

# Erratum: “A complete set of material properties of single domain 0.26Pb(In<sub>1/2</sub>Nb<sub>1/2</sub>)O<sub>3</sub>–0.46Pb(Mg<sub>1/3</sub>Nb<sub>2/3</sub>)O<sub>3</sub>–0.28PbTiO<sub>3</sub> single crystals” [Appl. Phys. Lett. 96, 012907 (2010)]

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In our publication<sup>1</sup> the reference frame corresponds to a reference frame rotated around the z-axis by 180° as described in Ref. 2. To avoid confusion, we provide a revised Table II for the complete data set based on the standard coordinate system (the 0° rotation one in Ref. 2).

Figures 2 and 3 were recalculated based on the standard coordinate system. The maximum values of  $d_{33}^*$ ,  $k_{33}^*$ ,  $\epsilon_{33}^*$ , and  $s_{33}^*$  occur at 60.7°, 57.8°, 90°, and 51.5°, respectively, from the poling direction [111]<sub>c</sub>.

TABLE II. (In standard coordinates) Measured and derived material properties of 0.26Pb(In<sub>1/2</sub>Nb<sub>1/2</sub>)O<sub>3</sub>–0.46Pb(Mg<sub>1/3</sub>Nb<sub>2/3</sub>)O<sub>3</sub>–0.28PbTiO<sub>3</sub> single-domain single crystal poled in [111]<sub>c</sub> (density:  $\rho=8102$  kg/m<sup>3</sup>).

Elastic stiffness constants, $c_{ij}$ (10 <sup>10</sup> N/m <sup>2</sup> )															
$c_{11}^E$ <sup>a</sup>	$c_{12}^E$	$c_{13}^E$	$c_{14}^E$	$c_{33}^E$	$c_{44}^E$ <sup>a</sup>	$c_{66}^E$ <sup>a</sup>	$c_{11}^D$	$c_{12}^D$	$c_{13}^D$	$c_{14}^D$	$c_{33}^D$ <sup>a</sup>	$c_{44}^D$	$c_{66}^D$		
20.96	8.29	6.46	-2.66	17.65	2.10	6.33	22.01	8.25	5.61	-1.08	19.81	6.68	6.88		
Elastic compliance constants, $s_{ij}$ (10 <sup>-12</sup> m <sup>2</sup> /N)															
$s_{11}^E$ <sup>a</sup>	$s_{12}^E$	$s_{13}^E$	$s_{14}^E$	$s_{33}^E$	$s_{44}^E$	$s_{66}^E$	$s_{11}^D$	$s_{12}^D$	$s_{13}^D$	$s_{14}^D$	$s_{33}^D$ <sup>a</sup>	$s_{44}^D$	$s_{66}^D$		
10.43	-6.40	-1.49	21.38	6.76	101.85	33.75	5.58	-1.87	-1.09	1.21	5.88	15.36	14.91		
Piezoelectric coefficients, $e_{i\lambda}$ (C/m <sup>2</sup> )				$d_{i\lambda}$ (10 <sup>-12</sup> C/N)				$g_{i\lambda}$ (10 <sup>-3</sup> Vm/N)				$h_{i\lambda}$ (10 <sup>8</sup> V/m)			
$e_{15}$	$e_{22}$	$e_{31}$	$e_{33}$	$d_{15}$	$d_{22}$	$d_{31}$	$d_{33}$	$g_{15}$	$g_{22}$	$g_{31}$	$g_{33}$	$h_{15}$	$h_{22}$	$h_{31}$	$h_{33}$
18.78	-6.48	-5.19	8.72	2190	-511	-34	74	3.93	-0.92	-0.55	1.20	24.39	-8.41	-9.74	16.38
Dielectric constants, $\epsilon(\epsilon_0)$						$\beta(10^{-4}/\epsilon_0)$				Electromechanical coupling factors $k_{i\lambda}$					
$\epsilon_{11}^S$ <sup>a</sup>	$\epsilon_{33}^S$ <sup>a</sup>	$\epsilon_{11}^T$ <sup>a</sup>	$\epsilon_{33}^T$ <sup>a</sup>	$\beta_{11}^S$	$\beta_{33}^S$	$\beta_{11}^T$	$\beta_{33}^T$	$k_{15}^a$	$k_{31}^a$	$k_{33}^a$	$k_t^a$	$k_{15}^a$	$k_{31}^a$	$k_{33}^a$	$k_t^a$
870	601	6286	702	11.49	16.63	15.91	14.24	0.92	0.13	0.36	0.33	0.92	0.13	0.36	0.33

