Алматы (7273)495-231
Ангарск (3955)60-70-56
Архангельск (8182)63-90-72
Астрахань (8512)99-46-04
Барнаул (3852)73-04-60
Белгород (4722)40-23-64
Благовещенск (4162)22-76-07
Брянск (4832)59-03-52
Владивосток (423)249-28-31
Владикавказ (8672)28-90-48
Владимир (4922)49-43-18
Волгоград (844)278-03-48
Вологда (8172)26-41-59
Воронеж (473)204-51-73
Екатеринбург (343)384-55-89
Иваново (4932)77-34-06
Ижевск (3412)26-03-58
Иркутск (395)279-98-46
Казань (843)206-01-48
Калининград (4012)72-03-81
Калуга (4842)92-23-67
Кемерово (3842)65-04-62
Киров (8332)68-02-04
Коломна (4966)23-41-49
Кострома (4942)77-07-48
Краснодар (861)203-40-90
Красноярск (391)204-63-61
Курск (4712)77-13-04
Курган (3522)50-90-47
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Магнитогорск (3519)55-03-13
Москва (495)268-04-70
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Набережные Челны (8552)20-53-41
Нижний Новгород (831)429-08-12
Новокузнецк (3843)20-46-81
Ноябрьск (3496)41-32-12
Новосибирск (383)227-86-73
Омск (3812)21-46-40
Орел (4862)44-53-42
Оренбург (3532)37-68-04
Пенза (8412)22-31-16
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Псков (8112)59-10-37
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Симферополь (3652)67-13-56
Смоленск (4812)29-41-54
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Ставрополь (8652)20-65-13
Сургут (3462)77-98-35
Сыктывкар (8212)25-95-17
Тамбов (4752)50-40-97
Тверь (4822)63-31-35
Тольятти (8482)63-91-07
Томск (3822)98-41-53
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Тюмень (3452)66-21-18
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Уфа (347)229-48-12
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Череповец (8202)49-02-64
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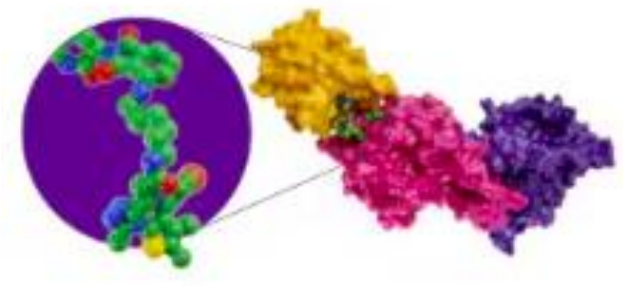
Технические характеристики на строительные блоки для разрушения белка, реагенты для клик-химии, смешивающие агенты и линкеры компании Sigma-Aldrich

Виды товаров: амины, карбоновые кислоты, азиды, алкины, гидроксилы, алкилгалогениды, бромы, малеимиды, эфиры, тиолы, дисульфиды, тетразины, библиотеки фрагментов, закодированных ДНК, целевые зонды селективные и характеризованные, биоортогональные репортеры теги, наборы для мониторинга пролиферации клеток, ингибиторы серингидролазы, азидные и алкиновые линкеры, дибензоциклооктины, тетразины или трансциклооктины, ацилтрифторбораты калия и др.

Protein Degrader Building Blocks



Targeted protein degradation (TPD) is a novel approach used in drug discovery to effectively infiltrate cell's proteasomal degradation pathway and quickly eliminate the target proteins from cells. This strategy has the potential to develop new therapeutics that differ considerably from classical inhibitors and is even viable for accessing "undruggable" protein targets and difficult-to-treat diseases.

Our diverse portfolio of protein degrader building blocks allows you to easily generate libraries of protein degraders that can be screened for an effective degradation of your target proteins from the cells. Our protein degraders, such as proteolysis targeting chimeras (PROTAC® molecules), are bifunctional molecules with three primary components:

Ligand at one end that targets the protein of interest (POI)

Second ligand at the opposite end that binds an E3 ligase

Crosslinker in the middle that joins the two ends

PROTEIN DEGRADER BUILDING BLOCK PORTFOLIO

Our protein degrader building blocks are permutations of the following components:

Ligands targeting the E3 ligase Cereblon (CRBN) or von Hippel–Lindau (VHL)

Crosslinkers with varied lengths and compositions

Conjugation sites with reactivity for common functional groups

ALL YOU NEED FOR SUCCESSFUL TARGETED PROTEIN DEGRADATION

The simultaneous degrader binding of two proteins brings the POI in close enough proximity for polyubiquitination to take place by the E2 enzyme associated to the E3 ligase, which flags the POI for degradation through the proteasome.

The precise design of small-molecule target degraders is crucial for successful protein degradation. Even slight alterations in ligands and crosslinkers can affect binding to the POI or E3 ligase. Thus, many analogs are synthesized, varying each structure slightly, and screened in cells to determine the optimal degrader for target degradation.

Our degrader building blocks are a collection of crosslinker-E3 ligand conjugates with a pendant functional group for covalent linkage to a target ligand to streamline synthesis. Since the same functional group is present across a series, one target ligand can be conjugated to several degrader building blocks simultaneously for facile library generation and subsequent screening.

ADVANTAGES OF OUR PROTEIN DEGRADERS

Compatibility: Linkers conjugate to common functional groups present on target ligands

Molecule design: The strategic variety of the combinations of linkers and ligands aids the design of target degraders

Synthetic time-saver: The E3 ligand-crosslinker conjugates decrease the amount of time spent on degrader synthesis

Library generation: Using degrader building blocks with the same conjugation site enables the simultaneous generation of several degraders via parallel synthesis

929352
(*S*,*R*,*S*)-AHPC-di-trimethylamide-dioxodisulfide-carbonate ester

930539
(*S*,*R*,*S*)-AHPC-Me-C_5-COOH
≥95%

T150
(−)-Thalidomide
>98%, solid

920703
((1-(*tert*-Butoxycarbonyl)piperidin-4-yl)methyl)proline

T151
(+)-Thalidomide
≥98% (HPLC), powder

T144
(±)-Thalidomide
≥98%, powder

744867
(1*R*,8*S*,9*s*)-Bicyclo[6.1.0]non-4-yn-9-ylmethyl *N*-succinimidyl carbonate
for Copper-free Click Chemistry

