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Corrigendum

Corrigendum to "First principles study of the conformations of cinchonidine on a Pt(111) surface" [J. Catal. 239 (2006) 220–226]

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The authors regret that within Table 1 of the article cited above the entries in the last line were not correct and should be replaced with the values given in the table below.

Table 1
Adsorption energies $E_{Ads.}$ calculated with respect of the Open(3) conformer of CD (in kcal/mol)

	SO(3)	SO(4)	SC(1)	SC(2)	SQB(1)	SQB(2)	T(1)	SVB(1)
$E_{Ads.}$	41.5	39.3	43.2	42.7	43.0	36.6	22.1	50.4

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