Corrigendum

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Crystal structure of 4-bromo-2-(1*H*-pyrazol-3-yl) phenol, $C_9H_7BrN_2O$

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In the paper by Jaćimović *et al.* [*Z. Kristallogr. NCS* **2017**; 232(3): 507–509] by an oversight, the crystal system of the reported structure was presented as triclinic instead of monoclinic.

The corrected Abstract of the paper *Z. Kristallogr. NCS* **2017**; 232(3): 507–509:

Abstract

$C_9H_7BrN_2O$, monoclinic, $C2/c$ (no. 15), $a = 16.255(3)$ Å, $b = 4.4119(9)$ Å, $c = 25.923(5)$ Å, $\beta = 107.99(3)°$, $V = 1768.2(7)$ Å^3, $Z = 8$, $R_{gt}(F) = 0.0450$, $wR_{ref}(F^2) = 0.0960$, $T = 150$ K.

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