Environmental Science Nano



CORRECTION

View Article Online



Cite this: Environ. Sci.: Nano, 2020, 7, 3627

Correction: Enhanced uptake of glyoxal at the acidic nanoparticle interface: implications for secondary organic aerosol formation

Qiuju Shi,^a Weina Zhang,^a Yuemeng Ji, [®]* Jiaxin Wang,^a Dandan Qin,^a Jiangyao Chen, ^{Dab} Yanpeng Gao, Guiying Li ^{Dab} and Taicheng An ^{Da}

DOI: 10.1039/d0en90056g

rsc li/es-nano

Correction for 'Enhanced uptake of glyoxal at the acidic nanoparticle interface: implications for secondary organic aerosol formation' by Qiuju Shi et al., Environ. Sci.: Nano, 2020, 7, 1126-1135, DOI: 10.1039/ D0EN00016G.

There is an error in Fig. 6 and therefore the graphical abstract, which incorporates Fig. 6, also includes this error. The error does not affect the results in this section. The corrected figures appear below.

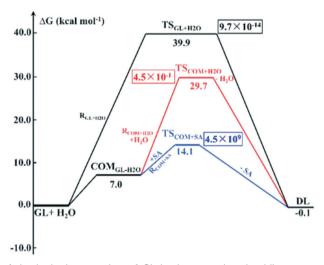
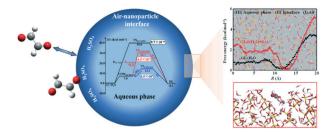


Fig. 6 The potential energy surface of the hydration reaction of GL in the neutral and acidic systems. The black line represents the direct $hydration\ reaction\ of\ GL\ (R_{GL+H_2O}).\ The\ red\ and\ blue\ lines\ represent\ the\ indirect\ hydration\ reaction\ of\ COM_{GL-H_2O}\ with\ H_2O\ and\ SA\ (R_{COM+H_2O}\ and\ sA\ (R_{COM+H_2O$ $R_{\text{COM+SA}}$), respectively. The unit of the rate constant is M^{-1} s⁻¹.

^a Guangdong Key Laboratory of Environmental Catalysis and Health Risk Control, Guangzhou Key Laboratory Environmental Catalysis and Pollution Control, School of Environmental Science and Engineering, Institute of Environmental Health and Pollution Control, Guangdong University of Technology, Guangzhou 510006, China. E-mail: jiym@gdut.edu.cn

^b Synergy Innovation Institute of GDUT, Shantou 515041, China



Graphical abstract

Correspondingly, in Section 3.3, on page 1131, the following sentence should be included in the third sentence of the third paragraph: "Herein, the discussed ΔG^{\ddagger} value of $R_{\text{COM+H,O}}/R_{\text{COM+SA}}$ is defined as the difference between the energies of TS_{COM+H₂O}/TS_{COM+SA} and COM_{GL-H₂O".}

Additionally, the second sentence in the fourth paragraph should be "Our calculated ΔG^{\dagger} value of $R_{\rm GL+H,O}$ agrees with the previous theoretical study (37.0 kcal mol⁻¹)⁶² which obtained the solvent energies based on the optimized gas-phase geometry". The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.