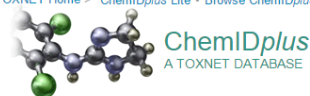


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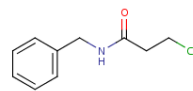
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Substance Name: Beclamide [INN:BAN:DCF]
RN: 501-68-8
InChIKey: JPYQFYIEOUVJDU-UHFFFAOYSA-N

Classification Code
 Drug / Therapeutic Agent

Molecular Formula
 C10-H12-Cl-N-O
Molecular Weight
 197.6638



All Links to Resources Names & Synonyms Registry Numbers Structure Descriptors Toxicity Physical Properties

Links to Resources**NLM Resources (File Locators)**

[DART](#) [RTECS](#) [DrugPortal](#) [PubMed Cancer](#)
[EINECS](#) [TOXLINE](#) [PubMed](#)
[MeSH](#) [PubChem](#) [PubMed Toxicology](#)

Other Resources (Internet Locators)

[NIST WebBook](#) [EPA ACToR](#)

Search for this InChIKey on the Web

Names and Synonyms**Name of Substance**

[Beclamide](#) [Beclamide \[INN:BAN:DCF\]](#)

Synonyms

[3-Chloro-N-\(phenylmethyl\)propanamide](#) [Benzchlorpropamide](#) [Chloroethylphenamide](#) [Neuracen](#)
[4-12-00-02234 \(Beilstein Handbook Reference\)](#) [Benzchlorpropamid](#) [Chloropropionamide](#) [Nidrane](#)
[Beclamid](#) [Benzchlorpropamide](#) [EINECS 207-927-1](#) [NSC 67062](#)
[Beclamida \[INN-Spanish\]](#) [Benzochlorpropamid](#) [Hibicon](#) [Nydran](#)
[Beclamide](#) [Benzylamide](#) [Klorakon](#) [Nydrane](#)
[Beclamidum](#) [BRN 2720702](#) [N-\(3-Chloropropionyl\)benzylamine](#) [Posedrin](#)
[Beclamidum \[INN-Latin\]](#) [Chlorakon](#) [N-Benzyl-3-chloropropionamide](#) [Posedrine](#)
[Beclamide](#) [Chlorakon \(VAN\)](#) [N-Benzyl-beta-chloropropionamide](#) [Sectar](#)
[Beclamidum](#) [Chlorakon \(VAN\)](#) [N-Benzyl-beta-chloropropionamide](#) [UNII-F5N0ALI65V](#)
[Beklamid](#) [N-Benzyl-beta-chloropropionamide](#)

Systematic Names

[Beclamide](#) [N-Benzyl-3-chloropropionamide](#) [Propanamide, 3-chloro-N-\(phenylmethyl\)- \(9Cl\)](#) [Propionamide, N-benzyl-3-chloro-](#)

Registry Numbers**CAS Registry Number**

[501-68-8](#)

System Generated Number

[0000501688](#)

Structure Descriptors**InChI**

1S/C10H12ClNO/c11-7-6-10(13)12-8-9-4-2-1-3-5-9/h1-5H,6-8H2,(H,12,13)

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InChIKey

JPYQFYIEOUVJDU-UHFFFAOYSA-N

[Search the web for this InChIKey](#)

Smiles

c1ccc(cc1)CNC(=O)CCCl

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Toxicity

Organism	Test Type	Route	Reported Dose (Normalized Dose)	Effect	Source
mammal (species unspecified)	LD50	unreported	1368mg/kg (1368mg/kg)		Pharmaceutical Chemistry Journal Vol. 28, Pg. 452, 1994.
mouse	LD50	intraperitoneal	650mg/kg (650mg/kg)		"Psychotropic Drugs and Related Compounds," 2nd ed., Usdin, E., and D.H. Efron, Washington, DC, 1972Vol. -, Pg. 384, 1972.

<http://chem.sis.nlm.nih.gov/chemidplus/rn/501-68-8>

Species (unspecified)	Route	Dose (mg/kg)	Notes	Source
mouse	LD50 intraperitoneal	650mg/kg (650mg/kg)		"Psychotropic Drugs and Related Compounds," 2nd ed., Usdin, E., and D.H. Efron, Washington, DC, 1972Vol. -, Pg. 384, 1972.
mouse	LD50 oral	1gm/kg (1000mg/kg)	BEHAVIORAL: ANTICONVULSANT	Pharmaceutical Chemistry Journal Vol. 14, Pg. 99, 1980.
rat	LD50 intraperitoneal	770mg/kg (770mg/kg)		Journal of Pharmacology and Experimental Therapeutics. Vol. 107, Pg. 403, 1953. Link to PubMed
rat	LD50 intravenous	770mg/kg (770mg/kg)		"Psychotropic Drugs and Related Compounds," 2nd ed., Usdin, E., and D.H. Efron, Washington, DC, 1972Vol. -, Pg. 384, 1972.
rat	LD50 oral	3200mg/kg (3200mg/kg)		Journal of Pharmacology and Experimental Therapeutics. Vol. 107, Pg. 403, 1953. Link to PubMed

Physical Properties

Physical Property	Value	Units	Temp (deg C)	Source
Melting Point	94	deg C		EXP
log P (octanol-water)	1.760	(none)		EST
Water Solubility	100	mg/L		EXP
Atmospheric OH Rate Constant	1.58E-11	cm ³ /molecule-sec	25	EST

Physical property data is provided to ChemIDplus by Syracuse Research Corporation.
See all available property data for this compound, including references.