Certificate of Analysis

Print Date: Feb 15th 2023

Product Name: JHU 37160

CAS Number: 2369979-68-8

IUPAC Name: 8-Chloro-11-(4-ethylpiperazin-1-yl)-4-fluoro-5H-dibenzo[b,e][1,4]diazepine

1. PHYSICAL AND CHEMICAL PROPERTIES

Batch Molecular Formula: **Batch Molecular Weight:**

Physical Appearance:

Solubility:

Storage:

Batch Molecular Structure:

C₁₉H₂₀CIFN₄.¹/₄H₂O 363.35 Yellow solid DMSO to 50 mM Store at RT

N. CI

2. ANALYTICAL DATA

HPLC: ¹H NMR: Mass Spectrum: Microanalysis:

Shows 99.8% purity Consistent with structure Consistent with structure

	Carbon Hydrogen Nitrogen				
Theoretical	62.81	5.69	15.42		
Found	62.75	5.62	15.71		

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

bio-techne.com	North America	China	Europe Middle East Africa	Rest of World
info@bio-techne.com techsupport@bio-techne.com	Tel: (800) 343 7475	info.cn@bio-techne.com Tel: +86 (21) 52380373	Tel: +44 (0)1235 529449	www.tocris.com/distributors Tel:+1 612 379 2956



www.tocris.com

Catalog No.: 7198 Batch No.: 1

TOCRIS a biotechne brand

Print Date: Feb 15th 2023

www.tocris.com

Product Name: JHU 37160

CAS Number: 2369979-68-8

IUPAC Name:

8-Chloro-11-(4-ethylpiperazin-1-yl)-4-fluoro-5H-dibenzo[b,e][1,4]diazepine

Description:

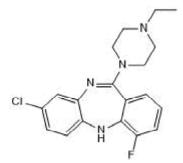
JHU 37160 is a high affinity and highly potent activator of hM_3D_q and hM_4D_i DREADDs (K_i values are 1.9 nM and 3.6 nM for hM_3D_q and hM_4D_i in vitro, respectively; EC₅₀ values are 0.2 nM and 18.5 nM for hM_4D_i and hM_3D_q in vitro, respectively). Displays approximately 25-fold higher affinity for hM_4D_i compared to DREADD agonist 21 (Cat. No. 5548). Selectively displaces [³H] clozapine from DREADDs in vivo, but not from other clozapine binding sites. Inhibits locomotor activity in mice expressing hM_3D_q and hM_4D_i in D_1 -expressing neurons, and increases hM_3D_q -stimulated locomotion in rats expressing hM_3D_q in THexpressing neurons. Brain penetrant in m... Please see product specific page on www.tocris.com for full description.

Physical and Chemical Properties:

Batch Molecular Formula: C₁₉H₂₀CIFN₄.¹/₄H₂O Batch Molecular Weight: 363.35 Physical Appearance: Yellow solid

Minimum Purity: ≥98%

Batch Molecular Structure:



References:

Bonaventura *et al* (2019) High-potency ligands for DREADD imaging and activation in rodents and monkeys. Nat.Commun. **10** 4627. PMID: 31604917.

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

bio-techne.com	North America	China	Europe Middle East Africa	Rest of World
info@bio-techne.com techsupport@bio-techne.com	Tel: (800) 343 7475	info.cn@bio-techne.com Tel: +86 (21) 52380373	Tel: +44 (0)1235 529449	www.tocris.com/distributors Tel:+1 612 379 2956

Catalog No.: 7198

1

Storage: Store at RT

Solubility & Usage Info: DMSO to 50 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 $45-60^{\circ}C$ water bath).

Information concerning product stability, particularly in solution, has rarely been reported and in most cases we can only offer a general guide. Our standard recommendations are:

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.