# **PRODUCT** INFORMATION



**ML-161** 

Item No. 15179

CAS Registry No.:	: 423735-93-7	
Formal Name:	2-bromo-N-[3-[(1-oxobutyl)amino]	
Synonyms:	phenyl]-benzamide AG-670, CID-1048267	
MF:	$C_{17}H_{17}BrN_2O_2$	
FW:	361.2	
Purity:	≥95%	
UV/Vis.:	λ <sub>max</sub> : 240 nm	Br
Supplied as:	A crystalline solid	~
Storage:	-20°C	
Stability:	≥4 years	
Information represents the product specifications. Batch specific analytical results are provided on each certificate of analysis.		

# Laboratory Procedures

ML-161 is supplied as a crystalline solid. A stock solution may be made by dissolving the ML-161 in the solvent of choice. ML-161 is soluble in organic solvents such as ethanol and DMSO, which should be purged with an inert gas. The solubility of ML-161 in these solvents is approximately 25 mg/ml.

# Description

ML-161 is an allosteric, reversible inhibitor of proteinase-activated receptor 1 (PAR1) on platelets, preventing surface expression of P-selectin induced by the peptide SFLLRN with an IC<sub>50</sub> value of 0.26  $\mu$ M.<sup>1</sup> It blocks platelet activation induced by thrombin as well as by SFLLRN but not by PMA (Item No. 10008014), U-46619 (Item No. 16450), or collagen.<sup>1</sup>

# Reference

1. Dockendorff, C., Aisiku, O., VerPlank, L., et al. Discovery of 1,3-diaminobenzenes as selective inhibitors of platelet activation at the PAR1 receptor. ACS Med. Chem. Lett. 3(3), 232-237 (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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