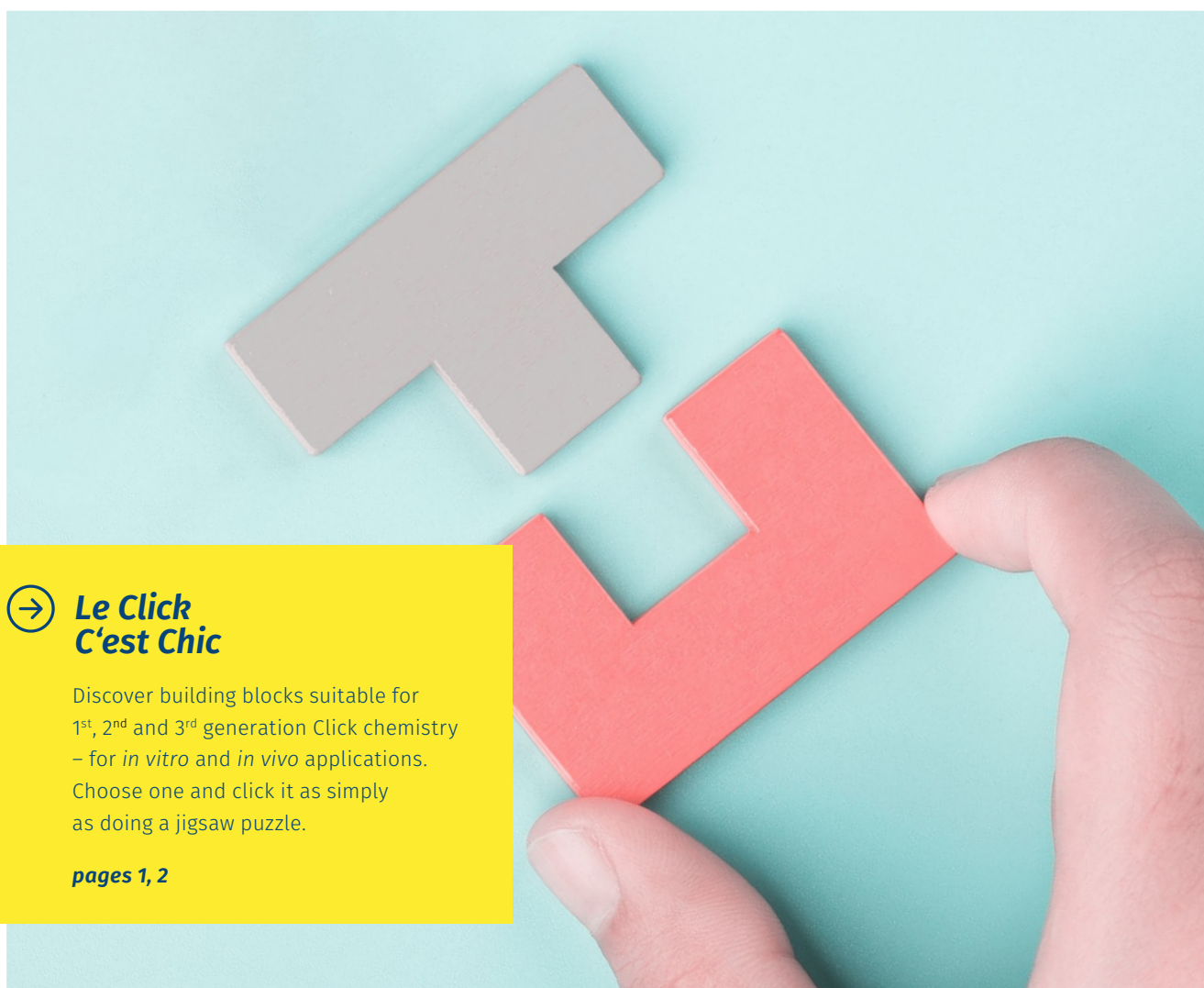




Iris
Biotech

CLICK CHEMISTRY

Three Generations



→ **Le Click C'est Chic**

Discover building blocks suitable for 1st, 2nd and 3rd generation Click chemistry – for *in vitro* and *in vivo* applications. Choose one and click it as simply as doing a jigsaw puzzle.

pages 1, 2

1st generation Click chemistry:
CuAAC

page 1

2nd generation Click chemistry:
SPAAC

page 1–2

3rd generation Click chemistry:
IEDDA

page 3



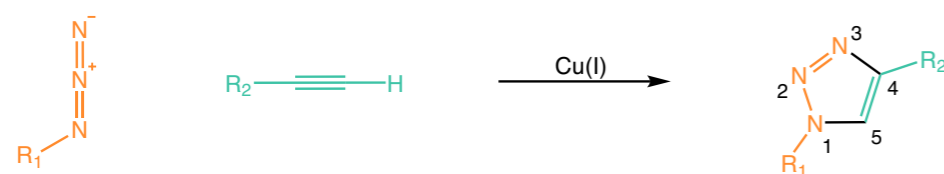
Version: IF4_3

Click Chemistry

Three Generations

1st Generation Click Chemistry: CuAAC

Alkynes and azides can undergo a Cu(I)-catalyzed azide-alkyne 1,3-dipolar cycloaddition (CuAAC) to afford 1,4-disubstituted 1,2,3-triazoles. Developed by K. Barry Sharpless and Morton Meldal, this type of chemical transformation was quickly dubbed "Click Chemistry". This reaction has found broad applicability in various fields and is as such currently the most widely used conjugation technique.



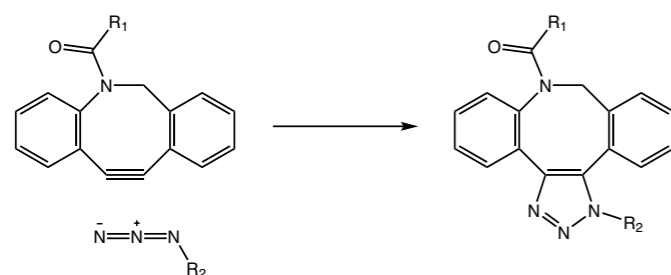
The copper-catalyzed azide-alkyne cycloaddition affords the 1,4 disubstituted isomer.

Worth noting is the fact that ruthenium is also able to catalyze a 1,3-dipolar cycloaddition between an azide and an alkyne affording the 1,5-disubstituted regioisomer instead.

However, the presence of copper limits *in vivo* applications due to high cell toxicity, undesired oxidation of proteins and the inhibition of luminescence properties of nanocrystals.

