

Product number: K8-5046
Product name: Seta-650-Azide

General Data

- Molecular Mass:** 969.10
- Solubility:** water, alcohol, DMF, DMSO
- Insoluble:** acetone, chloroform, toluene
- Storage:** Store in absence of light, desiccated and refrigerate

Description

- Hydrophilic, alkyne-reactive, long-lifetime reagent for click chemistry containing one azide function. Azides react with C≡C-triple bonds in either a Cu(I)-catalyzed or Cu-free 1,3-dipolar cycloaddition reaction to triazole.

Applications

- Click Chemistry reagent
- Fluorescence intensity and fluorescence polarization-based applications
- Resonance Energy Transfer (RET)

Advantages

- Perfectly suited for excitation with the 635-nm and 650-nm diode lasers
- Sensitive; high extinction coefficients and high quantum yields after covalent attachment to biomolecules
- Quantum yield is highly increased after covalent and non-covalent association with proteins
- pH-insensitive between pH 3 and pH 10
- Good aqueous solubility:** this label does not alter the solubility of the bioconjugate
- Photostability:** Higher photostability as compared to **Alexa Fluor™ 647** or **Cy5™**
- Low molecular weight:** **Seta** dyes do not add substantial mass to the conjugates
- Ideal for non-radioactive labeling of alkyne-modified proteins, DNA and oligonucleotides

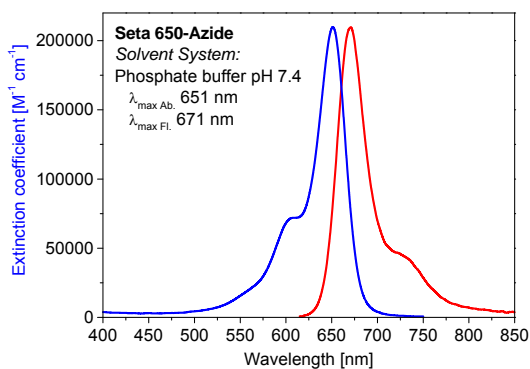
Spectral Data

Solvent System: phosphate buffer pH 7.4

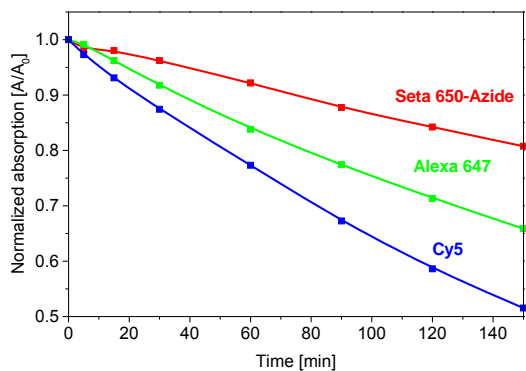
Sample	Absorption max. [nm]	Extinction Coefficient [M ⁻¹ cm ⁻¹]	Fluorescence max. [nm]	Quantum Yield ¹ [%]
Free dye	651	200,000	671	28

¹ Excitation at 610 nm

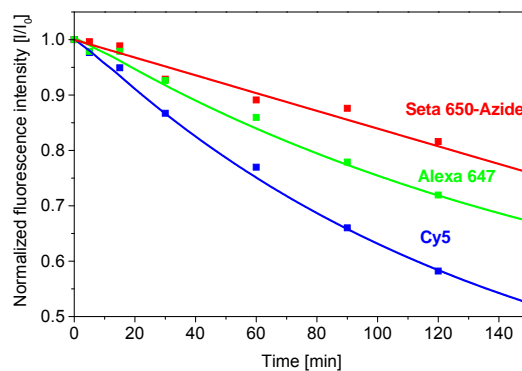
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Absorption and emission spectrum of a **Seta-650-Azide** in phosphate buffer (pH 7.4)



Decrease of the long-wavelength absorption band of **Seta-650-Azide** compared to **Cy5** and **Alexa Fluor 647**



Decrease of the fluorescence intensity of **Seta-650-Azide** compared to **Cy5** and **Alexa Fluor 647**