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## PKC€ Translocation Inhibitor II, PKCe141

**ALTERNATE NAME:** N-(3-Acetylphenyl)-9-amino-2,3-dihydro-[1,4]dioxino[2,3-g]thieno[2,3-b]quinoline-8-carboxamide; PKCe141

CATALOG #: AUR10281

## STRUCTURE:

MOLECULAR FORMULA: C22H17N3O4S

**MOLECULAR WEIGHT: 419.45** 

CAS#: N/A

APPEARANCE: Off-white solid

SOLUBILITY: DMSO (10 mg/ml)

**PURITY:** >95% by HPLC

STORAGE: Store at -2-8°C. Protect from air and light

**DESCRIPTION**: PKCe141 is a thienoquinoline derivative that dose-dependently inhibits PKCε-RACK2 interaction (IC<sub>50</sub> = 5.9 μM). Shown to occupy PKCε to RACK2 binding site in a reversible manner and block cellular phosphorylation of MARCKS and Elk-1 (IC<sub>50</sub> = 11.2 μM in HeLa cells). Prevents TPA-induced PKCε translocation to the plasma membrane, and reduces PKCε-induced cell migration, invasion and suppresses angiogenesis.

**HANDLING:** Do not take internally. Wear gloves and mask when handling the product! Avoid contact by all modes of exposure