2nd Generation Click Chemistry: SPAAC

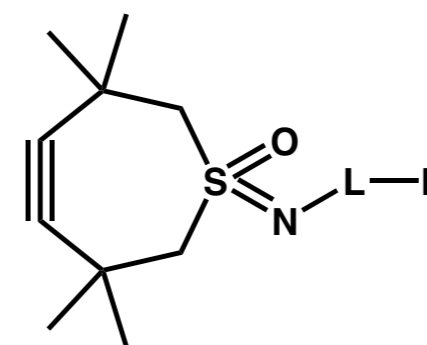
The so-called strain-promoted azide-alkyne cycloaddition (SPAAC) is widely utilized as copper-free click reaction in a broad range of research fields. However, when using cyclooctynes, their high reactivity comes with the drawback of causing thermal decomposition and/or oligomerization during storage and experimental handling. Furthermore, competitively occurring addition reaction with biological thiols, e.g. cysteine residues in proteins or peptides, diminish their biorthogonality. In contrast, the relatively new developed 4,8-diazacyclononynes (DACNs) possess high thermal and chemical stability while showing high reactivity and increased hydrophilicity as well as high selectivity towards ynophiles.



Schematic illustration of SPAAC.

CliCr®

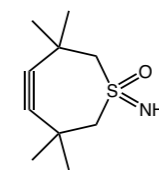

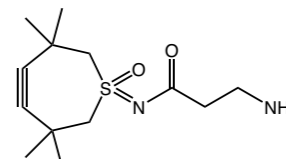

A superior class of reagents for metal-free click strain promoted conjugation with azides is based on CliCr®, which is based on the small-molecule TMTH-sulfoximine (TMTHSI). The 7-membered CliCr® ring can be conveniently functionalized with a variety of linkers, e.g. *via* acylation, sulfonylation, N-alkylation, or carbamoylation. Hence, CliCr® reagents can be used in diverse application, for example for the construction of Antibody-Drug Conjugates (ADCs), for *ex vivo* cell modification (e.g. CAR-T), for small molecule-drug conjugates, for oligonucleotide conjugates as well as for diagnostic labelling of a variety of agents.

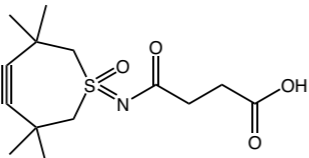

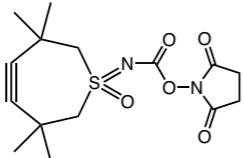



Chemical structure of the CliCr® base compound and its derivatization possibilities.

Key benefits are its fast reactivity and increased yield, its hydrophilicity compared to other metal-free click reagents combined with its small size as well as its broad compatibility.

CliCr® is provided under an intellectual property license from Cristal Therapeutics. The trademark CliCr® is the property of Cristal Therapeutics. For information on purchasing a license of CliCr® reagents, contact Cristal Therapeutics via Oxfordlaan 55, 6229 EV Maastricht (The Netherlands) or via info@crystaltherapeutics.com.

		Product details
RL-4180	CliCr® base compound	
	TMTH-Sulfoximine	
CAS-No.	2408481-82-1	
Formula	C ₁₀ H ₁₇ NOS	
Mol. weight	199,31 g/mol	
RL-4190	CliCr®-beta-Ala-NH₂*TFA	
	TMTH-sulfoximine beta-alanine amide TFA salt	
Formula	C ₁₃ H ₂₂ N ₂ O ₂ S*CF ₃ COOH	
Mol. weight	270,39*114,02 g/mol	

		Product details	
RL-4200	CliCr®-Suc		
TMTH-sulfoximine succinic acid			
CAS-No.	2479971-29-2		
Formula	C ₁₄ H ₂₁ NO ₅ S		
Mol. weight	299,39 g/mol		
RL-4330	CliCr®-Osu		
TMTH-sulfoximine succinimidyl ester			
CAS-No.	2408481-89-8		
Formula	C ₁₅ H ₂₀ N ₂ O ₅ S		
Mol. weight	340,39 g/mol		

References:

- *TMTHSI, a superior 7-membered ring alkyne containing reagent for strain-promoted azide-alkyne cycloaddition reactions*; J. Weterings, C. J. F. Rijcken, H. Veldhuis, T. Meulemans, D. Hadavi, M. Timmers, M. Honing, H. Ippel, R. M. J. Liskamp; **Chem. Sci.** 2020; **11**: 9011-9016. <https://doi.org/10.1039/d0sc03477k>
- *Exploring the Chemical Properties and Medicinal Applications of Tetramethylthiocycloheptyne Sulfoximine Used in Strain-Promoted Azide-Alkyne Cycloaddition Reactions*; M. Timmers, A. Kipper, R. Frey, S. Notermans, M. Voievudskyi, C. Wilson, N. Hentzen, M. Ringle, C. Bovino, B. Stump, C. J. F. Rijcken, T. Vermonden, I. Dijkgraaf, R. Liskamp; **Pharmaceuticals** 2023; **16**: 1155. <https://doi.org/10.3390/ph16081155>
- *Specific N-terminal attachment of TMTHSI linkers to native peptides and proteins for strain-promoted azide alkyne cycloaddition*; M. Timmers, W. Peeters, N. J. Hauwert, C. J. F. Rijcken, T. Vermonden, I. Dijkgraaf, R. M. J. Liskamp; **Chem. Commun.** 2023; <https://doi.org/10.1039/d3cc03397j>

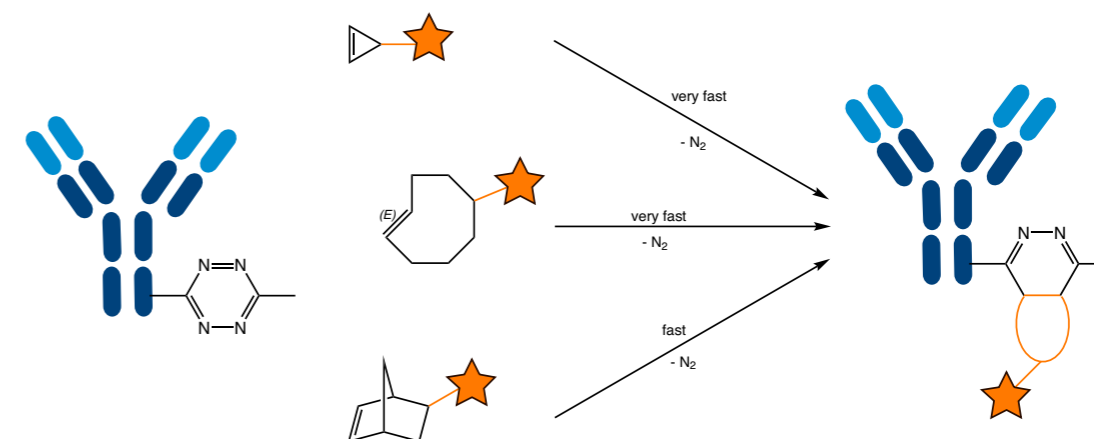


You need more information about the CliCr® technology?

Watch the recording of our online workshop!

