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Corrigendum

Corrigendum to "A conformational study of N-acetyl glucosamine derivatives utilizing residual dipolar couplings" [J. Magn. Reson. 212 (2011) 174–185]

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In Fig. 4 the denomination of the regions for the substances 1, 2, 3 and 5 has to be changed to that shown in the corrected figure. Due to the symmetric aglycones of that substances the conformers represented by the regions "A" and "B" as well as "C" and "D" are equal.

Consequently, in Table 5 in the column "Region" for substance 5 "A/B" changes to "A".

Table 5
Dihedral angles of the conformers obtained from the MD simulations.

Substance	Region[a]	Φ (°)	Ψ (°)
1	A	−87.3	172.9
	B	63.8	172.1
2	A	−45.2	−164.3
	B	74.4	−34.4
3	A	−56.1	173.7
	B	43.8	−164.6
4	A	−51.1	−0.6
	B	−68.4	164.1
	C	66.1	−37.9
	D	58.0	171.0
5	A	−79.2	173.4
6	A	−62.1	81.4
	B	−65.5	−105.1

[a] The letters in this column correspond to the letters shown in the Φ/Ψ-plots.

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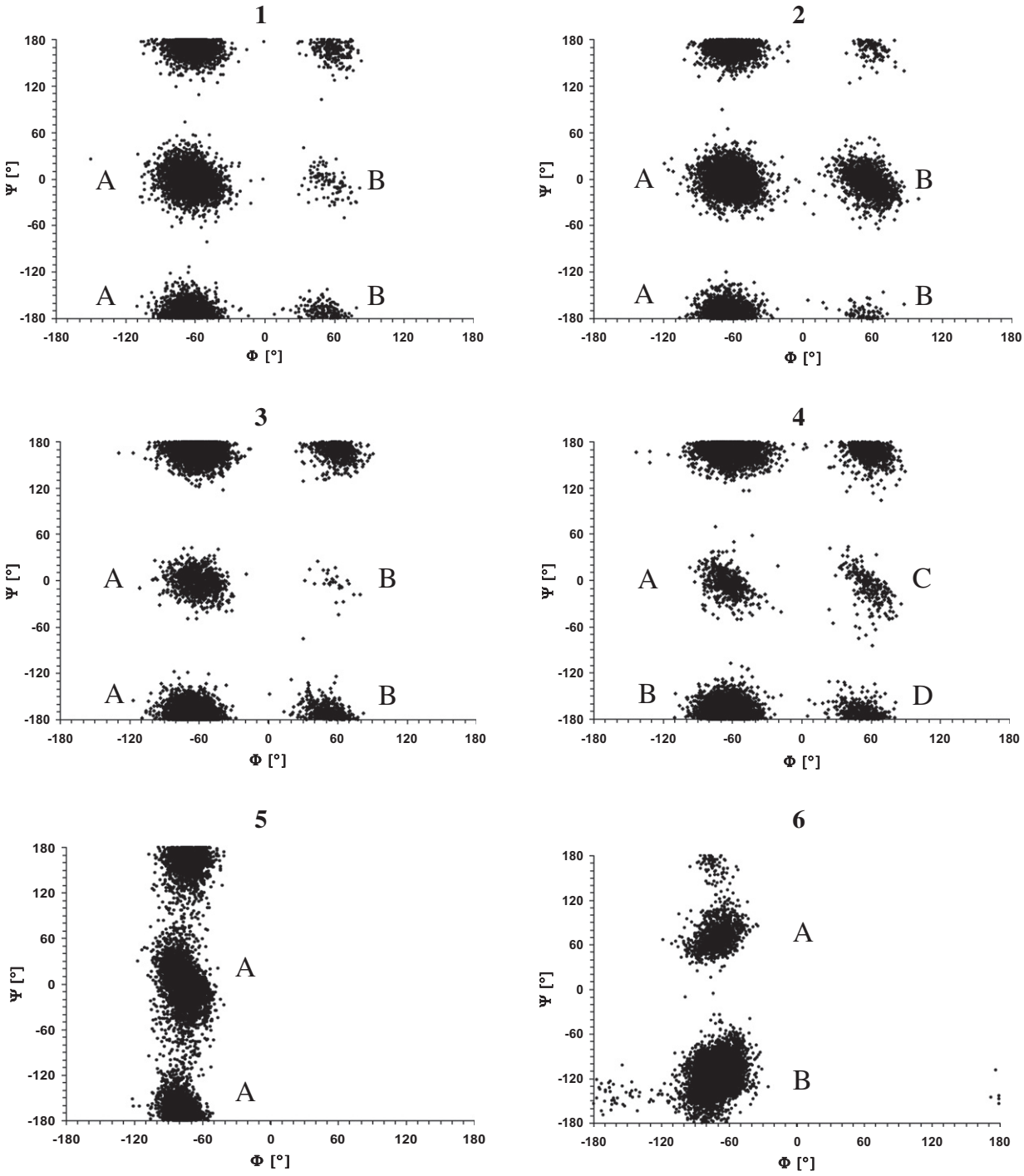

Fig. 4. Φ/Ψ-plots from the MD simulations.