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## DATASHEET

### SB 224289 hydrochloride

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#### Product overview

<b>Name</b>	SB 224289 hydrochloride
<b>Cat No</b>	HB1667
<b>Biological action</b>	Antagonist
<b>Purity</b>	>95%
<b>Description</b>	Selective 5-HT <sub>1B</sub> receptor antagonist

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#### Biological Data

<b>Biological description</b>	Selective 5-HT <sub>1B</sub> receptor antagonist (pK <sub>i</sub> = 8.2). Displays > 60-fold selectivity over 5-HT <sub>1D</sub> , 5-HT <sub>1A</sub> , 5-HT <sub>1E</sub> , 5-HT <sub>1F</sub> , 5-HT <sub>2A</sub> and 5-HT <sub>2C</sub> receptors. Active <i>in vivo</i> .
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#### Solubility & Handling

<b>Storage instructions</b>	Room temperature
<b>Solubility overview</b>	Soluble in DMSO (10mM)
<b>Important</b>	This product is for RESEARCH USE ONLY and is not intended for therapeutic or diagnostic use. Not for human or veterinary use.

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#### Chemical Data

<b>Chemical name</b>	1'-Methyl-5-[[[2'-methyl-4'-(5-methyl-1,2,4-oxadiazol-3-yl)biphenyl-4-yl]carbonyl]-2,3,6,7-tetrahydrospiro[furo[2,3-f]indole-3,4'-piperidine hydrochloride
<b>Molecular Weight</b>	557.09
<b>Molecular Formula</b>	C <sub>32</sub> H <sub>32</sub> N <sub>4</sub> O <sub>3</sub> ·HCl
<b>CAS Number</b>	180084-26-8
<b>PubChem identifier</b>	11226716
<b>SMILES</b>	<chem>O=C(C5=CC=C(C6=CC=C(C7=NOC(C)=N7)C=C6)C=C5)N2C1=CC3=C(OCC43CCN(CC4)C)C=C1CC2.Cl</chem>
<b>InChiKey</b>	GKKGKBYMDILCOF-UHFFFAOYSA-N

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