**3rd Generation Click Chemistry: IEDDA**

Tetrazine ligation represents the option for a copper-free, rapid, and fully biorthogonal type of Click chemistry. Mechanistically, this reaction proceeds *via* an inverse electron-demand Diels-Alder (IEDDA) cycloaddition reaction between a tetrazine and a strained alkene, such as a *trans*-cyclooctene (TCO), cyclopropane or norbornene, followed by a retro-Diels-Alder reaction under elimination of N₂, the latter rendering the reaction irreversible. As additional benefit, this reaction excels at very low concentrations and can be performed in aqueous media.



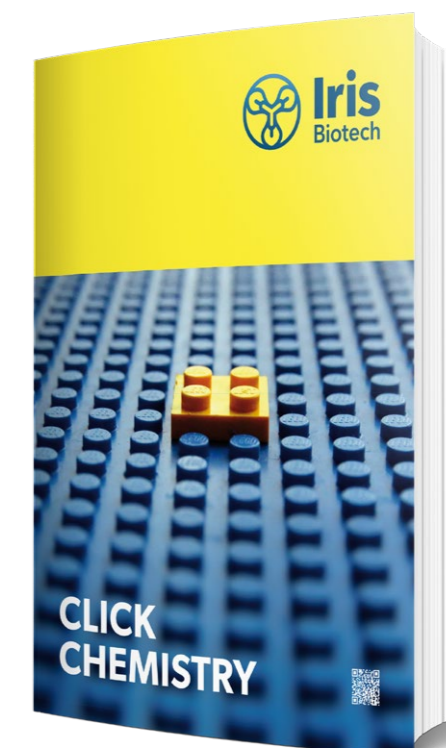
Common reaction partners for tetrazines.

There are two main types of tetrazines that are widely applied: 6-Methyl-substituted tetrazines (MeTz) and 6-hydrogen-substituted tetrazines (HTz). MeTz exhibit a high stability while still offering faster reaction kinetics with TCO derivatives than any other biorthogonal reaction pairs (approx. 1000 M⁻¹s⁻¹). HTzs show lower stability and less tolerance to harsh reaction conditions but offer even faster reaction kinetics (up to 30000 M⁻¹s⁻¹) for applications like *in vivo* imaging.



For more detailed information on traditional and copper-free Click conjugation, please see our brochure Click Chemistry.

Free Download is available on our website!

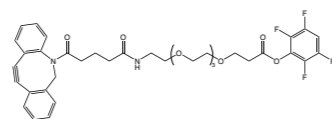


DBCO derivatives

PEG6740 DBCO-PEG(4)-TFP

Dibenzoazacyclooctyne-tetra(ethylene glycol)-propionyl 2,3,5,6-tetrafluorophenol ester

Formula $C_{37}H_{38}F_4N_2O_8$
Mol. weight 714,7 g/mol



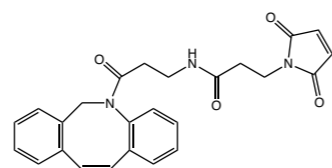
Product details



RL-2490 DBCO-mal

Dibenzoazacyclooctyne-maleimide

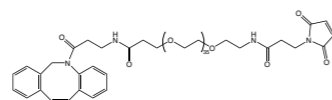
CAS-No. 1395786-30-7
Formula $C_{25}H_{21}N_3O_4$
Mol. weight 427,45 g/mol



PEG6785 DBCO-PEG(36)-MAL

Dibenzoazacyclooctyne-36(ethylene glycol)-maleimide

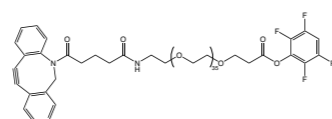
Formula $C_{100}H_{170}N_4O_{41}$
Mol. weight 2084,45 g/mol



PEG6765 DBCO-PEG(36)-TFP

Dibenzoazacyclooctyne-36(ethylene glycol)-propionyl 2,3,5,6-tetrafluorophenol ester

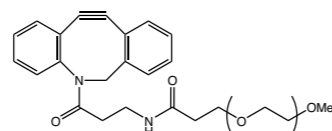
Formula $C_{101}H_{166}F_4N_2O_{40}$
Mol. weight 2124,41 g/mol



PEG7460 DBCO-PEG(24)-OMe

alpha-Methoxy-24(ethylene glycol)-amido-dibenzoazacyclooctyne

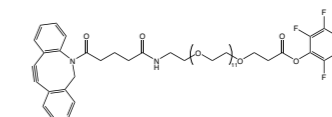
Formula $C_{68}H_{114}N_2O_{26}$
Mol. weight 1375,63 g/mol



PEG6750 DBCO-PEG(12)-TFP

Dibenzoazacyclooctyne-dodeca(ethylene glycol)-propionyl 2,3,5,6-tetrafluorophenol ester

Formula $C_{53}H_{70}F_4N_2O_{16}$
Mol. weight 1067,12 g/mol

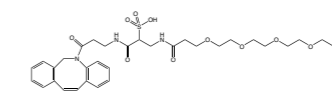


Product details

RL-2421 DBCO-Sulfo-PEG(4)-NH₂

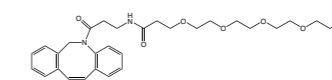
Dibenzoazacyclooctyne-tetra(ethylene glycol)amine

CAS-No. 2055198-05-3
Formula $C_{32}H_{42}N_4O_{10}S$
Mol. weight 674,76 g/mol

RL-2420 DBCO-PEG(4)-NH₂*TFA

Dibenzoazacyclooctyne-tetra(ethylene glycol)-amine trifluoro acetic acid salt

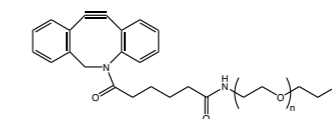
CAS-No. 1255942-08-5 net
Formula $C_{29}H_{37}N_3O_6 \cdot C_2F_3HO_2$
Mol. weight 523,62*114,02 g/mol



RL-2550 DBCO-mPEG (20kDa)

alpha-Dibenzoazacyclooctyne-omega-methoxy-poly(ethylene glycol)

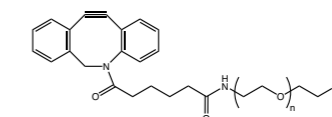
CAS-No. 2262541-53-5
Mol. weight 20000 Da



RL-2540 DBCO-mPEG (10kDa)

alpha-Dibenzoazacyclooctyne-omega-methoxy-poly(ethylene glycol)

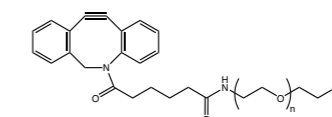
CAS-No. 2262541-53-5
Mol. weight 10000 Da



RL-2530 DBCO-mPEG (5kDa)

alpha-Dibenzoazacyclooctyne-omega-methoxy-poly(ethylene glycol)

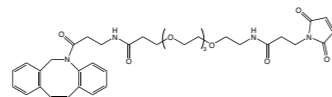
CAS-No. 2262541-53-5
Mol. weight 5000 Da



