

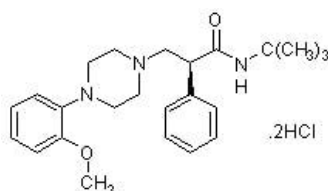
Product datasheet

(S)-WAY 100135, 5-HT_{1A} receptor antagonist ab146821

[2 References](#) [1 Image](#)

Overview

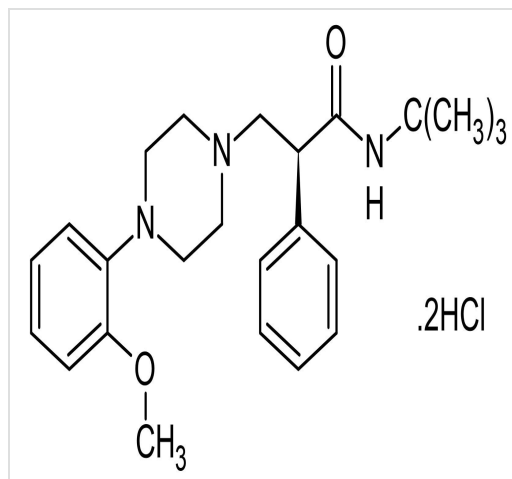
Product name	(S)-WAY 100135, 5-HT _{1A} receptor antagonist
Description	Potent, selective 5-HT _{1A} receptor antagonist
Purity	> 99%
CAS Number	149007-54-5
Chemical structure	



Properties

Chemical name	(S)-(-)- <i>N</i> - <i>tert</i> -Butyl-3-(4-(2-methoxyphenyl)piperazin-1-yl)-2-phenylpropanamide dihydrochloride
Molecular weight	468.47
Molecular formula	C ₂₄ H ₃₃ N ₃ O ₂ ·2HCl
PubChem identifier	14801907
Storage instructions	Store at room temperature.
Solubility overview	Soluble in water to 10 mM and in DMSO to 100 mM
Handling	Toxic, refer to SDS for further information Need more advice on solubility, usage and handling? Please visit our frequently asked questions (FAQ) page for more details.
SMILES	<chem>CC(C)(C)NC(=O)[C@H](CN1CCN(CC1)C2=CC=CC=C2OC)C3=CC=CC=C3.Cl.Cl</chem>
Source	Synthetic

Images



2D chemical structure image of ab146821, (S)-WAY 100135, 5-HT_{1A} receptor antagonist

Chemical Structure - (S)-WAY 100135, 5-HT_{1A} receptor antagonist (ab146821)

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