**BioVision** 05/14

PRODUCT: RG7388

**ALTERNATE NAME:** 4-[[(2R,3S,4R,5S)-3-(3-chloro-2-fluorophenyl)-4-(4-chloro-2-

fluorophenyl)-4-cyano-5-(2,2-dimethylpropyl)pyrrolidine-2-

carbonyl]amino]-3-methoxybenzoic acid; RG-7388

**CATALOG #:** 2577-1, 5

AMOUNT: 1 mg, 5 mg

STRUCTURE:

MOLECULAR FORMULA: C<sub>31</sub>H<sub>29</sub>Cl<sub>2</sub>F<sub>2</sub>N<sub>3</sub>O<sub>4</sub>

MOLECULAR WEIGHT: 616.50

**CAS No.** 1229705-06-9

**APPEARANCE:** White to off-white solid

SOLUBILITY: DMSO

PURITY: ≥99% by HPLC

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

**DESCRIPTION:** RG7388 is highly a potent and selective inhibitor of p53-

MDM2 interaction. RG7388 blocks p53–MDM2 binding and effectively activates the p53 pathway, leading to cell cycle arrest and/or apoptosis in cell lines expressing wild-type p53 and tumor growth inhibition or regression of osteosarcoma

xenografts in nude mice.

**REFERENCES:** Ding, Q., et al. (2013). J. Med. Chem. **56**, 5979-5983.

HANDLING:

Do not take internally. Wear gloves and mask when handling

the product! Avoid contact by all modes of exposure.

**RELATED PRODUCTS:** 

JNJ26854165 (Cat. No. 2479-5, 25) NSC-66811 (Cat. No. 2009-5) Nutlin-3 (Cat. No. 1842-1, 5) Pifithrin- $\alpha$  (Cat. No. 1554-10)

Pifithrin-α, cyclic, HBr (Cat. No. 2270-5)

USAGE: FOR RESEARCH USE ONLY! Not to be used in humans.