RL-2500 DBCO-PEG(4)-mal

Dibenzoazacyclooctyne-tetra(ethylene glycol)-maleimide

CAS-No. 1480516-75-3
 Formula $C_{36}H_{42}N_4O_9$
 Mol. weight 674,74 g/mol

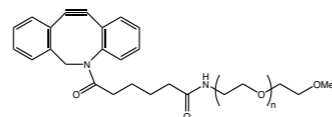


Product details

RL-2560 DBCO-mPEG (30kDa)

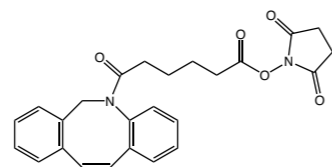
alpha-Dibenzoazacyclooctyne-omega-methoxy-poly(ethylene glycol)

CAS-No. 2262541-53-5
 Mol. weight 30000 Da

**RL-2440 DBCO-NHS**

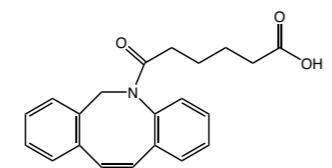
Dibenzoazacyclooctyne-carboxylic acid succinimidyl ester

CAS-No. 1384870-47-6
 Formula $C_{25}H_{22}N_2O_5$
 Mol. weight 430,45 g/mol

**RL-2430 DBCO-COOH**

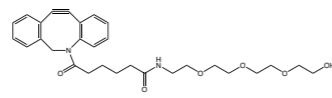
Dibenzoazacyclooctyne-carboxylic acid

CAS-No. 1425485-72-8
 Formula $C_{21}H_{13}NO_3$
 Mol. weight 333,38 g/mol

**RL-2510 DBCO-PEG(4)-OH**

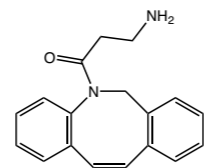
Dibenzoazacyclooctyne-tetra(ethylene glycol)

CAS-No. 1416711-60-8
 Formula $C_{29}H_{36}N_2O_6$
 Mol. weight 508,61 g/mol

**RL-2120 DBCO-NH₂**

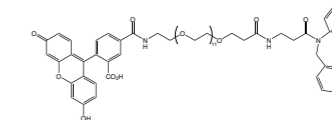
Dibenzocyclooctyne-amine

CAS-No. 1255942-06-3
 Formula $C_{18}H_{16}N_2O$
 Mol. weight 276,33 g/mol

**PEG6830 DBCO-dPEG(12)-(5)6-carboxyfluorescein**

Dibenzoazacyclooctyne-dodeca(ethylene glycol)-(5)6-carboxyfluorescein

Formula $C_{46}H_{39}N_3O_{10}$
 Mol. weight 1234,34 g/mol

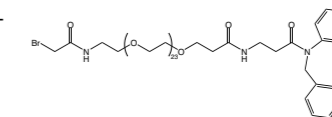


Product details

PEG6810 Bromoacetamido-PEG(24)-DBCO

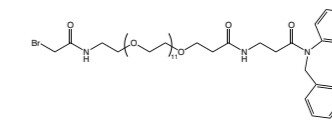
Bromoacetamido-24(ethylene glycol)-amido-dibenzoazacyclooctyne

Formula $C_{71}H_{118}BrN_3O_{27}$
 Mol. weight 1525,6 g/mol

**PEG6800 Bromoacetamido-PEG(12)-DBCO**

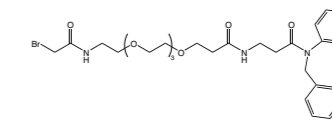
Bromoacetamido-dodeca(ethylene glycol)-amido-dibenzoazacyclooctyne

Formula $C_{47}H_{70}BrN_3O_{17}$
 Mol. weight 996,97 g/mol

**PEG6790 Bromoacetamido-PEG(4)-DBCO**

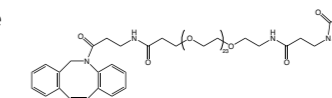
Bromoacetamido-tetra(ethylene glycol)-amido-dibenzoazacyclooctyne

Formula $C_{31}H_{38}BrN_3O_7$
 Mol. weight 644,55 g/mol

**PEG6780 DBCO-PEG(24)-MAL**

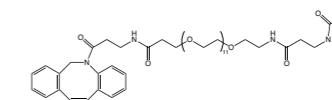
Dibenzoazacyclooctyne-24(ethylene glycol)-maleimide

Formula $C_{76}H_{122}N_4O_{29}$
 Mol. weight 1555,79 g/mol

**PEG6770 DBCO-PEG(12)-MAL**

Dibenzoazacyclooctyne-dodeca(ethylene glycol)-maleimide

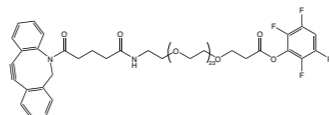
CAS-No. 2011777-01-6
 Formula $C_{52}H_{74}N_4O_{17}$
 Mol. weight 1027,16 g/mol



PEG6760 DBCO-PEG(24)-TFP

Dibenzoazacyclooctyne-24(ethylene glycol)-propionyl 2,3,5,6-tetrafluorophenol ester

CAS-No. 2754372-40-0
 Formula $C_{77}H_{118}F_4N_2O_{28}$
 Mol. weight 1595,75 g/mol

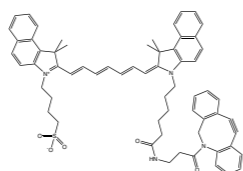


Product details

**RL-2870 ICG-DBCO**

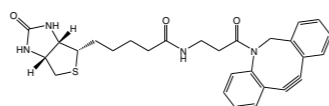
Indocyanine green dibenzoazacyclooctyne

Formula $C_{63}H_{64}N_4O_5S$
 Mol. weight 989,27 g/mol

**LS-4270 Biotin-DBCO**

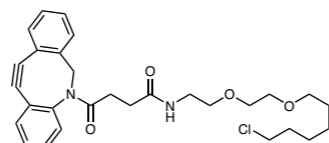
(3aS,4S,6aR)-N-[3-(11,12-Didehydridibenz[b,f]azocin-5(6H)-yl)-3-oxopropyl]hexahydro-2-oxo-1H-thieno[3,4-d]imidazole-4-pentanamide

CAS-No. 1418217-95-4
 Formula $C_{28}H_{30}N_4O_3S$
 Mol. weight 502,63 g/mol

**RL-3670 Halo-DBCO**

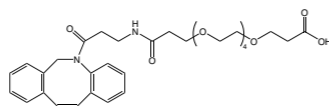
N-[2-[2-[(6-chlorohexyl)oxy]ethoxy]ethyl]-gamma-oxo-dibenz[b,f]azocine-5(6H)-butanamide

CAS-No. 1808119-16-5
 Formula $C_{29}H_{35}ClN_2O_4$
 Mol. weight 511,06 g/mol

**RL-2450 DBCO-PEG(5)-COOH**

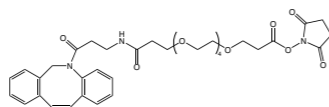
Dibenzoazacyclooctyne-penta(ethylene glycol)-propionic acid

