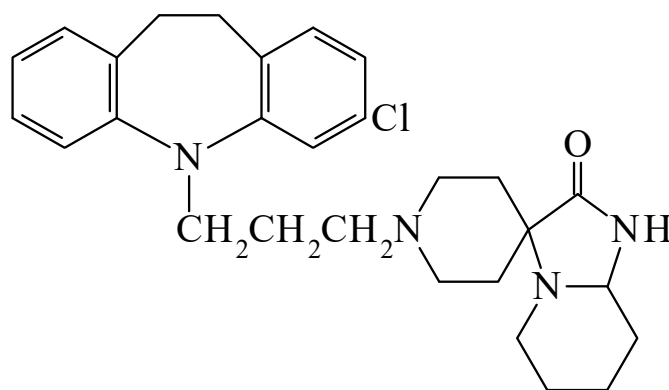




>>Entry Name: Mosapramine, INN

Synonym(s): 1'-[3-(3-Chloro-10,11-dihydro-5*H*-dibenz[*b,f*]azepin-5-yl)propyl]hexahydrospiro[imidazo[1,2-*a*]pyridine-3(2*H*),4'-piperidin]-2-one, 9Cl. 3-Chloro-5-[3-(2-oxo-1,2,3,5,6,7,8,8a-octahydroimidazo[1,2-*a*]pyridine-3-spiro-4'-piperidino)propyl]-10,11-dihydro-5*H*-dibenz[*b,f*]azepine. Clospipramine. Y 516



CAS Registry Number: 123435-19-8

Related CAS Registry Numbers(s): 120167-35-3 120181-35-3

Molecular Formula: C₂₈H₃₅ClN₄O

Molecular Weight: 479.063

Accurate Mass: 478.249938

Percentage Composition: C 70.20%; H 7.36%; Cl 7.40%; N 11.70%; O 3.34%

Use/Importance: Dopamine D₂-receptor antagonist. Neuroleptic

Development Status: Launched 1991 (Japan)

Calculated Log P Data: Log P 6.49 (uncertain value) (calc)

Other Data: Minor differences only in pharmacol. of enantiomers

>>Variant: (*R*)-form

Molecular Formula: C₂₈H₃₅ClN₄O

Molecular Weight: 479.063

Accurate Mass: 478.249938

Percentage Composition: C 70.20%; H 7.36%; Cl 7.40%; N 11.70%; O 3.34%

Physical Description: Cryst. (EtOAc)

Melting Point: Mp 145-147°

Optical Rotation: [α]_D −39.6 (c, 1 in CHCl₃) (98% ee)

>>Derivative: Hydrochloride (1:2)

Melting Point: Mp 270-272 dec.°

Optical Rotation: [α]_D +8.6 (c, 0.5 in H₂O)

>>Variant: (*S*)-form

Molecular Formula: C₂₈H₃₅ClN₄O

Molecular Weight: 479.063

Accurate Mass: 478.249938

Percentage Composition: C 70.20%; H 7.36%; Cl 7.40%; N 11.70%; O 3.34%

Physical Description: Cryst. (EtOAc)

Melting Point: Mp 142-143°

Optical Rotation: [α]_D +39.2 (c, 1 in CHCl₃) (97% ee)

>>Derivative: Hydrochloride (1:2)

Melting Point: Mp 269-270 dec.°

Optical Rotation: [α]_D −6 (c, 0.5 in H₂O)

>>Variant: (±)-form

CAS Registry Number: 89419-40-9

Molecular Formula: C₂₈H₃₅ClN₄O

Molecular Weight: 479.063

Accurate Mass: 478.249938

Percentage Composition: C 70.20%; H 7.36%; Cl 7.40%; N 11.70%; O 3.34%

>>Derivative: Hydrochloride (1:2)

Synonym(s): Mosapramine hydrochloride, JAN. Cremin