917397
(2-(Piperidin-4-yl)phenyl)methanamine dihydrochloride

761591
(4-(1,2,4,5-Tetrazin-3-yl)phenyl)methanamine hydrochloride
95%

764523
(*E*)-Cyclooct-4-enyl 2,5-dioxo-1-pyrrolidinyl carbonate

901490
(*S*,*R*,*S*)-AHPC hydrochloride
≥97%

918687
(*S*,*R*,*S*)-AHPC-alkyne-piperidine hydrochloride

920762
(*S*,*R*,*S*)-AHPC-benzyl-piperazine hydrochloride

901534
(*S*,*R*,*S*)-AHPC-C_6-CO_2H hydrochloride

905232
(*S*,*R*,*S*)-AHPC-C_6-PEG_3-butyl amine hydrochloride
≥95%

905380
(*S*,*R*,*S*)-AHPC-C_6-PEG_3-butyl chloride
≥95%

917818
(*S*,*R*,*S*)-AHPC-C_9-NH_2 hydrochloride
≥95%

920827
(*S*,*R*,*S*)-AHPC-methylamino-PEG_1-NH_2 hydrochloride

901503
(*S*,*R*,*S*)-AHPC-PEG_1-Alkyne

≥95%

903957
(*S*,*R*,*S*)-AHPC-PEG1-azide

901493
(*S*,*R*,*S*)-AHPC-PEG$_1$-NH$_2$ hydrochloride
≥95%

901517
(*S*,*R*,*S*)-AHPC-PEG$_2$-Alkyne
≥95%

904813
(*S*,*R*,*S*)-AHPC-PEG2-azide

906123
(*S*,*R*,*S*)-AHPC-PEG$_2$-butyl amine hydrochloride

905399
(*S*,*R*,*S*)-AHPC-PEG$_2$-butyl chloride

910600
(*S*,*R*,*S*)-AHPC-PEG$_2$-butyl CO$_2$H
≥95%

901488
(*S*,*R*,*S*)-AHPC-PEG$_2$-NH$_2$ hydrochloride
≥95%

901533
(*S*,*R*,*S*)-AHPC-PEG$_3$-Alkyne
≥95%

904651
(*S*,*R*,*S*)-AHPC-PEG$_3$-azide
≥95%

901511
(*S*,*R*,*S*)-AHPC-PEG$_3$-NH$_2$ hydrochloride
≥95%

901851
(*S*,*R*,*S*)-AHPC-PEG$_4$-Alkyne

901848
(*S*,*R*,*S*)-AHPC-PEG$_4$-NH$_2$ hydrochloride
≥95%

901855
(*S*,*R*,*S*)-AHPC-PEG$_5$-Alkyne

901850
(*S*,*R*,*S*)-AHPC-PEG$_5$-NH$_2$ hydrochloride

901873
(*S*,*R*,*S*)-AHPC-PEG$_6$-Alkyne

909378
(*S*,*R*,*S*)-AHPC-PEG$_6$-Azide
≥95%

905275
(*S*,*R*,*S*)-AHPC-PEG$_6$-butyl amine hydrochloride
≥95%

904864
(*S*,*R*,*S*)-AHPC-PEG$_6$-butyl azide
≥95%

901860
(*S*,*R*,*S*)-AHPC-PEG$_6$-NH$_2$ hydrochloride
≥95%

920606
(*S*,*R*,*S*)-AHPC-piperazine-pyridine-alkyne-NH$_2$ hydrochloride
921432
(*S*,*R*,*S*)-AHPC-pyridine-PEG$_1$-piperazine hydrochloride

921939
(*S*,*R*,*S*)-AHPC-pyrimidine-piperazine-PEG$_1$-NH$_2$ hydrochloride

920908
(*S*,*R*,*S*)-VL285 Phenol

920851
(*S*,*R*,*S*)-VL285 Phenol-C_2-NH_2 hydrochloride

922153
(*S*,*R*,*S*)-VL285 Phenol-C_3-piperazine hydrochloride

920878
(*S*,*R*,*S*)-VL285 Phenol-C_6-NH_2 hydrochloride

922161
(*S*,*R*,*S*)-VL285 Phenol-methylamino-PEG_1-NH_2 hydrochloride

920843
(*S*,*R*,*S*)-VL285 Phenol-PEG_1-piperazine hydrochloride

920894
(*S*,*R*,*S*)-VL285 Phenol-PEG_2-NH_2 hydrochloride

920886
(*S*,*R*,*S*)-VL285 Phenol-PEG_4-NH_2 hydrochloride

920835
(*S*,*R*,*S*)-VL285 Phenol-piperazine-pyridine-alkyne-NH_2 hydrochloride

901487
(*S*,*S*,*S*)-AHPC hydrochloride
≥97%

930717
1*H*-Isoindole-1,3(2*H*)-dione, 2-(2,6-dioxo-3-piperidinyl)-5-(4-piperidinyloxy) hydrochloride
≥95%

920487
1-((1-(*tert*-Butoxycarbonyl)azetidin-3-yl)methyl)-1*H*-pyrazole-3-carboxylic acid

920576
1-((1-(*tert*-Butoxycarbonyl)azetidin-3-yl)methyl)-1*H*-pyrazole-4-carboxylic acid

920649
1-((1-(*tert*-Butoxycarbonyl)piperidin-3-yl)methyl)-1*H*-pyrazole-3-carboxylic acid

923397
1-((2-(*tert*-Butoxycarbonyl)-5-oxa-2-azaspiro[3.4]octan-6-yl)methyl)azetidine-3-carboxylic acid

923389
1-((2-(*tert*-Butoxycarbonyl)-5-oxa-2-azaspiro[3.4]octan-6-yl)methyl)piperidine-4-carboxylic acid

920673
1-(1-(*tert*-Butoxycarbonyl)piperidin-3-yl)-1*H*-pyrazole-4-carboxylic acid

923346
1-(2-(*tert*-Butoxycarbonyl)-2-azaspiro[3.3]heptan-6-yl)-1*H*-pyrazole-4-carboxylic acid

901138
14-Azido-3,6,9,12-tetraoxatetradecan-1-amine

901155
17-Azido-3,6,9,12,15-pentaoxaheptadecan-1-amine

923273
2-((3a*R*,7a*R*)-1-(*tert*-Butoxycarbonyl)octahydro-6*H*-pyrrolo[2,3-c]pyridin-6-yl)thiazole-4-carboxylic acid

920525
2-((4-(*tert*-Butoxycarbonyl)-1,4-diazepan-1-yl)methyl)oxazole-4-carboxylic acid

920665
2-(1-(1-(*tert*-Butoxycarbonyl)piperidin-3-yl)-1*H*-pyrazol-4-yl)acetic acid

920533
2-(2-((1-(*tert*-Butoxycarbonyl)piperidin-4-yl)methyl)-1*H*-imidazol-1-yl)acetic acid

917419
2-(2-(Piperidin-4-yl)phenyl)ethan-1-amine dihydrochloride

920592
2-(2-(*tert*-Butoxycarbonyl)-5-oxo-2,6-diazaspiro[3.6]decan-6-yl)benzoic acid

923419
2-(3-((2-(*tert*-Butoxycarbonyl)-2-azaspiro[3.3]heptan-6-yl)oxy)phenyl)acetic acid

916641
2-(3-(Piperidin-4-yl)phenyl)ethan-1-amine dihydrochloride
95%

923427
2-(4-((2-(*tert*-Butoxycarbonyl)-2-azaspiro[3.3]heptan-6-yl)oxy)piperidin-1-yl)acetic acid

917362
2-(4-(1-(*tert*-butoxycarbonyl)piperidin-4-yl)phenyl)acetic acid
≥95%

923362
2-(4-(5-Bromopyridin-2-yl)-4,7-diazaspiro[2.5]octan-7-yl)acetic acid
≥90%

917605
2-(4-(Piperidin-4-yl)phenyl)acetic acid hydrochloride
≥95%

923354
2-(6-((4-(*tert*-Butoxycarbonyl)piperazin-1-yl)methyl)-5-oxa-2-azaspiro[3.4]octan-2-yl)acetic acid

920495
2-(6-(1-(*tert*-Butoxycarbonyl)azetidin-3-yl)pyrazin-2-yl)acetic acid

916854
2-(Piperidin-4-yl)benzoic acid hydrochloride
95%

912131
2-Chloro-1-(6-methoxy-1,2,3,4-tetrahydroquinolin-1-yl)ethan-1-one
≥95%

927708
2,2,11-trimethyl-4-oxo-3,8-dioxa-5,11-diazatetradecan-14-oic acid
≥95%

930237
2,5-dioxopyrrolidin-1-yl (2,5-dichlorobenzoyl)glycyl-L-leucinate
≥95%

930792
3-[1,3-Dihydro-4-(4-hydroxy-1-butyn-1-yl)-1-oxo-2*H*-isoindol-2-yl]-2,6-piperidinedione
≥95.0%

