

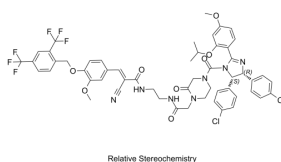
(rel)-PROTAC ERR α Degradar-1

Cat. No.: HY-128838A
Batch No.: 57881
Chemical Name: rel-(E)-N-(2-(2-(4-((4R,5S)-4,5-bis(4-chlorophenyl)-2-(2-isopropoxy-4-methoxyphenyl)-4,5-dihydro-1H

PHYSICAL AND CHEMICAL PROPERTIES

Molecular Formula: C₅₄H₄₉Cl₂F₆N₇O₈
Molecular Weight: 1108.91
Storage: -20°C, stored under nitrogen, away from moisture
* In solvent : -80°C, 6 months; -20°C, 1 month (stored under nitrogen, away from moisture)

Chemical Structure:



ANALYTICAL DATA

Appearance: Off-white to light yellow (Solid)
¹H NMR Spectrum: Consistent with structure
LCMS: Consistent with structure
Purity (LCMS): 95.89%
Conclusion: The product has been tested and complies with the given specifications.

Caution: Product has not been fully validated for medical applications. For research use only.

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