

# UbiQ

targeting the ubiquitin system

## K63 diUb-PA (human sequence, synthetic)

UbiQ code : UbiQ-114

Batch # : B01082015-001

Amount : 25 ug, lyophilized powder

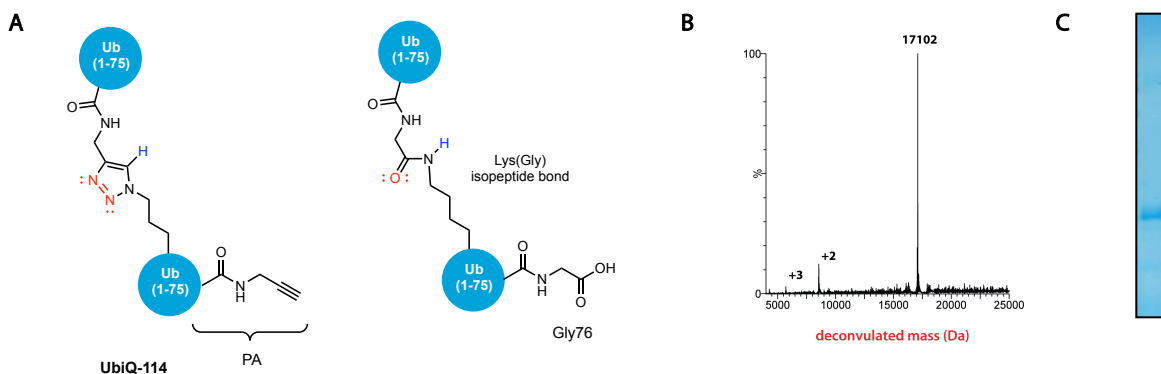
Purity :  $\geq 95\%$

Mol. Weight : 17.1 kDa

Storage : upon arrival, powder at  $-20^{\circ}\text{C}$ ; solution at  $-80^{\circ}\text{C}$ . Please avoid multiple freeze/thaw cycles.

## Productsheet

**Background.** K63 diUb-PA (UbiQ-114, Figure 1A) is an activity-based probe for deubiquitinating enzymes (DUBs). It is based on K63-linked that is modified on the C-terminus with a propargylamide (PA) electrophile. The isopeptide bond between the two Ub proteins is replaced by a DUB resistant triazole mimic. UbiQ-114 can be used for activity profiling experiments and determining DUB inhibitor specificity. The PA group has two unique capabilities: first, it forms a covalent linkage with (the active site Cys residue of) a DUB that can be cleaved by acid treatment (5% aq. TFA), allowing for proteomic analyses; secondly, it targets all three major DUB families: UCH, USP and OUT. Although Ub-PA based probes mainly target DUBs, the active-site Cys residue of certain HECT E3 ligases (HUWE1 and NEDD4) has been found to react with Ub-PA; as such Ub-PA based probes can also be used and evaluated for the study of HECT E3 ligases. For more details, please see reference 5 (Nair et al.).



**Figure 1.** A: structure UbiQ-114 and native isopeptide diUb. B: MS analysis UbiQ-114. C: SDS-PAGE analysis. 12% Bolt Bis-Tris Plus gel (Life technologies) and MES running buffer. CBB staining.

### important: sample preparation

- dissolve the powder in as little DMSO as possible (e.g., 20 - 40 mg/mL)
- add this DMSO stock slowly to milliQ (please note the order of addition)
- buffer as desired.

**Literature.** (1) Flierman et al. *Cell Chem Biol* **2016**, *23*, 472. (2) Ekkebus et al. *J Am Chem Soc* **2013**, *135*, 2867. (3) Sommer et al. *Bioorg Med Chem* **2013**, *21*, 2511. (4) de Jong et al. *ChemBioChem* **2012**, *13*, 2251. (5) Nair et al. *ACS Chem Biol* **2021**, *16*, 1615.