



Certificate of Analysis

www.tocris.com

Product Name: BSJ-03-204 Catalog No.: 6938 Batch No.: 1

CAS Number: 2349356-09-6

 $IUPAC \ Name: \ N-[4-[4-[6-[(6-Acetyl-8-cyclopentyl-7,8-dihydro-5-methyl-7-oxopyrido[2,3-\emph{a}]pyrimidin-2-yl)amino]-3-pyridinyl]-1-py$

piperazinyl]butyl]-2-[[2-(2,6-dioxo-3-piperidinyl)-2,3-dihydro-1,3-dioxo-1*H*-isoindol-4-yl]oxy]acetamide

1. PHYSICAL AND CHEMICAL PROPERTIES

Batch Molecular Formula: $C_{43}H_{48}N_{10}O_{8}.1\frac{1}{4}H_{2}O$

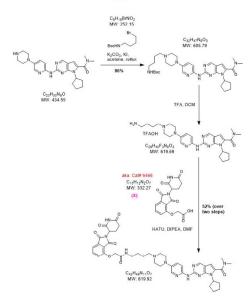
Batch Molecular Weight: 855.44

Physical Appearance: Yellow solid

Storage: DMSO to 20 mM Storage: Store at -20°C

Batch Molecular Structure:

5937



2. ANALYTICAL DATA

TLC: $R_f = 0.35$ (Dichloromethane:Methanol [9:1])

HPLC: Shows 97.7% purity

¹H NMR: Consistent with structure

Mass Spectrum: Consistent with structure

Microanalysis: Carbon Hydrogen Nitrogen

Theoretical 60.38 5.95 16.37 Found 59.98 5.85 16.04

Caution - Not Fully Tested • Research Use Only • Not For Human or Veterinary Use

www.tocris.com/distributors Tel:+1 612 379 2956



Product Information

Print Date: Mar 14th 2024

www.tocris.com

BSJ-03-204 1 **Product Name:** Catalog No.: 6938

CAS Number: 2349356-09-6

IUPAC Name: N-[4-[4-[6-[(6-Acetyl-8-cyclopentyl-7,8-dihydro-5-methyl-7-oxopyrido[2,3-d]pyrimidin-2-yl)amino]-3-pyridinyl]-1-

piperazinyl]butyl]-2-[[2-(2,6-dioxo-3-piperidinyl)-2,3-dihydro-1,3-dioxo-1H-isoindol-4-yl]oxy]acetamide

Description:

BSJ-03-204 is a selective Cdk4/6 Degrader (PROTAC®). Displays cereblon-dependent degradation with no effect on IKZF1/3 in Molt4 cells. Induces degradation of Cdk4/6 and G₁ cell cycle arrest, and inhibits cell proliferation in a mantle cell lymphoma (MCL) cell line. CDK4 antibodies validated for Simple Western™ (automated Western) instruments and Western Blot also available: Catalog # AF5254 and NBP1-31308. PROTAC® is a registered trademark of Arvinas Operations, Inc., and is used under license. Please see product specific page on www.tocris.com for full description.

Physical and Chemical Properties:

Batch Molecular Formula: C₄₃H₄₈N₁₀O₈.1½H₂O

Batch Molecular Weight: 855.44 Physical Appearance: Yellow solid

Minimum Purity: ≥98%

Batch Molecular Structure:

6937 C.H. BrNO MW: 252.15 C32H47N0O K₂CO₃, KI, TFA. DCM TFAOH C29H40F3N9O3 aka: Cat# 6466 C₁₅H₁₂N₂O₇ MW: 332.27 52% (over two steps HATU, DIPEA, DMF

Storage: Store at -20°C

Solubility & Usage Info:

DMSO to 20 mM

Stability and Solubility Advice:

Some solutions can be difficult to obtain and can be encouraged by rapid stirring, sonication or gentle warming (in a 45-60°C water bath).

Information concerning product stability, particularly in solution, has rarely been reported and in most cases we can only offer a general guide. *Unless contradicted by product-specific protocols or instructions, our standard recommendations apply:

SOLIDS: Provided storage is as stated on the product label and the vial is kept tightly sealed, the product can be stored for up to 6 months from date of receipt.

SOLUTIONS: We recommend that stock solutions, once prepared, are stored aliquoted in tightly sealed vials at -20°C or below and used within 1 month. Wherever possible solutions should be made up and used on the same day.

Licensing Information:

Sold under license from Dana-Farber Cancer Institute

Caution 19 Not Fully Tested Research Use Only • Not For Human or Veterinary Use

info.cn@bio-techne.com

bio-techne.com in far techne.com **North America** Tel: (800) 343 7475

Europe Middle East Africa

Rest of World

Tel: +44 (0)1235 529449 www.tocris.com/distributors techsupport@bio-techne.com Tel:+86 (21) 52380373

Jiang et al (2019) Development of dual and selective degraders of cyclin-dependent kinases 4 and 6. Angew.Chem.Int.Ed.Engl. 58

6321. PMID: 30802347.