Formula $C_{32}H_{40}N_2O_9$
 Mol. weight 596,67 g/mol

**RL-2460 DBCO-PEG(5)-NHS**

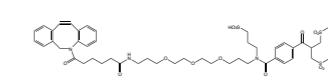
Dibenzoazacyclooctyne-penta(ethylene glycol)-propionic acid succinimidyl ester

CAS-No. 1378531-80-6
 Formula $C_{36}H_{43}N_3O_{11}$
 Mol. weight 693,74 g/mol

**RL-2480 DBCO-PEG(3)-BisSulfonThiol-Linker**

Dibenzoazacyclooctyne-PEG(3)-BisSulfon-Thiol-Linker

Formula $C_{59}H_{69}N_3O_{14}S_3$
 Mol. weight 1140,39 g/mol

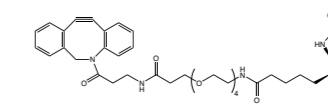


Product details

**RL-2520 Biotin-PEG(4)-DBCO**

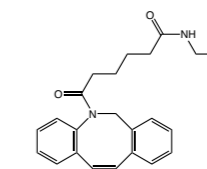
Dibenzoazacyclooctyne-tetra(ethylene glycol)-biotin

CAS-No. 1255942-07-4
 Formula $C_{39}H_{51}N_5O_8S$
 Mol. weight 749,92 g/mol

**RL-4020 DBCO-C6-Alkyne**

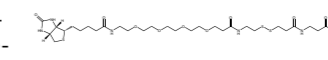
N-(propargylamidoadipoyl)-dibenzoazacyclooctyne

Formula $C_{24}H_{22}N_2O_2$
 Mol. weight 370,45 g/mol

**PEG8120 Biotin-PEG(4)-SS-DBCO**

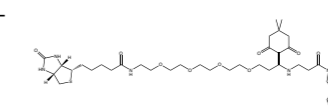
N-(2-((3-(3-(azadibenzocyclooctyn-1-yl)-3-oxopropylamino)-3-oxopropyl)disulfanyl)ethyl)-1-(5-((3aS,4S,6aR)-2-oxohexahydro-1H-thieno[3,4-d]imidazol-4-yl)pentanamido)-3,6,9,12-tetraoxapentadecan-15-amide

Formula $C_{44}H_{60}N_6O_9S_3$
 Mol. weight 913,18 g/mol

**PEG8140 Biotin-PEG(4)-Dde-DBCO**

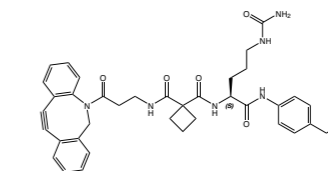
N-(15-(4,4-dimethyl-2,6-dioxocyclohexylidene)-19-oxo-19-(azadibenzocyclooctyn-1-yl)-3,6,9,12-tetraoxa-16-azanonadecyl)-5-((3aS,4S,6aR)-2-oxohexahydro-1H-thieno[3,4-d]imidazol-4-yl)pentanamide

CAS-No. 1807512-43-1
 Formula $C_{47}H_{61}N_5O_9S$
 Mol. weight 872,08 g/mol

**ADC1520 DBCO-cyclobutane-1,1-dicarboxamide-Cit-PAB**

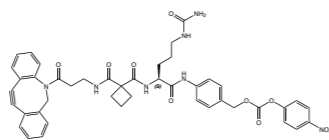
dibenzoazacyclooctyne-cyclobutane-1,1-dicarboxamide-citrullyl-(4-aminobenzyl alcohol)

CAS-No. 2576471-51-5
 Formula $C_{37}H_{40}N_6O_6$
 Mol. weight 664,75 g/mol



ADC1530 DBCO-cyclobutane-1,1-dicarboxamide-Cit-PAB-PNP

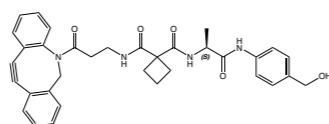
dibenzoazacyclooctyne-cyclobutane-1,1-dicarboxamide-citrullyl-(4-aminobenzyl)-(4-nitrophenyl)-carbonate

 CAS-No. 2576471-34-4
 Formula $C_{44}H_{43}N_7O_{10}$
 Mol. weight 829,85 g/mol


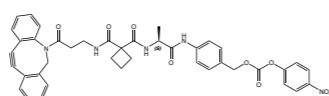
Product details

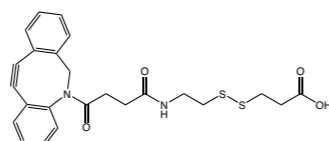

ADC1620 DBCO-cyclobutane-1,1-dicarboxamide-Ala-PAB

dibenzoazacyclooctyne-cyclobutane-1,1-dicarboxamide-alanyl-(4-aminobenzyl alcohol)

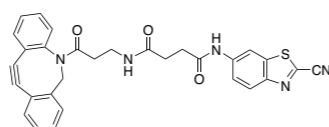
 CAS-No. 2576471-46-8
 Formula $C_{34}H_{34}N_4O_5$
 Mol. weight 578,66 g/mol

ADC1630 DBCO-cyclobutane-1,1-dicarboxamide-Ala-PAB-PNP

dibenzoazacyclooctyne-cyclobutane-1,1-dicarboxamide-alanyl-(4-aminobenzyl)-(4-nitrophenyl)-carbonate

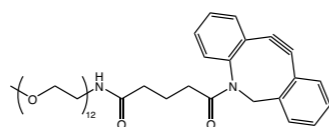
 CAS-No. 2576471-43-5
 Formula $C_{41}H_{37}N_5O_9$
 Mol. weight 743,76 g/mol

RL-4110 DBCO-Suc-SS-COOH

 CAS-No. 2749426-25-1
 Formula $C_{24}H_{24}N_2O_4S_2$
 Mol. weight 468,59 g/mol

RL-4310 DBCO-Suc-CBT

N1-(2-cyanobenzo[d]thiazol-6-yl)-N4-(3-(11,12-didehydro-5,6-dihydro-dibenzo[b,f]azocin-yl)-3-oxopropyl) succinamide

 Formula $C_{30}H_{23}N_5O_3S$
 Mol. weight 533,61 g/mol

PEG7465 Me-PEG(12)-DBCO

Methyl-12(ethylene glycol)-amido-dibenzoazacyclooctyne

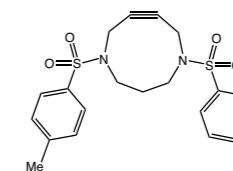
 Formula $C_{45}H_{68}N_2O_{14}$
 Mol. weight 861,04 g/mol

RL-4010 DBCO-SNAP

 Formula $C_{34}H_{31}N_2O_3$
 Mol. weight 585,67 g/mol

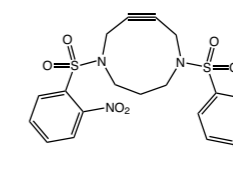

Product details


DACN derivatives
RL-2730 DACN(Tos2)