<sup>a</sup>Directly measured properties.

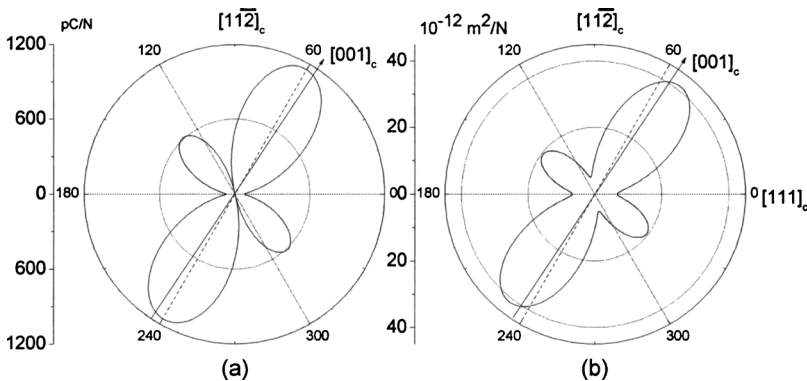


FIG. 2. Orientation dependence of piezoelectric constant  $d_{33}^*$  (a) and elastic compliance  $s_{33}^*$  (b), of single domain 0.26Pb(In<sub>1/2</sub>Nb<sub>1/2</sub>)O<sub>3</sub>–0.46Pb(Mg<sub>1/3</sub>Nb<sub>2/3</sub>)O<sub>3</sub>–0.28PbTiO<sub>3</sub> single crystal.

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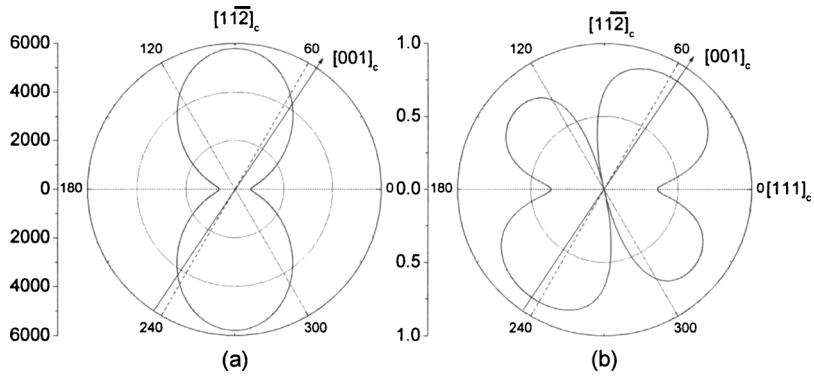


FIG. 3. Orientation dependence of dielectric constant  $\epsilon_{33}^*$  (a) and electromechanical coupling factor  $k_{33}^*$  (b) of single domain  $0.26\text{Pb}(\text{In}_{1/2}\text{Nb}_{1/2})\text{O}_3 - 0.46\text{Pb}(\text{Mg}_{1/3}\text{Nb}_{2/3})\text{O}_3 - 0.28\text{PbTiO}_3$  single crystal.

<sup>1</sup>X. Liu, S. Zhang, J. Luo, T. R. Shrout, and W. Cao, *Appl. Phys. Lett.* **96**, 012907 (2010).

<sup>2</sup>R. Zhang, B. Jiang, and W. Cao, *Appl. Phys. Lett.* **85**, 6380 (2004).