930822
3-[1,3-Dihydro-4-(5-hydroxy-1-pentyn-1-yl)-1-oxo-2*H*-isoindol-2-yl]-2,6-piperidinedione
≥95.0%

920711
3-(2-(*tert*-Butoxycarbonyl)-5-oxo-2,6-diazaspiro[3.6]decan-6-yl)benzoic acid

923257
3-(3-((7-(*tert*-Butoxycarbonyl)-7-azaspiro[3.5]nonan-2-yl)oxy)piperidin-1-yl)propanoic acid

920657
3-(7-(*tert*-Butoxycarbonyl)-2-ethyl-5,6,7,8-tetrahydroimidazo[1,2-a]pyrazin-3-yl)propanoic acid hydrochloride

920509
3-(7-(*tert*-Butoxycarbonyl)-5,6,7,8-tetrahydroimidazo[1,2-*a*]pyrazin-3-yl)propanoic acid

63179
3-(Maleimido)propionic acid N-hydroxysuccinimide ester
≥98.5% (HPLC)

358657
3-Maleimidopropionic acid *N*-hydroxysuccinimide ester
99%

901699
3-Mercaptopropanyl-*N*-hydroxysuccinimide ester
>95%

923311
4-((2-(*tert*-Butoxycarbonyl)-2-azaspiro[3.3]heptan-6-yl)oxy)picolinic acid

923400
4-((7-(*tert*-Butoxycarbonyl)-7-azaspiro[3.5]nonan-2-yl)oxy)benzoic acid

923265

4-((7-(*tert*-Butoxycarbonyl)-7-azaspiro[3.5]nonan-2-yl)oxy)picolinic acid

920614
4-(2-(*tert*-Butoxycarbonyl)-5-oxo-2,6-diazaspiro[3.6]decan-6-yl)benzoic acid

920622
4-(2-(*tert*-Butoxycarbonyl)-5-oxo-2,6-diazaspiro[3.6]decan-6-yl)picolinic acid

771058
4-(4′-Hydroxyphenylazo)benzoic acid
97%

49667
4-(Maleinimido)phenyl isocyanate
purum, ≥97.0% (CHN)

M6035
4-(*N*-Maleimidomethyl)cyclohexane-1-carboxylic acid 3-sulfo-*N*-hydroxysuccinimide ester sodium salt
powder

M5525
4-(*N*-Maleimidomethyl)cyclohexanecarboxylic acid *N*-hydroxysuccinimide ester
≥98%, powder

930695
4-Aminomethyl-2-(2,6-dioxopiperidin-3-yl)isoindole-1,3-dione hydrochloride
≥95%

63175
4-Maleimidobutyric acid *N*-hydroxysuccinimide ester
≥98.0% (HPLC)

930814
4-Pentynoic acid, 5-[2-(2,6-dioxo-3-piperidinyl)-2,3-dihydro-1-oxo-1*H*-isoindol-4-yl]
≥95.0%

928380
5-{4-[2-(2-{[(*tert*-butoxy)carbonyl]amino}ethoxy)ethyl]piperazin-1-yl}pyrimidine-2-carboxylic acid
≥95%

923338
5-((2-(*tert*-Butoxycarbonyl)-2-azaspiro[3.3]heptan-6-yl)oxy)pyrazine-2-carboxylic acid

920584
5-(1-(*tert*-Butoxycarbonyl)azetidin-3-yl)pyrazine-2-carboxylic acid

920630
5-(2-(*tert*-Butoxycarbonyl)-5-oxo-2,6-diazaspiro[3.6]decan-6-yl)picolinic acid

776173
5-Norbornene-2-acetic acid succinimidyl ester
97%

923370
6-((2-(*tert*-Butoxycarbonyl)-2-azaspiro[3.3]heptan-6-yl)oxy)pyrimidine-4-carboxylic acid

923303
6-((3a*S*,7a*R*)-1-(*tert*-Butoxycarbonyl)octahydro-6*H*-pyrrolo[2,3-c]pyridin-6-yl)pyridazine-3-carboxylic acid

920517
6-(4-(*tert*-Butoxycarbonyl)-1,4-diazepan-1-yl)pyrazine-2-carboxylic acid

M9794
6-Maleimidohexanoic acid *N*-hydroxysuccinimide ester
≥98%, powder

63177
6-Maleimidohexanoic acid *N*-hydroxysuccinimide ester
≥98.0% (HPLC)

451088
6-Mercapto-1-hexanol
97%

923958
6F,C5-Pomalidomide-4-piperidine-C_1-piperazine hydrochloride

917710
A1V1PF2-OEt
≥95%

917672
A1V1PF2-OEt-C10-NH_2 hydrochloride

917427
A1V1PF2-OEt-C_6-NH_2 hydrochloride

917923
A1V1PF2-OEt-PEG_1-NH_2 hydrochloride

916676
A1V1PF2-OEt-PEG_3-NH_2 hydrochloride

917222
A1V2PF1-NHEt
≥95%

917206
A1V2PF1-NHEt-C_{10}-NH_2

916943
A1V2PF1-NHEt-C_6-NH_2

917451
A1V2PF1-NHEt-PEG_1-NH_2
≥95%

917702
A1V2PF1-NHEt-PEG_3-NH_2

916714
A1V2PF2-NHEt
≥95%

916684
A1V2PF2-NHEt-C_{10}-NH_2
≥95%

917931
A1V2PF2-NHEt-C_6-NH_2

916935
A1V2PF2-NHEt-PEG$_1$-NH$_2$

917192
A1V2PF2-NHEt-PEG$_3$-NH$_2$
≥95%

797685
ADIBO-PEG4-acid
90% (HPLC)

SML2923
ARV-771
≥98% (HPLC)

