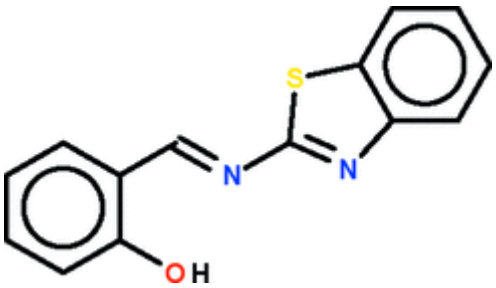

Acta Cryst. (2010). E**66**, o1826 [doi:10.1107/S1600536810023755]

A monoclinic modification of 2-[(1,3-benzothiazol-2-yl)iminomethyl]phenol

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Abstract: In the title Schiff base, $C_{14}H_{10}N_2OS$, the azomethine double bond is in an *E* configuration; the benzothiazolyl ring (r.m.s. deviation = 0.007 Å) is coplanar with the phenylene ring (r.m.s. deviation = 0.007 Å), the two rings being slightly bent at 2.6 (1)°. The hydroxy H atom forms an intramolecular hydrogen bond to the imino group. The bond dimensions of the monoclinic modification are similar to those of the orthorhombic modification [Liu *et al.* (2009). *Acta Cryst.* E**65**, o738].