



>>Entry Name: *N*-(4-Hydroxyphenyl)thiourea, 9CI

Synonym(s): 1-(*p*-Hydroxyphenyl)-2-thiourea, 8CI

CAS Registry Number: 1520-27-0

Molecular Formula: C₇H₈N₂OS

Molecular Weight: 168.219

Accurate Mass: 168.035733

Percentage Composition: C 49.98%; H 4.79%; N 16.65%; O 9.51%; S 19.06%

Physical Description: Plates (EtOH)

Melting Point: Mp 220-221 dec.°

Hazard & Toxicity: LD₅₀ (mus, ipr) 100 mg/kg

>>Derivative: Me ether

Synonym(s): *N*-(4-Methoxyphenyl)thiourea

CAS Registry Number: 2293-07-4

Molecular Formula: C₈H₁₀N₂OS

Molecular Weight: 182.246

Accurate Mass: 182.051383

Percentage Composition: C 52.72%; H 5.53%; N 15.37%; O 8.78%; S 17.59%

Melting Point: Mp 208-210°

Aldrich: 19936-2

Rare Chemicals Library: S40916-2

>>Derivative: Et ether

Synonym(s): *N*-(4-Ethoxyphenyl)thiourea

CAS Registry Number: 880-29-5

Molecular Formula: C₉H₁₂N₂OS

Molecular Weight: 196.273

Accurate Mass: 196.067033

Percentage Composition: C 55.08%; H 6.16%; N 14.27%; O 8.15%; S 16.34%

Physical Description: Cryst. (EtOH)

Melting Point: Mp 172°

References:

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Yokikawa, T. *et al.*, *Pharm. Bull.*, 1954, **1**, 54

Kopüchuk, I.I., *CA*, 1965, **63**, 2913

Hughes, J.L. *et al.*, *J. Med. Chem.*, 1975, **18**, 1077

Agarwal, S. *et al.*, *Indian J. Chem., Sect. B*, 1987, **26**, 642, (*synth*)