914223
Biotin-SLF

913979
Biotin-Thalidomide
≥95%

917478
BocA1V1PF2
≥95%

917184
BocA1V1PF2-OC$_{10}$-NH$_2$ hydrochloride

916927
BocA1V1PF2-OC$_6$-NH$_2$ hydrochloride
≥95%

917435
BocA1V1PF2-OPEG$_1$-NH$_2$ hydrochloride
≥95%

917680
BocA1V1PF2-OPEG$_3$-NH$_2$ hydrochloride

916978
BocA1V2PF1
≥95%

916706
BocA1V2PF1-NHC$_{10}$-NH$_2$

917966
BocA1V2PF1-NHC$_6$-NH$_2$

916951
BocA1V2PF1-NHPEG$_1$-NH$_2$

917214
BocA1V2PF1-NHPEG$_3$-NH$_2$
≥85%

917974
BocA1V2PF2

917699
BocA1V2PF2-NHC$_{10}$-NH$_2$
≥85%

917443
BocA1V2PF2-NHC$_6$-NH$_2$

917958
BocA1V2PF2-NHPEG$_1$-NH$_2$

916692
BocA1V2PF2-NHPEG$_3$-NH$_2$
≥95%

902675
BocNH-PEG$_6$-acid

901572
Bromo-PEG$_2$-acid
≥95%

901568
Bromo-PEG$_3$-acid

911690
C5 Lenalidomide
≥95%

920754
C5 Lenalidomide-benzyl-piperazine hydrochloride

917303
C5 Lenalidomide-C$_6$-PEG$_1$-C$_3$-PEG$_1$-Butyl NH$_2$ hydrochloride
≥95%

911720
C5 Lenalidomide-C$_9$-NH$_2$ hydrochloride
≥95%

927732
C5 Lenalidomide-difluoroPEG$_1$-C$_4$-piperazine Hydrochloride
≥95%

919896
C5 Lenalidomide-dipiperazine-NH$_2$ hydrochloride

920800
C5 Lenalidomide-methylamino-PEG1-NH$_2$ hydrochloride
≥95%

911739
C5 Lenalidomide-PEG$_1$-NH$_2$ hydrochloride
≥95%

919969
C5 Lenalidomide-PEG$_2$-Butyl NH$_2$ hydrochloride
≥95%

911755
C5 Lenalidomide-PEG$_5$-NH$_2$ hydrochloride

921327
C5 Lenalidomide-PEG$_6$-Butyl NH$_2$ hydrochloride
≥95%

917729
C5 Lenalidomide-piperazine-pyridine-alkyne-NH_2 hydrochloride
≥95%

920797
C5 Lenalidomide-pyridine-PEG_1-piperazine hydrochloride
≥95%

921300
C5 Lenalidomide-pyrimidine-piperazine-PEG_1-NH_2 hydrochloride
≥95%

911704
C5 Lenalidomine-C_3-NH_2 hydrochloride
≥95%

911712
C5 Lenalidomine-C_6-NH_2 hydrochloride
≥95%

929441
C5-Pomalidomide-piperazine hydrochloride
≥95%

919381
CCW16
≥95%

919403
CCW16-C4-BocNH
95%

919411
CCW16-C6-BocNH

919489
CCW16-C6-PEG_3-butyl-BocNH

919438
CCW16-C9-BocNH
≥95%

919446

CCW16-PEG$_1$-BocNH

919470
CCW16-PEG$_2$-butyl-BocNH
≥95%

919454
CCW16-PEG$_3$-BocNH
≥95%

919462
CCW16-PEG$_5$-BocNH
≥95%

SML2687
dBET1
≥98% (HPLC)

SML2683
dBET6
≥98% (HPLC)

SML2911
dBRD9 Hydrochloride
≥97% (HPLC)

761516
Dibenzocyclooctyne-acid
95%, storage temp.:-20°C

761540
Dibenzocyclooctyne-amine
for Copper-free Click Chemistry

761524
Dibenzocyclooctyne-*N*-hydroxysuccinimidyl ester
for Copper-free Click Chemistry

759902
Dibenzocyclooctyne-PEG4-acid
for Copper-free Click Chemistry

761982
Dibenzocyclooctyne-PEG4-alcohol
for Copper-free Click Chemistry

764019
Dibenzocyclooctyne-PEG4-*N*-hydroxysuccinimidyl ester
≥90%

762040
Dibenzocyclooctyne-sulfo-*N*-hydroxysuccinimidyl ester
for Copper-free Click Chemistry

SML2601
dTAG-13
≥98% (HPLC)

SML2684
dTRIM24
≥98% (HPLC)

SML2950
E7820
≥98% (HPLC)

925101
E7820-C_3-NH_2 hydrochloride

925128
E7820-C6-NH_2 hydrochloride

924997
E7820-C_9-NH_2 hydrochloride

925004
E7820-PEG_1-NH_2 hydrochloride

925039
E7820-PEG_3-NH_2 hydrochloride

925012
E7820-PEG_5-NH_2 hydrochloride

929484
FBnG-C3-PEG_1-C3-NH_2 hydrochloride
≥95%

929301
FBnG-C3-PEG$_3$-C3-NH$_2$ hydrochloride
≥95%

929328
FBnG-C3-PEG$_5$-C3-NH$_2$ hydrochloride
≥95%

914738
KB02-SLF
≥95%

914975
KB03-SLF
≥95%

913715
KB05-SLF

901558
Lenalidomide
≥95%

SML2283
Lenalidomide
≥98% (HPLC)

918091
Lenalidomide-Photoswitch2-NH$_2$ hydrochloride
≥95%

920681
Lithium 2-(3-(4-(*tert*-butoxycarbonyl)-1,4-diazepan-1-yl)pyrazin-2-yl)acetate

M3884
Maleimidoacetic acid N-hydroxysuccinimide ester
≥95%

930210
***N*-(2,5-Dichlorobenzoyl)glycyl-L-leucine**
≥95.0%

929476
***N*-Acetyl-*S*-(2-amino-9-(4-fluorobenzyl)-6-oxo-6,9-dihydro-1*H*-purin-8-yl)-L-cysteine**
≥95%

745073
***N*-[(1*R*,8*S*,9*s*)-Bicyclo[6.1.0]non-4-yn-9-ylmethyloxycarbonyl]-1,8-diamino-3,6-dioxaoctane**
for Copper-free Click Chemistry

56951
***N*-(2-Aminoethyl)maleimide trifluoroacetate salt**
≥95% (HPLC), ≥98% (T)

773263
N-(2-Hydroxyethyl)maleimide
97%

100919
***N*-(4-Bromobutyl)phthalimide**
98%

911798
***N*-(4-Bromophenyl)-*N*-phenylacrylamide**
≥95%

901494
***N*-Methylated pomalidomide**
≥98%

N6287
Nutlin-3
≥98% (HPLC), powder

SML0580
Nutlin-3a
≥98% (HPLC)

929433
Opto-pomalidomide-C_2-NH_2 hydrochloride
≥95%

929379
Opto-thalidomide-*O*-acetamide-C_4-NH_2 hydrochloride

P0018
Pomalidomide
≥98% (HPLC)

