

Bay 11-7085

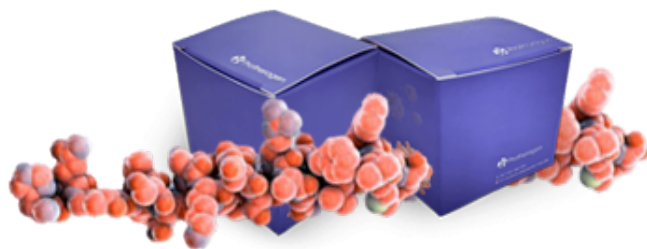
Cat. No.:	OB0225WX-2682
Appearance:	Solid
Purity:	≥99%
Identity:	Confirmed by NMR, and HPLC.
Size:	10 mg; 25 mg; 50 mg; 100 mg

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

Product Overview

Description	BAY 11-7085 is an irreversible antagonist of TNF-alpha-induced I-kappa-B-alpha p phosphorylation, characterized by a half-maximal inhibitory concentration (IC50) of 10 micromolar.
Synonym	bay 11-7085; 196309-76-9; (E)-3-((4-(tert-Butyl)phenyl)sulfonyl)acrylonitrile; (E)-3-(4-tert-butylphenyl)sulfonylacrylonitrile; bay11-7085; BAY 11 -7083; (E)-3-(4-tert-butylphenyl)sulfonylprop-2-enenitrile; MFCD01862602; BAY-117085; (E)-3-(4-tert-butylphenyl)sulfonyl-2-propenenitrile; BAY-11-708 5; (2E)-3-[[4-(1,1-dimethylethyl)phenyl]sulfonyl]-2-propenenitrile; 3-((4-(1,1-dimethylethyl)phenyl)sulfonyl)-2-propenenitrile; BAY 11-7085; (E)-3-[[4-(1,1-dimethylethyl)phenyl]sulfonyl]-2-propenenitrile; BAY 117083; BAY 11708 5; 2-Propenenitrile, 3-[[4-(1,1-dimethylethyl)phenyl]sulfonyl]-, (2E)-; (E)-3-(4-tert-butylbenzenesulfonyl)-acrylonitrile; (E)-3-(4-tert-butylphenyl)sulfonyl-2-propenenitrile; (E)-3-(4-tert-butylphenyl)sulfonyl-2-propenenitrile; (2E)-3-[(4-tert-butylphenyl)sulfonyl]prop-2-enenitrile; (2E)-3-(4-tert-butylbenzene-1-sulfonyl)prop-2-enenitrile
CAS No.	196309-76-9
Compound CID	5353432
Formula	C ₁₃ H ₁₅ NO ₂ S
Formula Weight	249.33

Specification



IUPAC Name	(E)-3-(4- <i>tert</i> -Butylphenyl)sulfonylprop-2-enenitrile
InChI	InChI=1S/C13H15NO2S/c1-13(2,3)11-5-7-12(8-6-11)17(15,16)10-4-9-14/h4-8,10H,1-3H3/b10-4+
InChI Key	VHKZGNPOHPFPER-ONNFQVAWSA-N
SMILES string	<chem>CC(C)(C)C1=CC=C(C=C1)S(=O)(=O)/C=C/C#N</chem>
Stability	3 years in powder form.
Storage	Storage at -20°C.
Applications	Bay 11-7085 is applied in molecular biology research to study the irreversible inhibition of IκBα phosphorylation and its subsequent impact on preventing NF-κB-mediated pro-inflammatory gene expression.

Library Information

Target	IκB/IKK
Receptor	IκBα
Pathways	NF-κB
Plate Number	L6700-08
Empty Location	a1-h1; a12-h12
Container	96-well plate
Formulation	10mM DMSO
DMSO Max Solubility	50 mg/mL; 200.54 mM
Ethanol Max Solubility	50 mg/mL
ALogP	2.635
HBA_Count	2
HBD_Count	0
Rotatable Bond	3