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## DATASHEET

LUF7487

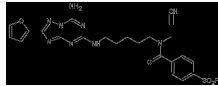
### Product overview

<b>Name</b>	LUF7487
<b>Cat No</b>	HB9585
<b>Biological description</b>	Novel, covalent Affinity-Based Probe (AfBP) which is suitable for click-conjugation (e.g. with the Cyanine-3 fluorophore) for visualization applications (e.g. SDS-PAGE). Irreversibly binds to the A <sub>2</sub> AR (apparent pK <sub>i</sub> values at A <sub>2a</sub> AR are 8.41 (after 0.5h) and 8.82 (after 3h), where a K <sub>i</sub> indicates a covalent mode of action).
<b>Biological action</b>	<i>Sold under license from the Oncode Institute and Universiteit Leiden</i> Other
<b>Purity</b>	>95%
<b>Description</b>	Novel, covalent Affinity-Based Probe (AfBP). Clickable antagonist for the A <sub>2</sub> AR.

### 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

<b>Chemical name</b>	4-[5-[[7-Amino-2-(furan-2-yl)-[1,2,4]triazolo[1,5-a][1,3,5]triazin-5-yl]amino]pentyl-prop-2-ynylcarbamoyl]benzenesulfonyl fluoride
<b>Molecular Weight</b>	526.5
<b>Chemical structure</b>	
<b>Molecular Formula</b>	C <sub>23</sub> H <sub>23</sub> FN <sub>8</sub> O <sub>4</sub> S
<b>PubChem identifier</b>	145953622
<b>SMILES</b>	<chem>C#CCN(CCCCCN1=NC2=NC(=NN2C(=N1)N)C3=CC=CO3)C(=O)C4=CC=C(C=C4)S(=O)(=O)F</chem>
<b>InChiKey</b>	KGDQUCQGDCWHAI-UHFFFAOYSA-N
<b>Licensing details</b>	Sold under license from the Oncode Cancer Institute and Universiteit Leiden

### References

#### An Affinity-Based Probe for the Human Adenosine A(2A) Receptor.

Yang X et al (2018) Journal of medicinal chemistry 61

**PubMedID** [30080404](#)