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(2Z)-3-(4-Chloroanilino)-1-(5-hydroxy-3-methyl-1-phenyl-1*H*-pyrazol-4-yl)but-2-en-1-one A. M. Asiri, A. O. Al-Youbi, K. A. Alamry, H. M. Faidallah, S. W. Ng and E. R. T. Tiekink

Abstract: With the exception of the terminal benzene rings, the atoms in the title compound, $C_{20}H_{18}ClN_3O_2$, are approximately coplanar (r.m.s. deviation = 0.0495 Å). The benzene/chlorobenzene rings form dihedral angles of 3.02 (4) and 41.59 (5)°, respectively, with this plane. The hydroxy, amino and carbonyl groups all lie to the same side of the molecule, enabling the formation of intramolecular O-H---O and N-H----O hydrogen bonds that close S(6) rings. The configuration about the 2-butene bond is Z. Supramolecular chains mediated by C-H---Cl interactions and aligned along the c axis are found in the crystal packing. These assemble into layers that are connected by weak π -- π interactions between centrosymmetrically related chlorobenzene rings [3.8156 (9) Å].