930660
Pomalidomide 4′-PEG$_3$-amine hydrochloride
≥95%

923931
Pomalidomide-4-piperidine-C$_1$-piperazine hydrochloride

920746
Pomalidomide-benzyl-piperazine hydrochloride

930652
Pomalidomide-C$_2$-NH$_2$ hydrochloride
≥95%

929336
Pomalidomide-C$_5$-phosphoramidite

901500
Pomalidomide-C$_3$-CO$_2$H
≥95%

911658
Pomalidomide-C$_3$-NH$_2$ hydrochloride
≥95%

901496
Pomalidomide-C$_6$-CO$_2$H
≥98%

911666
Pomalidomide-C$_6$-NH$_2$ hydrochloride
≥95%

904880
Pomalidomide-C$_6$-PEG$_1$-C$_3$-PEG$_1$-butyl azide
≥95%

906050
Pomalidomide-C$_6$-PEG$_1$-C$_3$-PEG$_1$-butyl iodide
≥95%

904686
Pomalidomide-C$_6$-PEG$_3$-butyl azide
≥95%

904708
Pomalidomide-C_6-PEG_3-butyl iodide
≥95%

901525
Pomalidomide-C_9-CO_2H
≥95%

920738
Pomalidomide-C9-NH_2 hydrochloride

927678
Pomalidomide-difluoroPEG_1-C_4-piperazine Hydrochloride
≥95%

915572
Pomalidomide-dipiperazine-NH_2 hydrochloride
≥95%

921319
Pomalidomide-methylamino-PEG1-NH_2 hydrochloride

930547
Pomalidomide-PEG_1-C_2-azide
≥95.0%

930563
Pomalidomide-PEG_2-C_2-azide
≥95%

930520
Pomalidomide-PEG_2-C_2-NH_2 hydrochloride
≥95%

901523
Pomalidomide-PEG_1-Alkyne
≥98%

903825
Pomalidomide-PEG1-azide

901527

Pomalidomide-PEG$_1$-CO$_2$H
≥95%

☐
901516
Pomalidomide-PEG$_1$-NH$_2$ hydrochloride
≥95%

☐
901529
Pomalidomide-PEG$_2$-Alkyne
≥95%

☐
903833
Pomalidomide-PEG2-azide

☐
910449
Pomalidomide-PEG$_2$-butyl CO$_2$H
≥95%

☐
904724
Pomalidomide-PEG$_2$-butyl iodide
≥95%

☐
901526
Pomalidomide-PEG$_2$-CO$_2$H
95%

☐
901513
Pomalidomide-PEG$_2$-NH$_2$ hydrochloride
≥95%

☐
901531
Pomalidomide-PEG$_3$-Alkyne
≥95%

☐
904678
Pomalidomide-PEG3-azide

☐
901504
Pomalidomide-PEG$_3$-CO$_2$H
≥95%

☐
901495
Pomalidomide-PEG$_3$-NH$_2$ hydrochloride
≥95%

909386
Pomalidomide-PEG$_4$-Azide
≥95%

901824
Pomalidomide-PEG$_4$-CO$_2$H

901830
Pomalidomide-PEG$_4$-NH$_2$ hydrochloride
≥95%

901834
Pomalidomide-PEG$_5$-Alkyne
≥95%

909394
Pomalidomide-PEG$_5$-azide
≥95%

901828
Pomalidomide-PEG$_5$-CO$_2$H
≥95%

901831
Pomalidomide-PEG$_5$-NH$_2$ hydrochloride

909408
Pomalidomide-PEG$_6$-azide
≥95%

905224
Pomalidomide-PEG$_6$-butyl amine hydrochloride
≥95%

903434
Pomalidomide-PEG$_6$-butyl azide
≥95%

910481
Pomalidomide-PEG$_6$-butyl CO$_2$H
≥95%

903442
Pomalidomide-PEG_6-butyl iodide
≥95%

901829
Pomalidomide-PEG_6-CO_2H

901832
Pomalidomide-PEG_6-NH_2 hydrochloride
≥98%

930555
Pomalidomide-piperazine-acetic acid
≥95%

923877
Pomalidomide-piperazine-C_1-4-piperidine hydrochloride

923885
Pomalidomide-piperazine-piperidine-4-carboxamide hydrochloride

930709
Pomalidomide-piperazine-propanoic acid
≥95.0%

930571
Pomalidomide-piperidine-carboxylic acid
≥95%

920789
Pomalidomide-pyridine-PEG_1-piperazine hydrochloride

901567
Propargyl-PEG_1-acid
≥95%

901566
Propargyl-PEG_2-acid

901565
Propargyl-PEG_3-acid

42002
Sulfo-N-succinimidyl 4-maleimidobutyrate sodium salt
~90%

901577
t-Boc-*N*-amido-PEG_2-CH_2CO_2H
≥95%

901575
t-Boc-*N*-amido-PEG_3-CH_2CO_2H
≥95%

790435
TCO PEG4 succinimidyl ester

790451
TCO-amine HCl salt

790443
TCO-PEG3-maleimide

920541
***tert*-Butyl 1-oxa-4,9-diazaspiro[5.6]dodecane-9-carboxylate**

920568
***tert*-Butyl 3-(5,6,7,8-tetrahydroimidazo[1,2-a]pyrazin-3-yl)propanoate**

917400
***tert*-Butyl 4-(3-(hydroxymethyl)phenyl)piperidine-1-carboxylate**
≥95%

900913
Tetrazine-PEG5-NHS ester
≥95%

SML2686
THAL-SNS-032
≥98% (HPLC)

919888
Thalidomide-Photoswitch3-NH_2 hydrochloride
≥95%

930628
VH 032 amide-alkyl C_3-acid
≥95%

930636
VH 032 amide-alkyl C_5-acid
≥95%

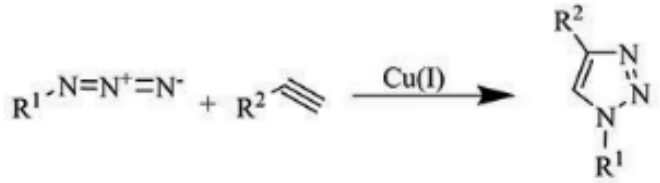
930644
VH 032 amide-PEG_2-acid
≥95%

930601
VH032-cyclopropane-F
≥95%

930598
VH032-OH
≥95%

SML1896
VH298
≥98% (HPLC)

Click Chemistry Reagents



Our extensive portfolio of click chemistry reagents offers a variety of azides, alkynes, catalysts, and ligands to help accelerate your research in the exciting arena of "click" chemistry. Click chemistry is a term coined by Barry Sharpless to describe chemical reactions that are modular, efficient, wide in scope, provide very high yields, and generate only inoffensive byproducts. The most well-known example of a "click" reaction is the Copper(I)-catalyzed Azide-Alkyne 1,3-dipolar Cycloaddition (CuAAC), which yields a 1,4-disubstituted five-membered 1,2,3-triazole ring.

This reaction between azides and alkynes offers high yields and involves functionalities that can be introduced relatively easily in a variety of molecules such as synthetic polymers, fluorophores, small molecules or into specific locations in biomolecules. An advantage of this reaction for biological purposes is that the azide and alkyne functional groups are largely

inert, or biorthogonal, towards biological molecules and aqueous environments. “Click” chemistry continues to gain popularity and is used in a variety of research fields with significant contributions to the fields of chemical biology, polymer chemistry, bioconjugation, and drug discovery.

AMINO ACID AZIDES/ALKYNES FOR CLICK CHEMISTRY

Peptide synthesis using natural and non-natural amino acids is a powerful tool in the development of therapeutics and in understanding biological chemistry. We offer a variety of Fmoc- and Boc-protected azido amino acids for your peptide or amino acid-based chemical ligation needs, such as azide-alkyne cycloaddition reactions and Staudinger ligations.

AZIDE SOURCES FOR CLICK CHEMISTRY

Incorporating azido functional groups into organic molecules is becoming an increasingly important task as these moieties continue to impact organic chemistry as well as biology in uses ranging from amino group protection to chemical ligation. We offer a wide selection of azide sources, from sodium azide to diphenyl phosphoryl azide, to facilitate azide synthesis and the preparation of tailor-made organic azides.

ORGANIC AZIDES FOR CLICK CHEMISTRY

Since the preparation of the first organic azide, phenyl azide, by Peter Griess in 1864, this energy-rich and versatile class of compounds has enjoyed considerable interest. Completely new perspectives have emerged, notably the use of organic azides for peptide synthesis, combinatorial synthesis, heterocycle synthesis, and the ligation or modification of biopolymers. The most prominent fields of application today are azide-alkyne cycloadditions and different variants of the Staudinger ligation. The azido group can also be used as a protecting group for primary amines, especially in sensitive substrates such as complex carbohydrates or peptide nucleic acids (PNA) and coordination compounds, as azides are stable to alkene metathesis conditions.

PEG AZIDES FOR CLICK CHEMISTRY

PEG polymers contain numerous inherently favorable biological characteristics including high water solubility and a lack of toxicity and immunogenicity. Thus, the chemical modification of biologically active compounds, such as peptides, antibody fragments, enzymes, or small molecules with polyethylene glycol chains, referred to as “PEGylation”, often leads to improved pharmacokinetics and biological function in many applications. Our PEG azides are ideal starting materials for the synthesis of PEG derivatives via azide-alkyne cycloaddition or Staudinger ligation.

TRIFUNCTIONAL PROBE BUILDING BLOCKS FOR CLICK CHEMISTRY

Small-molecule probes are widely used in chemical biology research for target ID/validation and the interrogation of biological systems. We have compiled a collection of trifunctional building blocks to facilitate the design and synthetic development of chemical probes. Each contains three components: a connectivity group, a reactive group, and a bio-orthogonal handle for downstream applications. Not only does the collection enable simultaneous incorporation of reactive groups, but the connectivity group (e.g. amine) can be leveraged

to prepare libraries of probe analogs, allowing the biologist to screen for the optimal probe for a given assay.

