

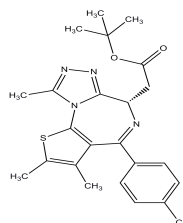
Bromodomain Inhibitor, (+)-JQ1

ALTERNATE NAME: *t*-Butyl-2-((6*S*)-4-(4-chlorophenyl)-2,3,9-trimethyl-6*H*-thieno[3,2-*f*][1,2,4]triazolo[4,3-*a*][1,4]diazepin-6-yl)acetate

CATALOG #: 2070-1, 5

AMOUNT: 1 mg, 5 mg

STRUCTURE:



MOLECULAR FORMULA: C₂₃H₂₅ClN₄O₂S

MOLECULAR WEIGHT: 456.99

CAS NUMBER: 1268524-70-4

APPEARANCE: Off-white solid

SOLUBILITY: DMSO (10 mM)

PURITY: ≥98% by HPLC

STORAGE: Store at -20 °C. Protect from air and moisture.

DESCRIPTION: Cell-permeable. A potent and highly specific inhibitor for the BET (bromodomain and extra-terminal) family of bromodomains. (+)-JQ1 binds to BRD4 bromodomains 1 and 2 with K_d values of ~ 50 and 90 nM, respectively. The binding is competitive with acetyl lysine. (+)-JQ1 can be a useful chemical probe to investigate the role of BET bromodomains in the transcriptional regulation of oncogenesis.

REFERENCE: Flippakopoulos, P., *et al.* (2010). *Nature* **468**, 1067-1073.

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

RELATED PRODUCTS:

- AMI-1 (Cat. No. 1943-5, 25)
- 5-Aza-2'-deoxycytidine (Cat. No. 1754-10, 50)
- 3-Deazaneplanocin (Cat. No. 2060-250, 1000)
- BIX01294 (Cat. No. 1678-5, 25)
- RG108 (Cat. No. 1679-10, 30)
- UNC0638 (cat. No. 1933-1, 5)