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Letters to the Editor

LETTER TO EDITOR

Discussion on the comments on DFT calculation, biological activity, anion sensing studies and crystal structure of (*E*)-4-chloro-2-((pyridin-2-ylimino)-methyl)phenol by Nuray Yıldırım, Neslihan Demir, Gökhan Alpaslan, Bahadir Boyacıoğlu, Mustafa Yıldız and Huseyin Ünver, published in the *Journal of the Serbian Chemical Society*, Volume 83, Issue 6, 2018, pp. 707–721

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We re-evaluated the X-ray data carefully. Since we worked on the Schiff bases of 2-aminopyridine and 3-aminopyridine, the possibility that the crystals have been confused cannot be ruled out.

Therefore, we agree with Marko V. Rodić on B structure and would like to thank for his careful consideration. The data and studies in the paper https://doi.org/10.2298/JSC180925082R are correct.

We really regret this undesirable confusion.

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