TETRAZINE/STRAINED-ALKENES FOR CLICK CHEMISTRY

The reaction of 1,2,4,5 tetrazines with strained alkenes has been applied as a rapid bio-orthogonal click chemistry reaction for biological labeling and cell detection applications among others. This reaction precedes rapidly via an inverse electron demand [4 + 2] Diels-Alder cycloaddition to yield a stable covalent linkage without the need for a catalyst and the only byproduct is dinitrogen. When using trans-cyclooctene, this reaction is orders of magnitude faster than azide-cyclooctyne-based click chemistry and has therefore found use in applications where low concentrations of reactants are desired or faster kinetics are needed.

COPPER-FREE CLICK CHEMISTRY

Cu-free cycloadditions offer efficient ligation reactions useful for a variety of bioconjugation applications. When working with live cells these reactions offer the advantage of not having the cytotoxicity associated with Cu-catalyzed cycloadditions. We offer a suite of cyclooctynes, tetrazines and strained-alkene reagents for use in a variety of applications.

STAUDINGER LIGATION

The reaction between an azide and a phosphine forming an aza-ylide was first reported in 1919 by Nobel Prize laureate Herrmann Staudinger. It has found widespread application in chemical synthesis and is valuable as a highly chemoselective ligation method for the preparation of bioconjugates. Both reactive functionalities involved in the Staudinger ligation reaction are bio-orthogonal and readily combine at room temperature in aqueous environments. These conditions make it possible to exploit the Staudinger ligation in complex cellular and organismal environments in the investigation of various processes in chemical biology. Our portfolio offers phosphine ligands for your various Staudinger ligation and conjugation applications.

744867
(1*R*,8*S*,9*s*)-Bicyclo[6.1.0]non-4-yn-9-ylmethyl *N*-succinimidyl carbonate
for Copper-free Click Chemistry

709492
(1*S*,4*S*,8*S*)-5-Benzyl-2-isobutyl-8-methoxy-1,8-dimethylbicyclo[2.2.2]octa-2,5-diene
97%

T511293
(2′*S*)-2′-Deoxy-2′-fluoro-5-ethynyluridine, (F-ara-EdU)
Aldrich^{CPR}

761591
(4-(1,2,4,5-Tetrazin-3-yl)phenyl)methanamine hydrochloride
95%

900750
(4-(2-(2-(2-Aminoethoxy)ethoxy)ethoxy)phenyl)(4-(prop-2-yn-1-yloxy)phenyl)methanone HCl salt
≥95%

900603
(4-(2-(2-Aminoethoxy)ethoxy)phenyl)(4-(prop-2-yn-1-yloxy)phenyl)methanone
≥95%

900599
(4-(Aminomethyl)phenyl)(4-(prop-2-yn-1-yloxy)phenyl)methanone
95%

900598
(4-(Bromomethyl)phenyl)(4-(prop-2-yn-1-yloxy)phenyl)methanone
≥95%

900615
(4-Hydroxyphenyl)(4-(prop-2-yn-1-yloxy)phenyl)methanone
≥95%

764523
(*E*)-Cyclooct-4-enyl 2,5-dioxo-1-pyrrolidinyl carbonate

901503
(*S*,*R*,*S*)-AHPC-PEG$_1$-Alkyne
≥95%

903957
(*S*,*R*,*S*)-AHPC-PEG1-azide

901517
(*S*,*R*,*S*)-AHPC-PEG$_2$-Alkyne
≥95%

904813
(*S*,*R*,*S*)-AHPC-PEG2-azide

901533
(*S*,*R*,*S*)-AHPC-PEG$_3$-Alkyne
≥95%

904651
(*S*,*R*,*S*)-AHPC-PEG$_3$-azide
≥95%

901851
(*S*,*R*,*S*)-AHPC-PEG$_4$-Alkyne

901855
(*S*,*R*,*S*)-AHPC-PEG$_5$-Alkyne

901873
(*S*,*R*,*S*)-AHPC-PEG$_6$-Alkyne

909378
(*S*,*R*,*S*)-AHPC-PEG$_6$-Azide
≥95%

904864
(*S*,*R*,*S*)-AHPC-PEG$_6$-butyl azide
≥95%

714291
(*S*)-5-Azido-2-(Fmoc-amino)pentanoic acid
≥97.0% (HPLC)

513989
1-Azido-1-deoxy-β-D-galactopyranoside
97%

514004
1-Azido-1-deoxy-β-D-glucopyranoside

514012
1-Azido-1-deoxy-β-D-lactopyranoside
97%

901152
1-Azido-2-(2-(2-ethoxyethoxy)ethoxy)ethane
≥95%

901132
1-Azido-2-(2-(2-methoxyethoxy)ethoxy)ethane

765953
1,11-Diazido-3,6,9-trioxaundecane

901136

1,17-Diazido-3,6,9,12,15-pentaoxaheptadecane

17758
11-Azido-3,6,9-trioxaundecan-1-amine
technical, ≥90% (GC)

901155
17-Azido-3,6,9,12,15-pentaoxaheptadecan-1-amine

ALD00024
2-(2-Propyn-1-yloxy)ethanesulfonyl fluoride
96%

679011
2-(Diphenylphosphino)terephthalic acid 1-methyl 4-pentafluorophenyl diester
97%

900858
2-(Prop-2-yn-1-yloxy)-4-(3-(trifluoromethyl)-3*H*-diazirin-3-yl)benzoic acid
≥95%

670995
2-Acetamido-2-deoxy-β-D-glucopyranosyl azide
≥98% (HPLC)

712795
2-Azido-2-deoxy-D-glucose
≥97.0% (HPLC)

ALD00160
2-Thiophenesulfonyl fluoride
95%

ALD00148
2,3-Dihydro-1,4-benzodioxin-6-sulfonyl fluoride
95%

919950
3-(([1,1′-Biphenyl]-4-ylmethyl)thio)-6-methyl-1,2,4,5-tetrazine
>98%

900614
3-(4-(Prop-2-yn-1-yloxy)benzoyl)benzoic acid
≥95%

690104
3-(4-Azidophenyl)propionic acid
≥97.0% (T)

900605
3-(Bromomethyl)-5-((trimethylsilyl)ethynyl)benzenesulfonyl fluoride
≥95%

ALD00168
3-(Fluorosulfonyl)-2-thiophenecarboxylic acid methyl ester
95%

762016
3-Azido-1-propanamine
≥95%

776130
3-Azido-1-propanol
≥96%

900859
3-Bromo-5-((trimethylsilyl)ethynyl)benzenesulfonyl fluoride
≥95%

498289
3-Ethynylaniline
≥98%

ALD00094
3-Methoxybenzenesulfonyl fluoride

ALD00364
4-(2-Bromoacetyl)benzenesulfonyl fluoride
95% (GC)

