PRODUCT INFORMATION



SR 16832

Item No. 27632

CAS Registry No.: 2088135-12-8

Formal Name: 2-chloro-N-(6-methoxy-4-quinolinyl)-5-

nitro-benzamide

MF: C₁₇H₁₂CIN₃O₄

357.7 FW: **Purity:** ≥98% λ_{max} : 233 nm A crystalline solid UV/Vis.: Supplied as:

-20°C Storage: Stability: ≥4 years

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

Laboratory Procedures

SR 16832 is supplied as a crystalline solid. A stock solution may be made by dissolving the SR 16832 in the solvent of choice, which should be purged with an inert gas. SR 16832 is soluble in the organic solvent DMSO.

Description

SR 16832 is a dual-site covalent antagonist of peroxisome proliferator-activated receptor γ (PPARγ).¹ It inhibits MRL20-induced allosteric activation of PPARy in a reporter assay using HEK293T cells when used at a concentration of 5 μM. SR 16832 also reduces basal activity of PPARγ and inhibits binding of docosahexaenoic acid (DHA; Item No. 90310) to PPARy in a time-resolved FRET (TR-FRET) assay.

Reference

1. Brust, R., Lin, H., Fuhrmann, J., et al. Modification of the orthosteric PPARy covalent antagonist scaffold yields an improved dual-site allosteric inhibitor. ACS Chem. Biol. 12(4), 969-978 (2017).

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

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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