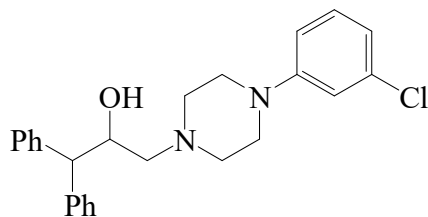




>>Entry Name: 4-(3-Chlorophenyl)- α -(diphenylmethyl)-1-piperazineethanol, 9Cl

Synonym(s): 3-[4-(3-Chlorophenyl)piperazin-1-yl]-1,1-diphenyl-2-propanol. BRL 15572



Related CAS Registry Numbers(s): 193611-72-2

Molecular Formula: C₂₅H₂₇ClN₂O

Molecular Weight: 406.954

Accurate Mass: 406.18119

Percentage Composition: C 73.79%; H 6.69%; Cl 8.71%; N 6.88%; O 3.93%

Biological Use/Importance: 5-HT_{1D} receptor antagonist

Physical Description: No phys. props. found

References:

Price, G.W. *et al.*, *Naunyn-Schmiedeberg's Arch. Pharmacol.*, 1997, **356**, 312-320, (*pharmacol*)

Schlicker, E. *et al.*, *Naunyn-Schmiedeberg's Arch. Pharmacol.*, 1997, **356**, 321-327, (*pharmacol*)