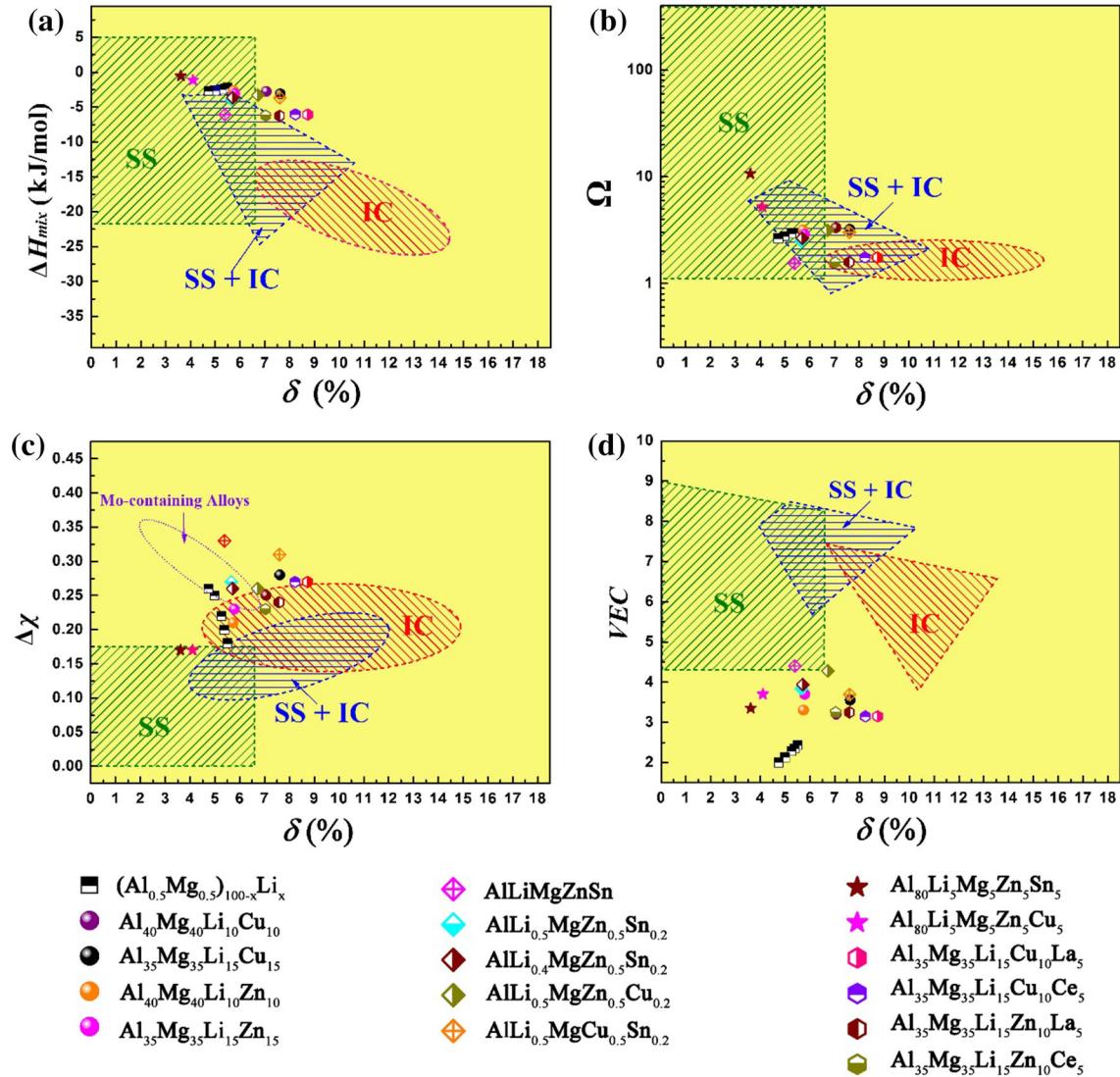


Fig. 5. Phase constituent prediction maps: (a)  $\delta - \Delta H_{\text{mix}}$ ; (b)  $\delta - \Omega$ ; (c)  $\delta - \Delta \chi$ ; and (d)  $\delta - \text{VEC}$  plots for multiprincipal component alloys in this work overlaid on cross-hatched regions developed in previous HEA investigations. (For  $(\text{Al}_{0.5}\text{Mg}_{0.5})_{100-x}\text{Li}_x$ ,  $x = 5, 10, 15, 25$  and  $33.33$ ).

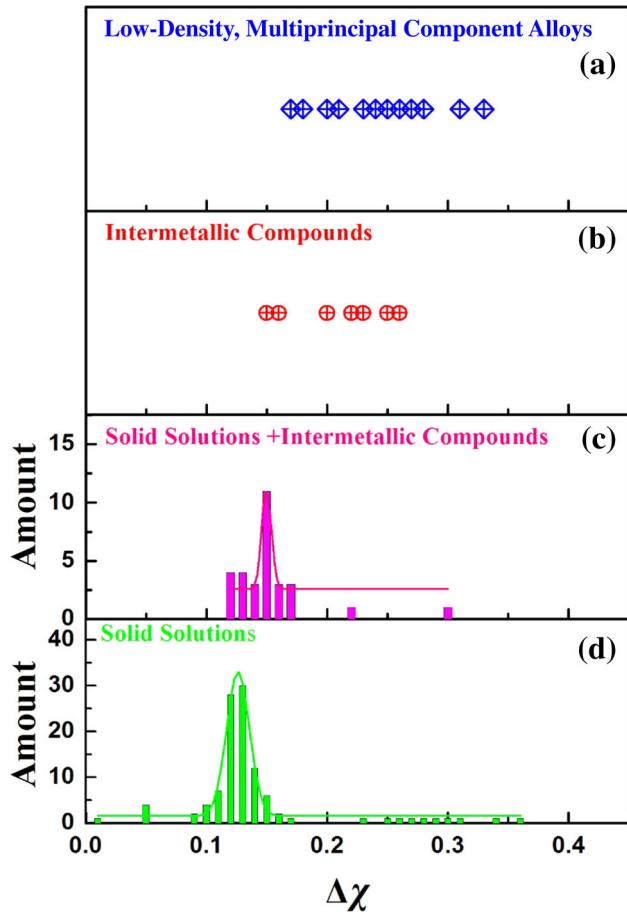


Fig. 6. Values and frequency distributions (c and d) of electronegativity difference ( $\Delta\chi$ ) for: (a) low-density, multiprincipal component alloys (this work); (b) intermetallic compounds; (c) both solid solutions and intermetallic compounds; and (d) solid solution phases.