900602
4-(4-(Prop-2-yn-1-yloxy)benzoyl)benzoic acid
≥95%

ALD00176
4-(Acetylamino)benzenesulfonyl fluoride
95%

404764
4-Acetamidobenzenesulfonyl azide

97%

359556
4-Azidoaniline hydrochloride
97%

340138
4-Carboxybenzenesulfonazide
97%

481122
4-Ethynylaniline
97%

ALD00130
4-Methoxybenzenesulfonyl fluoride
95%

152854
4-Methoxybenzyloxycarbonyl azide
95%

ALD00132
4-Phenoxybenzenesulfonyl fluoride
95%

900613
5-(4-(4-(Prop-2-yn-1-yloxy)benzoyl)phenoxy)pentanoic acid
≥95%

ALD00158
5-Chloro-2-thiophenesulfonyl fluoride
95%

809705
5-DBCO-PEG$_4$-dUTP, 10mM Aqueous Solution

909475
5-Ethynyl uridine
≥95%

T511307
5-Ethynyl-2′-deoxycytidine, (EdC)
Aldrich^{CPR}

776173
5-Norbornene-2-acetic acid succinimidyl ester
97%

712752
6-Azido-6-deoxy-D-galactose
≥98.0% (HPLC)

712760
6-Azido-6-deoxy-D-glucose
≥95% (HPLC)

8.51097
6-Azido-hexanoic acid
Novabiochem®

797685
ADIBO-PEG4-acid
90% (HPLC)

901982
Alkyne Agarose

764205
Alkyne-PEG4-maleimide
90% (HPLC)

764167
Alkyne-PEG5-acid

764191
Alkyne-PEG5-*N*-hydroxysuccinimidyl ester

JKA4095
Alkyne-PEG5K-Alkyne
average M_n 5,000

764248
Amino-PEG4-alkyne

900957
Azide agarose

762024
Azide-PEG3-biotin conjugate

QBD10524
Azido-dPEG®$_{11}$-amine

QBD10513
Azido-dPEG®$_{12}$-acid

QBD10505
Azido-dPEG®$_{12}$-NHS ester

QBD10569
Azido-dPEG ®$_{12}$-TFP ester
>90%

QBD10525
Azido-dPEG®$_{23}$-amine

QBD10514
Azido-dPEG®$_{24}$-acid

QBD10571
Azido-dPEG ®$_{24}$-TFP ester

QBD10526
Azido-dPEG ®$_{35}$-amine
>90%

QBD10573
Azido-dPEG ®$_{36}$-TFP ester

QBD10502
Azido-dPEG®$_{4}$-acid

QBD10501
Azido-dPEG®$_{4}$-NHS ester

QBD10567
Azido-dPEG ®$_{4}$-TFP ester

QBD10523
Azido-dPEG®$_7$-amine

QBD10512
Azido-dPEG®$_8$-acid

QBD10503
Azido-dPEG®$_8$-NHS ester

244546
Azidomethyl phenyl sulfide
95%

349488
Azidotrimethyltin(IV)
97%

900891
Azo biotin-azide

900912
Biotin picolyl azide

QBD11811
Biotin-dPEG ®$_{12}$-DBCO

764213
Biotin-PEG4-alkyne
for copper catalyzed click labeling

QBD11372
Bis-dPEG ®$_{11}$-DBCO

712221
Boc-propargyl-Gly-OH
≥98.0% (HPLC)

QBD11204

Bromoacetamido-dPEG ®$_{11}$-azide

QBD11223
Bromoacetamido-dPEG ®$_{12}$-amido-DBCO

QBD11205
Bromoacetamido-dPEG ®$_{23}$-azide

QBD11224
Bromoacetamido-dPEG ®$_{24}$-amido-DBCO

QBD11217
Bromoacetamido-dPEG ®$_{3}$-azide

QBD11221
Bromoacetamido-dPEG ®$_{4}$-amido-DBCO

909505
C8-Alkyne-dU-CEP

212865
Copper(I) bromide
98%

205540
Copper(I) iodide
98%

326755
Copper(II) acetate
98%

C1297
Copper(II) sulfate
ReagentPlus®, ≥99%

209198
Copper(II) sulfate pentahydrate
ACS reagent, ≥98.0%

777331
Cy3-alkyne
for copper catalyzed click labeling

777315
Cy3-azide
90% (HPLC)

777366
DBCO-Cy3
for Copper-free Click Chemistry

777374
DBCO-Cy5
for Copper-free Click Chemistry

QBD11812
DBCO-dPEG ®$_{12}$-carboxyfluorescein

QBD10593
DBCO-dPEG ®$_{12}$-MAL

QBD11366
DBCO-dPEG ®$_{12}$-TFP ester

QBD11383
DBCO-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

QBD10594
DBCO-dPEG ®$_{24}$-MAL

QBD11370
DBCO-dPEG ®$_{24}$-TFP ester

QBD11362
DBCO-dPEG ®$_{4}$-TFP ester

QBD11815
DBCO-TFP ester

901991
Dde biotin picolyl azide

901765
Dde biotin-azide

761516
Dibenzocyclooctyne-acid
95%, storage temp.:-20°C

761540
Dibenzocyclooctyne-amine
for Copper-free Click Chemistry

760668
Dibenzocyclooctyne-maleimide
for Copper-free Click Chemistry

761524
Dibenzocyclooctyne-*N*-hydroxysuccinimidyl ester
for Copper-free Click Chemistry

759902
Dibenzocyclooctyne-PEG4-acid
for Copper-free Click Chemistry

761982
Dibenzocyclooctyne-PEG4-alcohol
for Copper-free Click Chemistry

760749
Dibenzocyclooctyne-PEG4-biotin conjugate
for Copper-free Click Chemistry

760676
Dibenzocyclooctyne-PEG4-maleimide
for Copper-free Click Chemistry

764019
Dibenzocyclooctyne-PEG4-*N*-hydroxysuccinimidyl ester
≥90%

762040
Dibenzocyclooctyne-sulfo-*N*-hydroxysuccinimidyl ester
for Copper-free Click Chemistry

178756
Diphenyl phosphoryl azide
97%

910147
Fluorescein Azide
≥95%

F5312
Fmoc-(*R*)-propargyl-Ala-OH

F5437
Fmoc-(*S*)-propargyl-Ala-OH

714151
Fmoc-β-azido-Ala-OH
≥98.0% (HPLC)

808520
γ-[(6-Azidohexyl)-imido]-ATP sodium salt

808512
γ-[(Propargyl)-imido]-ATP sodium salt

900610
Hydroxyl di-benzophenone $(PEG)_3$ alkyne

925225
KB02yne

925144
KB05yne
≥95%

QBD10596
m-dPEG ®$_{12}$-DBCO

QBD10540
m-dPEG ®$_{24}$-Azide (Azido-m-dPEG ®$_{24}$)

689807
Methoxypolyethylene glycol azide
PEG average M_n 2,000

689475
Methoxypolyethylene glycol azide

900914
Methyltetrazine-NHS ester
≥95%

JKA13019
mPEG12-Azide

JKA3177
mPEG5K-Alkyne
average M_n 5,000

8.51021
N_3-$(PEG)_7$-COOH (33 atoms)
Novabiochem®

745073
***N*-[(1*R*,8*S*,9*s*)-Bicyclo[6.1.0]non-4-yn-9-ylmethyloxycarbonyl]-1,8-diamino-3,6-dioxaoctane**
for Copper-free Click Chemistry

708925
***N*-[[2-(Diphenylphosphino)phenyl]methylene]-3-methyl-L-valine sodium salt**
95%

923818
***N*-(But-3-yn-1-yl)-2-(((2-oxo-2*H*-chromen-7-yl)oxy)methyl)acrylamide**

900915
***N*-Azidoacetylgalactosamine-tetraacylated**

900917
***N*-Azidoacetylmannosamine-tetraacylated**
95%

778923
***N*-Boc-4-pentyne-1-amine**
≥95% (GC)

