



Certificate of Analysis

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Product Name: CST 905 Catalog No.: 7745 Batch No.: 1

 $IUPAC\ Name: (2S,4R)-1-((S)-2-(2-(4-(4-(3-(3-((N-Ethyl-N-methylsulfamoyl)amino)-2,6-difluorobenzoyl)-1H-pyrrolo[2,3-b]pyridin-5-(2S,4R)-1-((S)-2-(2-(4-(4-(3-(3-((N-Ethyl-N-methylsulfamoyl)amino)-2,6-difluorobenzoyl)-1H-pyrrolo[2,3-b]pyridin-5-(2S,4R)-1-((S)-2-(2-(4-(4-(3-(3-((N-Ethyl-N-methylsulfamoyl)amino)-2,6-difluorobenzoyl)-1H-pyrrolo[2,3-b]pyridin-5-(2S,4R)-1-((S)-2-(2-(4-(4-(3-(3-((N-Ethyl-N-methylsulfamoyl)amino)-2,6-difluorobenzoyl)-1H-pyrrolo[2,3-b]pyridin-5-(2S,4R)-1-((S)-2-(2-(4-(4-(3-(3-((N-Ethyl-N-methylsulfamoyl)amino)-2,6-difluorobenzoyl)-1H-pyrrolo[2,3-b]pyridin-5-(2S,4R)-1-(2S,4R$

yl)phenyl)piperazin-1-yl)acetamido)-3,3-dimethylbutanoyl)-4-hydroxy-N-(4-(4-methylthiazol-5-yl)benzyl)pyrrolidine-2-

carboxamide

1. PHYSICAL AND CHEMICAL PROPERTIES

Batch Molecular Formula: $C_{51}H_{58}F_2N_{10}O_7S_2$

Batch Molecular Weight: 1025.2

Physical Appearance: Yellow solid

Solubility: DMSO to 100 mM Storage: Store at -20°C

Batch Molecular Structure:

2. ANALYTICAL DATA

HPLC: Shows 98.0% purity

¹H NMR: Consistent with structure

Mass Spectrum: Consistent with structure

Microanalysis: Carbon Hydrogen Nitrogen

Theoretical 59.75 5.7 13.66 Found 58.77 5.86 13.19

9449 www.tocris.com/distributors Tel:+1 612 379 2956



Product Information

Print Date: Mar 14th 2024

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yl)phenyl)piperazin-1-yl)acetamido)-3,3-dimethylbutanoyl)-4-hydroxy-N-(4-(4-methylthiazol-5-yl)benzyl)pyrrolidine-2-

carboxamide

Description:

CST 905 is a potent and selective BRAF V600E Degrader (PROTAC $^{\otimes}$) (DC $_{50}$ = 18 nM). CST 905 comprises a ligand for the von Hippel Lindau (VHL) E3 ligase joined by a linker to a PLX-derived ligand. Unlike other BRAF V600E Degraders, CST 905 does not activate ERK. B-Raf antibody validated for Simple Western TM (automated Western) instruments and Western Blot also available: Catalog # AF3424. PROTAC $^{\otimes}$ 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: C₅₁H₅₈F₂N₁₀O₇S₂

Batch Molecular Weight: 1025.2 Physical Appearance: Yellow solid

Minimum Purity: ≥98%

Batch Molecular Structure:

Storage: Store at -20°C

CAUTION - This product is light sensitive and we recommend that the solid material and any solutions obtained are protected from exposure to light.

Solubility & Usage Info:

DMSO to 100 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.

References:

Miller et al (2022) Encoding BRAF inhibitor functions in protein degraders. RSC Med.Chem. 13 731. PMID: 35814929.