



## Product Information

### CCR inhibitor ML604086, Purity $\geq 98\%$

**Cat. No.:** X23-10-ZQ689

**Size:** 5 mg; 10 mg; 25 mg; 50 mg; 100 mg

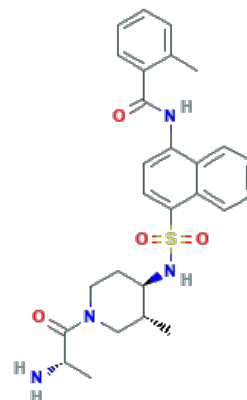
**MDL:** MFCD32201065

**CAS Number:** 850330-18-6

**Compound CID:** 57564362

**Synonym:** ML604086; 850330-18-6; Amino[sulfonyl]-1-naphthalenyl]-2-methylpiperidiny]; CCR inhibitor

**This product is for research use only and is not intended for diagnostic use.**



### Product Information

<b>Description</b>	ML604086 specifically binds to chemokine receptors. This binding blocks the receptors and prevents the corresponding chemokines from exerting their effects on immune cells, leading to reduced chemotaxis and cell activation.
<b>Molecular Weight</b>	508.63
<b>Molecular Formula</b>	C <sub>27</sub> H <sub>32</sub> N <sub>4</sub> O <sub>4</sub> S
<b>IUPAC Name</b>	<i>N</i> -[4-[[[(3 <i>R</i> ,4 <i>R</i> )-1-[(2 <i>S</i> )-2-Aminopropanoyl]-3-methylpiperidin-4-yl]sulfonyl]naphthalen-1-yl]-2-methylbenzamide
<b>InChI</b>	InChI=1S/C27H32N4O4S/c1-17-8-4-5-9-20(17)26(32)29-24-12-13-25(22-11-7-6-10-21(22)24)36(34,35)30-23-14-15-31(16-18(23)2)27(33)19(3)28/h4-13,18-19,23,30H,14-16,28H2,1-3H3,(H,29,32)/t18-,19+,23-/m1/s1
<b>InChI Key</b>	FABRBEGQMXBELT-SELNLUPBSA-N
<b>Canonical SMILES</b>	CC1CN(CCC1NS(=O)(=O)C2=CC=C(C3=CC=CC=C32)NC(=O)C4=CC=CC=C4C)C(=O)C(C)N
<b>Isomeric SMILES</b>	C[C@@H]1CN(CC[C@H]1NS(=O)(=O)C2=CC=C(C3=CC=CC=C32)NC(=O)C4=CC=CC=C4C)C(=O)[C@H](C)N
<b>Form</b>	Lyophilized powder
<b>Purity</b>	$\geq 98\%$
<b>Identity</b>	Confirmed by NMR/HPLC/MS.
<b>Stability</b>	The product is stable for three years when stored at the recommended temperature in lyophilized powder.
<b>Applications</b>	ML604086 can be used to explore its function as an inhibitor in biological pathways, providing insight into its therapeutic potential.



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**Storage**

Store at -20°C.

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