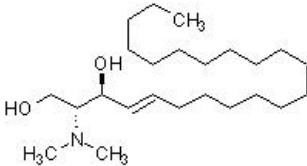


Product datasheet

N,N-Dimethyl-D-erythro-sphingosine (PKC inhibitor),
Sphingosine kinase and PKC inhibitor ab141237

Overview

Product name	N,N-Dimethyl-D-erythro-sphingosine (PKC inhibitor), Sphingosine kinase and PKC inhibitor
Description	Sphingosine kinase and PKC inhibitor
Biological description	Inhibitor of sphingosine kinase (IC ₅₀ =5 µM) and PKC (IC ₅₀ =12 µM). Blocks conversion of sphingosine to sphingosine-1-phosphate. Induces apoptosis and EGFR autophosphorylation. Produced <i>via</i> ceramide catabolism. Active <i>in vitro</i> .
Purity	> 98%
CAS Number	119567-63-4
Chemical structure	

Properties

Chemical name	(E,2S,3R)-2-(Dimethylamino)octadec-4-ene-1,3-diol
Molecular weight	327.55
Molecular formula	C ₂₀ H ₄₁ NO ₂
PubChem identifier	5282309
Storage instructions	Store at -20°C (desiccating conditions).
Solubility overview	Supplied in isopropanol (5 mg/ml)
Handling	Need more advice on solubility, usage and handling? Please visit our frequently asked questions (FAQ) page for more details.
SMILES	CCCCCCCCCCCC/C=C/[C@H]([C@H](CO)N(C)C)O
Source	Synthetic

Please note: All products are "FOR RESEARCH USE ONLY. NOT FOR USE IN DIAGNOSTIC PROCEDURES, NOT FOR USE IN HUMANS"

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- We provide support in Chinese, English, French, German, Japanese and Spanish
- Extensive multi-media technical resources to help you
- We investigate all quality concerns to ensure our products perform to the highest standards

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