# **PRODUCT** INFORMATION



PF-04753299

Item No. 36502

CAS Registry No.: Formal Name:	1289620-49-0 N-hydroxy-αR-methyl-α-(methylsulfonyl)- [1,1'-biphenyl]-4-butanamide	
MF:	$C_{18}H_{21}NO_4S$	
FW:	347.4	CH OH
Purity:	≥98%	V V V N
Supplied as:	A solid	o=s=o H
Storage:	-20°C	
Stability:	≥4 years	
Information represents the product specifications. Batch specific analytical results are provided on each certificate of analysis.		

# Laboratory Procedures

PF-04753299 is supplied as a solid. A stock solution may be made by dissolving the PF-04753299 in the solvent of choice, which should be purged with an inert gas. PF-04753299 is soluble in the organic solvent DMSO at a concentration of approximately 34.7 mg/ml.

# Description

PF-04753299 is an inhibitor of LpxC (IC<sub>50</sub> = 1.37 nM), a bacterial deacetylase involved in the biosynthesis of LPS lipid A.<sup>1</sup> It is selective for LpxC over a panel of 16 human matrix metalloproteinases (MMPs; K<sub>s</sub> = >1  $\mu$ M for all). PF-04753299 is active against *P. aeruginosa*, K. pneumoniae, E. coli, E. aerogenes, C. freundii, and A. baumannii (MIC<sub>90</sub>s = 4, 16, 2, 4, 8, and 32 µg/ml, respectively). In vivo, PF-04753299 reduces splenic bacterial burden in a mouse model of systemic *P. aeruginosa* infection (ED<sub>50</sub> = 35 mg/kg).

# Reference

1. Brown, M.F., Reilly, U., Abramite, J.A., et al. Potent inhibitors of LpxC for the treatment of Gram-negative infections. J. Med. Chem. 55(2), 914-923 (2012).

WARNING THIS PRODUCT IS FOR RESEARCH ONLY - NOT FOR HUMAN OR VETERINARY DIAGNOSTIC OR THERAPEUTIC USE.

## SAFETY DATA

This material should be considered hazardous until further information becomes available. Do not ingest, inhale, get in eyes, on skin, or on clothing. Wash thoroughly after handling. Before use, the user must review the complete Safety Data Sheet, which has been sent via email to your institution.

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1180 EAST ELLSWORTH RD ANN ARBOR, MI 48108 · USA PHONE: [800] 364-9897 [734] 971-3335 FAX: [734] 971-3640 CUSTSERV@CAYMANCHEM.COM WWW.CAYMANCHEM.COM