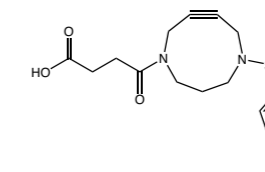
N,N'-bis(p-toluenesulfonyl)-4,8-diazacyclononyne

 CAS-No. 1797508-57-6
 Formula $C_{21}H_{24}N_2O_4S_2$
 Mol. weight 432,56 g/mol

RL-2710 DACN(Tos,Ns)

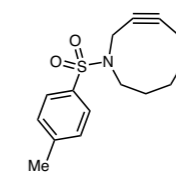
N-(o-nitrobenzenesulfonyl)-N'-(p-toluenesulfonyl)-4,8-diazacyclononyne

 CAS-No. 1797508-58-7
 Formula $C_{20}H_{21}N_2O_6S_2$
 Mol. weight 463,53 g/mol

RL-2720 DACN(Tos,Suc-OH)

N-succinoyl-N'-(p-toluenesulfonyl)-4,8-diazacyclononyne

 CAS-No. 2109751-68-8
 Formula $C_{18}H_{22}N_2O_5S$
 Mol. weight 378,44 g/mol

RL-2735 DACN(Tos)*HCl

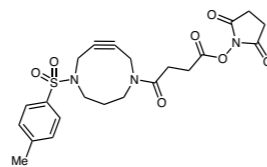
N-(p-toluenesulfonyl)-4,8-diazacyclononyne hydrochloride

 CAS-No. 2331322-18-8
 Formula $C_{14}H_{18}N_2O_2S^*HCl$
 Mol. weight 278,37*36,46 g/mol


RL-2725 DACN(Tos,Suc-NHS)

N-(succinoyl-NHS ester)-N'-(p-toluenesulfonyl)-4,8-diazacyclononyne

CAS-No. 2411082-26-1
Formula $C_{22}H_{25}N_3O_5S$
Mol. weight 475,52 g/mol

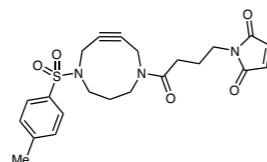


Product details


RL-3630 DACN(Tos,Mal)

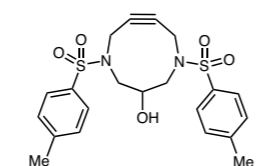
N-(maleimidobutyryl)-N'-(p-toluenesulfonyl)-4,8-diazacyclononyne

CAS-No. 2411082-28-3
Formula $C_{22}H_{25}N_3O_5S$
Mol. weight 443,52 g/mol


RL-2737 DACN(Tos2,6-OH)

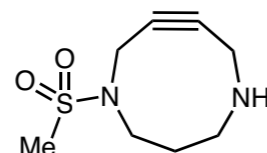
4,8-Bis(p-toluenesulfonyl)-4,8-diazacyclononyne-6-ol

CAS-No. 2109751-74-6
Formula $C_{21}H_{24}N_2O_5S_2$
Mol. weight 448,55 g/mol


RL-3600 DACN(Ms)*HCl

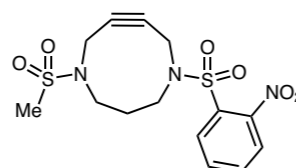
N-(Mesyl)-4,8-diazacyclononyne hydrochloride

CAS-No. 2331322-16-6
Formula $C_8H_{14}N_2O_2S^+HCl$
Mol. weight 202,27*36,46 g/mol


RL-3610 DACN(Ms,Ns)

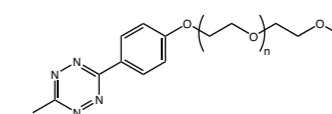
N-(Mesyl)-N'-(2-nosyl)-4,8-diazacyclononyne

CAS-No. 2411082-25-0
Formula $C_{14}H_{17}N_3O_6S_2$
Mol. weight 387,43 g/mol


Tetrazine derivatives
RL-2390 MeTz-mPEG (10kDa)

alpha-Methyltetrazine-omega-methoxy-poly(ethylene glycol)

Mol. weight 10000 Da

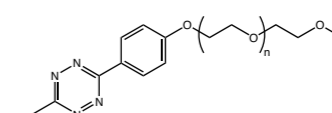


Product details


RL-2400 MeTz-mPEG (20kDa)

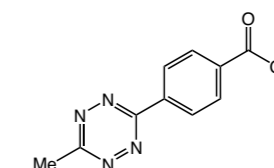
alpha-Methyltetrazine-omega-methoxy-poly(ethylene glycol)

Mol. weight 20000 Da


RL-2130 (Me)Tz-benzoic acid

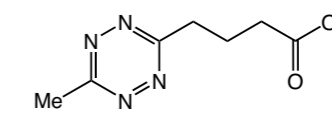
4-(6-methyl-1,2,4,5-tetrazin-3-yl)benzoic acid

CAS-No. 1345866-66-1
Formula $C_{10}H_8N_4O_2$
Mol. weight 216,2 g/mol


RL-2140 (Me)Tz-butanoic acid

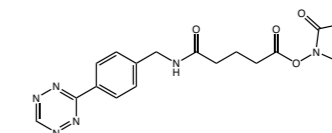
4-(6-methyl-1,2,4,5-tetrazin-3-yl)butanoic acid

CAS-No. 1923268-81-8
Formula $C_7H_{10}N_4O_2$
Mol. weight 182,18 g/mol


RL-2240 Bz-Tz-NHS

2,5-dioxopyrrolidin-1-yl 5-(4-(1,2,4,5-tetrazin-3-yl)benzylamino)-5-oxopentanoate

CAS-No. 1244040-64-9
Formula $C_{18}H_{18}N_6O_5$
Mol. weight 398,37 g/mol

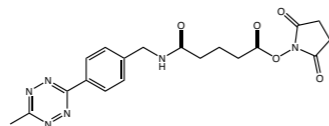


RL-2230 Bz-(Me)Tz-NHS

2,5-dioxopyrrolidin-1-yl 5-(4-(6-methyl-1,2,4,5-tetrazin-3-yl)benzylamino)-5-oxopentanoate

CAS-No. 1454558-58-7

Mol. weight 412,41 g/mol



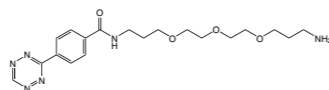
Product details

**RL-2590 Tz-benzoyl-TOTA*TFA**

Tz-benzoyl-TOTA*TFA

Formula $C_{19}H_{28}N_6O_4 \cdot C_2H_3F_3O_2$

Mol. weight 404,46*114,02 g/mol

**RL-2250 Bz-Tz-PEG(5)-NHS**

2,5-dioxopyrrolidin-1-yl 1-(4-(1,2,4,5-tetrazin-3-yl)phenyl)-3-oxo-6,9,12,15,18-pentaoxa-2-azahenicosan-21-oate

