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Corrigendum

Corrigendum to "Crystal structures of α and β forms of poly(tetramethylene succinate)" (Polymer 2000:41;4719–27)☆

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The authors regret that errors were included in the β form molecular coordinates in Table 4, p. 4722. The correct table is shown below. These changes do not alter the results and discussion of this paper.

Table 4
The final fractional atomic coordinates of the α and β forms of PTMS

Atom	α form			β form		
	x	*y*	*z*	*x*	*y*	*z*
C_1	0.0839	0.0711	0.5389	0.0401	0.0372	0.5700
C_2	−0.0117	0.1387	0.6360	0.0049	−0.0916	0.6496
C_3	−0.0195	0.0832	0.8352	−0.0286	−0.0630	0.8347
C_4	0.0690	−0.0216	0.9564	0.0697	0.0284	0.9670
O_1	0.0696	0.0351	0.7495	0.0852	−0.0075	0.7753
O_2	−0.1536	0.1964	0.8164	−0.1948	−0.1766	0.7837
H_1a	0.0392	0.1534	0.4566	−0.1116	0.1372	0.5383
H_1b	0.3308	0.0504	0.6104	0.2728	0.0802	0.6432
H_2a	0.1104	0.2419	0.6821	0.1563	−0.1914	0.6803
H_2b	−0.2591	0.1590	0.5696	−0.2280	−0.1343	0.5754
H_4a	0.3190	−0.0237	1.0322	0.0077	0.1541	0.9370
H_4b	−0.0138	−0.1308	0.9106	0.3143	0.0186	1.0552

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