PRODUCT INFORMATION



Arry-380 analog

Item No. 24186

CAS Registry No.: 937265-83-3

Formal Name: 6-[5-[[[2-(methylsulfonyl)ethyl]amino]

> methyl]-2-furanyl]-N-[3-methyl-4-([1,2,4] triazolo[1,5-a]pyridin-7-yloxy)phenyl]-4-

quinazolinamine

MF: $C_{29}H_{27}N_7O_4S$

FW: 569.6 **Purity:**

UV/Vis.: λ_{max} : 334, 361 nm Supplied as: A crystalline solid

-20°C Storage: Stability: ≥4 years

Information represents the product specifications. Batch specific analytical results are provided on each certificate of analysis.

Laboratory Procedures

Arry-380 analog is supplied as a crystalline solid. A stock solution may be made by dissolving the arry-380 analog in the solvent of choice. Arry-380 analog is soluble in organic solvents such as DMSO and dimethyl formamide, which should be purged with an inert gas. The solubility of arry-380 analog in these solvents is approximately 25 mg/ml.

Arry-380 analog is sparingly soluble in aqueous buffers. For maximum solubility in aqueous buffers, arry-380 analog should first be dissolved in DMSO and then diluted with the aqueous buffer of choice. Arry-380 analog has a solubility of approximately 0.1 mg/ml in a 1:10 solution of DMSO:PBS (pH 7.2) using this method. We do not recommend storing the aqueous solution for more than one day.

Description

Arry-380 analog was designed as an inhibitor of EGFR (ErbB1).¹

Reference

1. Lyssikatos, J.P., Hicks, J.M., Marmsater, F.P., et al. N4-phenyl-quinazoline-4-amine derivatives and related compounds as ErbB type I receptor tyrosine kinase inhibitors for the treatment of hyperproliferative diseases. Array Biopharma. US20170252317A1 (2017).

WARNING
THIS PRODUCT IS FOR RESEARCH ONLY - NOT FOR HUMAN OR VETERINARY DIAGNOSTIC OR THERAPEUTIC USE.

This material should be considered hazardous until further information becomes available. Do not ingest, inhale, get in eyes, on skin, or on clothing. Wash thoroughly after handling. Before use, the user must review the complete Safety Data Sheet, which has been sent via email to your institution.

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