Applied Physics A (2022) 128:601
https://doi.org/10.1007/s00339-022-05629-3

Applied Physics A
Materials Science & Processing

CORRECTION

Correction to: Structural and crystallographic characterization of grain boundaries coarse particles in an Al–Mg-Si alloy, using convergent beam electron diffraction

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Published online: 22 June 2022

Correction to: Applied Physics A (2022) 128:386
https://doi.org/10.1007/s00339-022-05518-9

The email address of the third author, A. Redjaïmia is forgotten.

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- In the abstract line 5 and page 2 of 10 the Sect. 2 s paragraph line 9 from the up: The sentence should read:

…transmission electron microscopy using convergent beam electron diffraction (CBED) in microdiffraction mode…

- The presentation of Table 1, Table 3, and Table 4 was incorrect. The correct Table 1, Table 3, and Table 4 are given below.
- Page 2 of 10, Sect. 2 line 1 from the bottom: The sentence should read:

Al–Mg-Si.

- Page 4 of 10, third paragraph line 3 from the up: Should read:

Tables II and VII.

- Page 4 of 10, third paragraph line 6 from the up the Ideal symmetries should be underlined and written without spaces: Should read:

The ideal symmetries {(2mm); 2mm} and {(6); 3} of the ZOLZ and the WP, …

- Page 4 of 10, third paragraph line 9 from the up: Should read:

… group $m\bar{3}$ or $2/m\bar{3}$ in ….

- Page 5 of 10, first column line 2 from the bottom: Should read:

… Fig. 6b …

- Page 5 of 10, second column line 1 from the up: Should read:

… Fig. 6a …

- Page 6 of 10, in the caption, to Fig. 5 the Ideal symmetries should be underlined and written without spaces. The correct version is given below.

Figure 5 Microdiffraction patterns from GBCP recorded along two zone axes a) ⟨001⟩ and b) ⟨111⟩, respectively. These two diffraction patterns exhibit the following symmetry: a) {(4mm); (2mm)}; {4mm; 2mm} ZOLZ and WP symmetries along ⟨001⟩ zone axis, b) {(6mm), (6)}; {3m3} ZOLZ and WP symmetries along ⟨111⟩ zone axis

The original article can be found online at https://doi.org/10.1007/s00339-022-05518-9.

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Table 1 Orientation relationships (OR) variants between the cubic α − Al(Mn, Fe)Si phase and the aluminum matrix (m)

Phase	*OR*	Ref
$\alpha - Al(Mn,Fe)Si$	$\left[1\bar{1}1\right]_{\alpha}//\left[1\bar{1}1\right]_{m}$	[43]
(AA 3003 Alloy)	$\left(5\overline{27}\right)_{\alpha}//(011)_{m}$	[44]
$\alpha - Al(Mn,Fe)Si$	$\left[1\bar{1}1\right]_{\alpha}//\left[1\bar{1}1\right]_{m}$	[45]
(3xxx Alloy)	$\left(5\overline{27}\right)_{\alpha}//(011)_{m}$	
$\alpha - Al(Mn,Fe)Si$	$\left[11\bar{1}\right]_{\alpha}//\left[11\bar{1}\right]_{m}$	
(DC-cast AA 3003 Alloy)	$(527)_{\alpha}//(011)_{m}$	
	$\left[\bar{2}50\right]_{\alpha}//\left[11\bar{1}\right]_{m}$	
	$(527)_{\alpha}//(011)_{m}$	

Table 3 Microdifraction data for symmetry analysis

	ZOLZ (Zero Order Laue Zone)		WP (Whole pattern)	
Zone axis	Net symmetry	Ideal symmetry	Net symmetry	Ideal symmetry
⟨001⟩	(4*mm*)	<u>2*mm*</u>	4*mm*	<u>2*mm*</u>
⟨111⟩	(6*mm*)	<u>6</u>	3*m*	<u>3</u>

Table 4 Determination of the intersection point group H whose elements are common to the GBCP and the aluminium matrix, the number of variants n which is the ratio between the orders of the point groups of the matrix and the point group H

Orientation relationships	Superimposed symmetries	Shared symmetries	$H_h = G^{m\bar{3}m}_{m=48} \cap G^{m\bar{3}}_{24}$	$n = \frac{m}{h}$
$\left(\bar{2}\boldsymbol{OO}\right)_{\boldsymbol{GBCP}}//\left(\bar{2}20\right)_{\boldsymbol{M}}$	$2/m$ on $2/m$	$2/m$	$m\bar{3}m \cap \boldsymbol{m}\bar{3} = 2/m$	$48/4 = 12$
$(\boldsymbol{O2O})_{\boldsymbol{GBCP}}//\left(\overline{11}1\right)_{\boldsymbol{M}}$	$2/m$ on $\bar{3}$	1		
$[001]_{\boldsymbol{GBCP}}//[112]_{\boldsymbol{M}}$	$2/m$ on 1	1		

- The presentation of pinacoid writing was incorrect on; Page 7 of 10, second column line 2 from the bottom, page 8 of 10 first column line 7 from the up, and page 9 of 10 in the conclusion: The correct version is: {(010) or (*h*0*l*)}.