# **PRODUCT** INFORMATION



## 7-deaza-8-bromo NAD<sup>+</sup> (sodium salt)

Item No. 38479

Formal Name:	3-(aminocarbonyl)-1-[5-O- [hydroxy(phosphonooxy)phosphinyl]- $\beta$ -D-ribofuranosyl]-pyridinium, P' $\rightarrow$ 5'-	2
	ester with 6-bromo-7-β-D-ribofuranosyl-	С
	7H-pyrrolo[2,3-d]pyrimidin-4-amine,	
Synonym:	7-deaza-8-bromo Nicotinamide adenine	
	dinucleotide N N O	
MF:	$C_{22}H_{26}BrN_6O_{14}P_2 \bullet Na$	
FW:	763.3 H <sub>N</sub> N <sup></sup> OH	
Purity:	≥95% <sup>2</sup> <u> </u>	
Supplied as:	A solid	
Storage:	-80°C	
Stability:	≥2 years	

Information represents the product specifications. Batch specific analytical results are provided on each certificate of analysis.

## Description

7-deaza-8-bromo NAD<sup>+</sup> is a brominated and deazaadenine derivative of the signaling molecule and cofactor NAD<sup>+</sup> (Item No. 16077).<sup>1</sup> It has been used as a synthetic intermediate in the synthesis of cyclic and membrane-permeable diphosphate modulators of transient receptor potential melastatin 2 (TRPM2).

## Reference

1. Moreau, C., Kirchberger, T., Swarbrick, J.M., et al. Structure-activity relationship of adenosine 5'-diphosphoribose at the transient receptor potential melastatin 2 (TRPM2) channel: Rational design of antagonists. J. Med. Chem. 56(24), 10079-100102 (2013).

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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