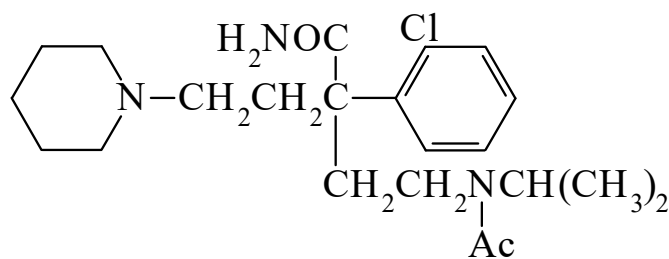




>>Entry Name: Bidisomide, INN, USAN

Synonym(s): α -[2-[Acetyl(1-methylethyl)amino]ethyl]- α -(2-chlorophenyl)-1-piperidinebutanamide, 9Cl. α -*o*-(Chlorophenyl)- α -[2-(*N*-isopropylacetamido)ethyl]-1-piperidinebutyramide. SC 40230



CAS Registry Number: 103810-45-3

Related CAS Registry Numbers(s): 121530-69-6 121530-70-9

Molecular Formula: C₂₂H₃₄ClN₃O₂

Molecular Weight: 407.982

Accurate Mass: 407.233954

Percentage Composition: C 64.77%; H 8.40%; Cl 8.69%; N 10.30%; O 7.84%

Use/Importance: Antiarrhythmic (Class Ia/Ib) agent. Cardiac depressant

pKa Value: pK_a 9.3

Experimental Log P Data: Log P 0.5 (exp)

Calculated Log P Data: Log P 1.92 (calc)

>>Variant: (±)-form

CAS Registry Number: 116078-65-0

Molecular Formula: C₂₂H₃₄ClN₃O₂

Molecular Weight: 407.982

Accurate Mass: 407.233954

Percentage Composition: C 64.77%; H 8.40%; Cl 8.69%; N 10.30%; O 7.84%

Melting Point: Mp 140-141°

References:

Desai, B.N. *et al.*, *J. Med. Chem.*, 1988, **31**, 2158, (*synth, pharmacol, pmr*)

Frederick, L.G. *et al.*, *Cardiovasc. Res.*, 1989, **23**, 897, (*pharmacol*)

Martin, C.L. *et al.*, *Drug Dev. Res.*, 1989, **17**, 51; 119, (*activity*)

Page, R.L. *et al.*, *Clin. Pharmacol. Ther. (St. Louis)*, 1992, **51**, 371, (*pharmacol, man*)

Cook, C.S. *et al.*, *Xenobiotica*, 1993, **23**, 1299, (*pharmacokinet*)

Komori, S., *Heart Vessels*, 1995, **10**, 7, (*pharmacol*)