CAS-No. 1682653-80-0

Formula $C_{27}H_{36}N_6O_{10}$

Mol. weight 604,61 g/mol

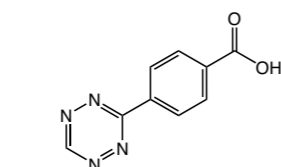
**RL-2580 Tz-benzoic acid**

4-(1,2,4,5-tetrazin-3-yl)benzoic acid

CAS-No. 1345866-65-0

Formula $C_8H_6N_4O_2$

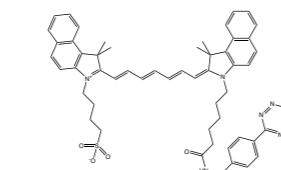
Mol. weight 202,17 g/mol

**RL-2860 ICG-Tz**

Indocyanine green tetrazine

Formula $C_{54}H_{57}N_7O_4S$

Mol. weight 900,14 g/mol

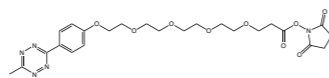
**RL-2330 MeTz-PEG(4)-NHS**

Methyltetrazine-PEG(4)-propanoyl succinimidyl ester

CAS-No. 1802907-92-1

Formula $C_{24}H_{31}N_5O_9$

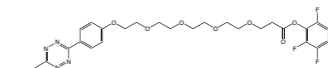
Mol. weight 533,53 g/mol

**RL-3905 MeTz-PEG(4)-STP**

sodium 2,3,5,6-tetrafluoro-4-((1-(4-(6-methyl-1,2,4,5-tetrazin-3-yl)phenoxy)-3,6,9,12-tetraoxapentadecan-15-oyl)oxy)benzenesulfonate

Formula $C_{26}H_{27}F_4N_4NaO_{10}S$

Mol. weight 686,56 g/mol



Product details

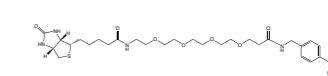
**LS-4290 Biotin-PEG(4)-MeTz**

N-(4-(6-methyl-1,2,4,5-tetrazin-3-yl)benzyl)-1-(5-((3aS,4S,6aR)-2-oxohexahydro-1H-thieno[3,4-d]imidazol-4-yl)pentanamido)-3,6,9,12-tetraoxapentadecan-15-amide

CAS-No. 1962919-31-8

Formula $C_{31}H_{46}N_8O_5S$

Mol. weight 674,82 g/mol

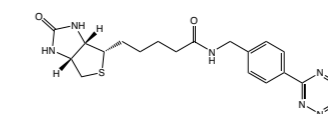
**LS-4280 Biotin-MeTz**

N-(4-(6-methyl-1,2,4,5-tetrazin-3-yl)benzyl)-5-((3aS,4S,6aR)-2-oxohexahydro-1H-thieno[3,4-d]imidazol-4-yl)pentanamide

CAS-No. 1802883-51-7

Formula $C_{26}H_{25}N_7O_2S$

Mol. weight 427,53 g/mol

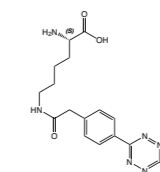
**HAA9170 H-Lys(MeTz-PhAc)-OH*TFA**

N-(2-(4-(6-methyl-1,2,4,5-tetrazin-3-yl)phenyl)acetyl)-L-lysine TFA salt

CAS-No. 2578384-82-2 (net)

Formula $C_{17}H_{22}N_6O_3 \cdot CF_3COOH$

Mol. weight 358,40*114,02 g/mol

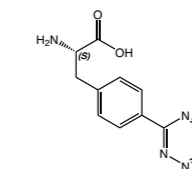
**HAA9470 H-L-Phe(4-MeTz)-OH*TFA**

(S)-2-amino-3-(4-(6-methyl-1,2,4,5-tetrazin-3-yl)phenyl)propanoic acid trifluoroacetic acid salt

CAS-No. 1698038-85-5 net

Formula $C_{12}H_{13}N_5O_2 \cdot CF_3COOH$

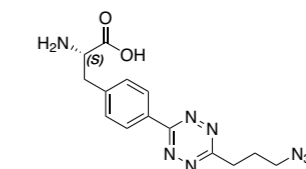
Mol. weight 259,27*114,02 g/mol

**HAA9480 H-L-Phe(4-Azido-PrTz)-OH*TFA**

(S)-2-amino-3-(4-(6-(3-azidopropyl)-1,2,4,5-tetrazin-3-yl)phenyl)propanoic acid trifluoroacetic acid salt

Formula $C_{14}H_{16}N_8O_2 \cdot CF_3COOH$

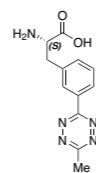
Mol. weight 328,34*114,02 g/mol



HAA9490 H-L-Phe(3-MeTz)-OH*TFA

(S)-2-amino-3-(3-(6-methyl-1,2,4,5-tetrazin-3-yl)phenyl)propanoic acid trifluoroacetic acid salt

CAS-No. 2036323-75-6 net
 Formula $C_{12}H_{13}N_5O_2 \cdot CF_3COOH$
 Mol. weight 259,27*114,02 g/mol

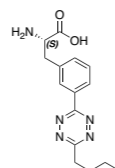


Product details

**HAA9500 H-L-Phe(3-BuTz)-OH*TFA**

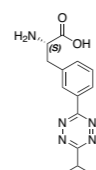
(S)-2-amino-3-(3-(6-butyl-1,2,4,5-tetrazin-3-yl)phenyl)propanoic acid trifluoroacetic acid salt

CAS-No. 2036323-83-6 net
 Formula $C_{15}H_{19}N_5O_2 \cdot CF_3COOH$
 Mol. weight 301,35*114,02 g/mol

**HAA9510 H-L-Phe(3-iPrTz)-OH*TFA**

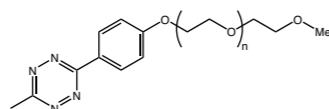
(S)-2-amino-3-(3-(6-isopropyl-1,2,4,5-tetrazin-3-yl)phenyl)propanoic acid trifluoroacetic acid salt

CAS-No. 2421119-12-0 net
 Formula $C_{14}H_{17}N_5O_2 \cdot CF_3COOH$
 Mol. weight 287,32*114,02 g/mol

**RL-2380 MeTz-mPEG (5kDa)**

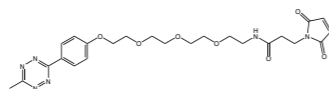
alpha-Methyltetrazine-omega-methoxy-poly(ethylene glycol)

