

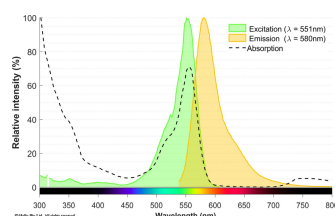
# DATASHEET

## Janelia Fluor® 549, Azide

### Product overview

<b>Name</b>	Janelia Fluor® 549, Azide
<b>Cat No</b>	HB7988
<b>Biological description</b>	<p>Cell-permeable, yellow fluorescent dye with an azide reactive group for copper-free click chemistry.</p> <p>Suitable for confocal microscopy and super resolution microscopy (SRM) including techniques such as dSTORM (both live and fixed cells) and STED. Also suitable for flow cytometry.</p> <p>Janelia Fluor® 549 is 2 x brighter than TMR and Cy3 <i>in vitro</i> and live-cell experiments.</p> <p><b>Spectrally similar dyes:</b> Alexa Fluor® 546, Alexa Fluor® 555, BDY TMR-X, Atto 550, CF 555, TAMRA, Cyanine 3</p>
<b>Alternative names</b>	JF549, Azide
<b>Biological action</b>	Dyes & stains
<b>Description</b>	Yellow dye supplied as an azide for click chemistry. Suitable for dSTORM, STED, confocal microscopy, live cell imaging and flow cytometry.

### Images



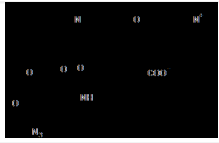
### Biological Data

<b>Application notes</b>	<p><b>#Protocol 1: Measurement of excitation and emission spectra of Janelia Fluor® 549, azide</b></p> <ul style="list-style-type: none"><li>Janelia Fluor® 549, azide was prepared at 1µm in PBS.</li><li>Spectra were generated on a Tecan Infinite M200 PRO using the following parameters:<ul style="list-style-type: none"><li>Excitation: Recording at 638nm while exciting between 280nm and 610nm</li><li>Emission: Exciting at 509nm while recording between 535nm and 800nm</li><li>Absorbance: Measured between 300 and 800nm</li></ul></li></ul>
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### Solubility & Handling

<b>Storage instructions</b>	-20°C
<b>Solubility overview</b>	Soluble in DMSO
<b>Important</b>	This product is for RESEARCH USE ONLY and is not intended for therapeutic or diagnostic use. Not for human or veterinary use

## Chemical Data

Chemical name	3,6-Di-1-azetidinyI-9-[5-[[[2-[2-[2-azidoethoxy]ethoxy]ethoxy]ethyl]carbamoyl]-2-carboxyphenyl]xanthylum, inner salt
Molecular Weight	654.71
Chemical structure	
Molecular Formula	C <sub>35</sub> H <sub>38</sub> N <sub>6</sub> O <sub>7</sub>
PubChem identifier	137919860
SMILES	<chem>O=C(NCCOCCOCCOCCN=[N+]=[N-])C1=CC=C(C([O-])=O)C(C(C2=CC=C(N3CCC3)C=C2O4)=C(C=C/5)C4=CC5=[N+]6CCC\6)=C1</chem>
Source	Synthetic
InChiKey	XDSXVSGQYCACTI-UHFFFAOYSA-N
Licensing details	Sold under license from the Howard Hughes Medical Institute, Janelia Research Campus
Excitation	549 nm
Emission	571 nm

## References

### A general method to improve fluorophores for live-cell and single-molecule microscopy.

Grimm JB et al (2015) Nature methods 12

PubMedID [25599551](#)