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Erratum

Ab initio Study, Including Electron Correlation, of the Electronic Structures, the Dipole Moments, the Static Polarizabilities and of the Harmonic Fields of H_2CO, H_2CS and H_2SiO

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There is one error in Table 2. The values for the CH-bond length in H_2CO calculated with the CEPA-PNO method (basis *A*, basis *B*) have to be interchanged.

H_2CO: CEPA-PNO

	basis *A*	basis *B*
r_{CH}[Å]	1.116	1.104

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