Mol. weight 5000 Da

**RL-2340 MeTz-PEG(4)-mal**

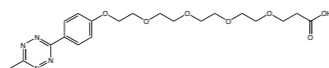
Methyltetrazine-PEG(4)-maleimide

CAS-No. 1802908-02-6
 Formula $C_{24}H_{30}N_6O_7$
 Mol. weight 514,53 g/mol

**RL-2310 MeTz-PEG(4)-COOH**

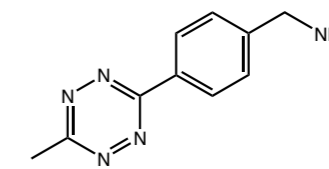
Methyltetrazine-PEG(4)-acid

CAS-No. 1802907-91-0
 Formula $C_{20}H_{28}N_4O_7$
 Mol. weight 436,56 g/mol

**RL-2360 MeTz-Bzl-NH₂*HCl**

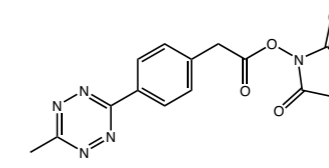
Methyltetrazine-benzylamine*HCl

CAS-No. 1596117-29-1
 Formula $C_{10}H_{11}N_5 \cdot HCl$
 Mol. weight 201,23*36,46 g/mol

**RL-2320 MeTz-PhAc-NHS**

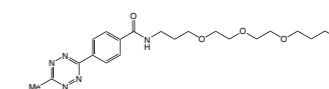
Methyltetrazine-phenylacetyl succinimidyl ester

CAS-No. 1644644-96-1
 Formula $C_{15}H_{13}N_5O_4$
 Mol. weight 327,29 g/mol

**RL-2110 MeTz-Phenyl-TOTA*TFA**

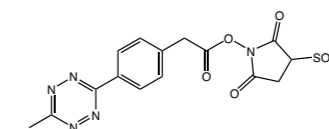
4-(6-methyl-1,2,4,5-tetrazin-3-yl)-N-(4,7,10-trioxatri-decane-13-amine)benzamide trifluoroacetate salt

CAS-No. 2250433-74-8 (net)
 Formula $C_{20}H_{30}N_6O_4 \cdot CF_3CO_2H$
 Mol. weight 418,49*114,02 g/mol

**RL-3915 MeTz-PhAc-Sulfo-NHS**

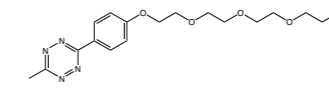
sodium 1-(2-(4-(6-methyl-1,2,4,5-tetrazin-3-yl)phenyl)acetoxyl)-2,5-dioxopyrrolidine-3-sulfonate

CAS-No. 1821017-46-2
 Formula $C_{15}H_{12}N_5NaO_7$
 Mol. weight 429,34 g/mol

**RL-2370 MeTz-PEG(4)-NH₂*HCl**

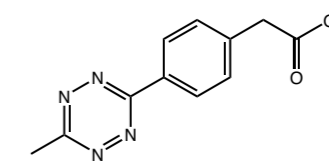
Methyltetrazine-PEG(4)-amine HCL salt

CAS-No. 1802908-05-9 net
 Formula $C_{17}H_{25}N_5O_4 \cdot HCl$
 Mol. weight 363,41*HCl g/mol

**RL-2300 MeTz-PhAcOH**

Methyltetrazine-phenylacetic acid

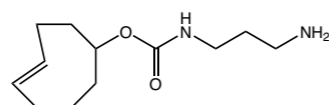
CAS-No. 1380500-88-8
 Formula $C_{11}H_{10}N_4O_2$
 Mol. weight 230,22 g/mol



TCO derivatives

TCO1060 TCO-NH₂*HCl*trans*-Cyclooctene-amine hydrochloride

CAS-No. 1800507-94-1
 Formula C₁₂H₂₂N₂O₂*HCl
 Mol. weight 226,32*36,45 g/mol



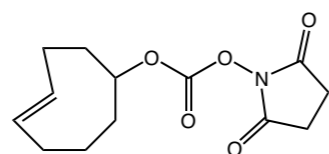
Product details



TCO1000 TCO-NHS

trans-Cyclooctene succinimidyl carbonate

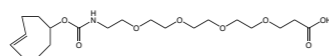
CAS-No. 1191901-33-3
 Formula C₁₃H₁₇NO₅
 Mol. weight 267,28 g/mol



TCO1040 TCO-PEG(4)-COOH

trans-Cyclooctene-PEG(4)-Acid

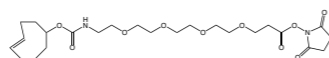
CAS-No. 1802913-21-8
 Formula C₂₀H₃₅NO₈
 Mol. weight 417,49 g/mol



TCO1010 TCO-PEG(4)-NHS

trans-Cyclooctene-PEG(4)-carboxy succinimidyl ester

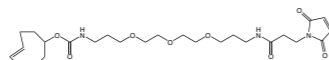
CAS-No. 1621096-79-4
 Formula C₂₄H₃₈N₂O₁₀
 Mol. weight 514,57 g/mol



TCO1050 TCO-PEG(3)-mal

trans-Cyclooctene-PEG(3)-maleimide

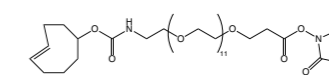
CAS-No. 1809356-72-6
 Formula C₂₆H₄₁N₃O₈
 Mol. weight 523,62 g/mol



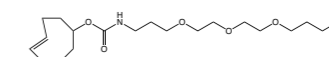
TCO1020 TCO-PEG(12)-NHS

trans-Cyclooctene-PEG(12)-carboxy succinimidyl ester

CAS-No. 2185016-39-9
 Formula C₄₀H₇₀N₂O₁₈
 Mol. weight 866,99 g/mol

TCO1070 TCO-PEG(3)-NH₂*HCl*trans*-Cyclooctene-PEG(3)-amine

CAS-No. 2028288-77-7
 Formula C₁₉H₃₆N₂O₅*HCl
 Mol. weight 372,51*36,46 g/mol

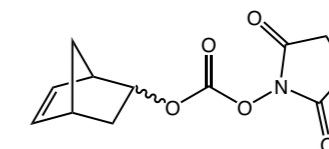


Norbornene derivatives

RL-2080 Norbornene-NHS

(Norbornene-2-yl)-N-hydroxysuccinimidylcarbonate

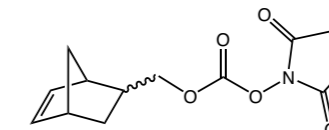
CAS-No. 1888335-48-5
 Formula C₁₂H₁₃NO₅
 Mol. weight 251,24 g/mol



RL-2090 Norbornene-methyl-NHS

(Norbornene-2-methyl)-N-hydroxysuccinimidylcarbonate

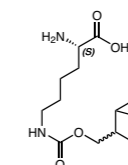
CAS-No. 1986791-87-0
 Formula C₁₃H₁₅NO₅
 Mol. weight 265,26 g/mol



HAA9235 H-L-Lys(Norbornene-methoxycarbonyl)-OH*HCl

N-epsilon-(norbornene-methoxycarbonyl)-L-lysine hydrochloride

CAS-No. 1378916-76-7 net
 Formula C₁₅H₂₄N₂O₄*HCl
 Mol. weight 296,37*36,46 g/mol



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Notes



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