JKA5086
N3-PEG3500-NHS
average M_n 3,500

JKA5239
N3-PEG5K-NH2
TFA Salt, average M_n 5,000

JKA5088
N3-PEG5K-NHS
average M_n 5,000

808490
N^6-Propargyl-ATP sodium salt

76318
***O*-(2-Aminoethyl)-*O'*-(2-azidoethyl)heptaethylene glycol**
≥90% (oligomer purity)

77787
***O*-(2-Aminoethyl)-*O'*-(2-azidoethyl)nonaethylene glycol**
≥90% (oligomer purity)

76172
***O*-(2-Aminoethyl)-*O'*-(2-azidoethyl)pentaethylene glycol**
≥90% (oligomer purity)

71613
O-(2-Azidoethyl)-O-[2-(diglycolyl-amino)ethyl]heptaethylene glycol
≥90% (oligomer purity)

689440
O-(2-Azidoethyl)heptaethylene glycol
≥95% (oligomer purity)

901767
PC biotin-PEG3-NHS ester

807737
Photo-reactive Clickable *cis*-Sterol Probe

756598
Poly(ethylene glycol) bisazide
average M_n 1,100

445878
Poly(ethylene glycol) dimethyl ether
average M_n ~250

445908
Poly(ethylene glycol) dimethyl ether

average M_n ~2,000

445894
Poly(ethylene glycol) dimethyl ether
average M_n ~1,000

445886
Poly(ethylene glycol) dimethyl ether
average M_n ~500, contains 100 ppm BHT as stabilizer

305413
Poly(ethylene glycol) distearate
average M_n ~930

689696
Polyoxyethylene bis(azide)
average M_n 2,000

904880
Pomalidomide-C_6-PEG_1-C_3-PEG_1-butyl azide
≥95%

904686
Pomalidomide-C_6-PEG_3-butyl azide
≥95%

901523
Pomalidomide-PEG_1-Alkyne
≥98%

903825
Pomalidomide-PEG1-azide

901529
Pomalidomide-PEG_2-Alkyne
≥95%

903833
Pomalidomide-PEG2-azide

901531
Pomalidomide-PEG_3-Alkyne
≥95%

904678
Pomalidomide-PEG3-azide

909386
Pomalidomide-PEG$_4$-Azide
≥95%

901834
Pomalidomide-PEG$_5$-Alkyne
≥95%

909394
Pomalidomide-PEG$_5$-azide
≥95%

909408
Pomalidomide-PEG$_6$-azide
≥95%

903434
Pomalidomide-PEG$_6$-butyl azide
≥95%

764221
Propargyl-*N*-hydroxysuccinimidyl ester

901566
Propargyl-PEG$_2$-acid

902667
Propargyl-PEG6-acid

P50919
Propargylamine hydrochloride
95%

760706
Sulfo-dibenzocyclooctyne-biotin conjugate

for Copper-free Click Chemistry

790435
TCO PEG4 succinimidyl ester

790451
TCO-amine HCl salt

790443
TCO-PEG3-maleimide

900927
TCO-PNB ester
≥95%

QBD11371
Tetra(-dPEG ®$_{11}$-DBCO)PE

651664
Tetrabutylammonium azide

172405
Tetraethylene glycol dimethyl ether
≥99%

473839
Tetrafluoro-2-(tetrafluoro-2-iodoethoxy)ethanesulfonyl fluoride
97%

900913
Tetrazine-PEG5-NHS ester
≥95%

678937
Tris[(1-benzyl-1*H*-1,2,3-triazol-4-yl)methyl]amine
97%

762342
Tris(3-hydroxypropyltriazolylmethyl)amine
95%

900331
VPhos

95%

Chemical Biology Tools

We make ingenuity accessible for you so you can focus on bridging the gap between chemistry and biology. Breakthrough in your work and lead the scientific community into an emerging field backed by our best-in-class solutions and industry leadership.

DEXTERITY KITS

Dexterity Kits are an easy, all-in-one method to perform high sensitivity proximity labeling for any protein as long as there is a known antibody or small molecule specific for their protein of interest (POI). These protein labeling kits allow you to perform precise micromapping within 4 nm of your POI. Dexterity Kits provide:

- Facile incorporation of small molecules as an amide linkage, either with an amine or acid group
- Short labeling radius
- Benign method of activation (visible light)
- Catalytic signal amplification

DNA-ENCODED LIBRARIES (DELS)

DNA Encoded Fragment Library

The fragment DELs from our partner DyNAbind provides an entire screening library in a single tube, making your drug discovery process faster, more effective, and less expensive. We've lowered the barriers to DyNAbind DEL technology to provide:

- Off-the-shelf, affordable access to DEL technology
- Dynamic fragment library for a revolutionary approach to screening with DELs
- High-level quality control of every fragment component maximizes data reliability and validation of good hits
- Rule-of-Three used in library design ensures compounds are in favorable fragment space with good ligand efficiency and plenty of room for optimization
- Ability to perform and validate initial DEL screening internally before engaging service provider
- A library with maximized diversity, ready to deploy against nearly any druggable target

10 Million Compound DNA-Encoded Library (DEL)

This DNA-encoded library from our partner DyNAbind contains a library of 10 million small molecule compounds each tagged with a DNA barcode. This DEL provides:

- 10 Million compound, off-the-shelf, affordable DEL
- Fits in a single microcentrifuge tube
- No sophisticated robotics required to use the kit
- Rapid assay by binding, washing, PCR, and next gen sequencing
- Free analysis on our website portal
- Structures are revealed for 50 hits or up to 60 with ties at no additional charge
- The label license allows any use of hits in research and development or manufacturing at no additional charge

CHEMICAL PROBES AND TOOLS

Probe Building Blocks: Design Your Own Probe

With our selection of reactive groups, scaffolds, linkers and tags, you can design probes to include functionalities compatible for:

- Target engagement
- Enrichment
- Fluorescence
- Bioorthogonal chemistries

Target-Based Probes: Selective and Characterized

Choose from probes with high-specificity for one target or subset of a protein to provide insight into biological processes, protein activity and localization, and cell signaling networks.

Bioorthogonal Reporter Tags: Investigate Without Interference

Probes containing an alkyne, azide or tetrazine handle are paired with bioorthogonally compatible fluorescent or biotinylated tags

Baseclick EdU Cell Proliferation Monitoring Kits: Rapid Quantification of Cell Growth

Detect cell proliferation especially for *in vivo* experiments with the EdU nucleoside:

- Nontoxic to animals
- Directly incorporated into replicating DNA
- Contains an alkyne handle for click chemistry labeling with azide florescent probes

Boger Serine Hydrolase Inhibitors: An Irreversible Bind

Access a library of molecules from the Dale Boger lab with broad spectrum inhibition of the serine hydrolase family. Tune the compounds to a desired reactivity and selectivity through functional group attachment at an amine site and develop inhibitors to specific subclasses of serine hydrolase enzymes.

Quorum Sensing Modulators: Interfere with Bacterial Signaling

Quorum sensing is a chemical communication mechanism used by many common bacteria. New quorum sensing modulators developed by the Helen Blackwell laboratory are some of the most potent inhibitors and activators reported for Gram-negative species.

BIOCONJUGATION

Crosslinkers: A Variety of Linkers for a Variety of Biological Needs

We offer crosslinkers that vary in length, solubility and functional group reactivity.

- Common bioconjugation linkers including NHS esters for amines or maleimide for thiols
- Improved thiol coupling with APN reactivity
- Enhanced solubility with sulfo-NHS esters and PEGylated linkers
- Cleavable crosslinkers also available

We also provide crosslinkers designed for click chemistry reactivity.

- Azide and alkyne linkers for copper click chemistry
- Dibenzocyclooctyne (DBCO), azide, tetrazine or transcyclooctyne for copper-free click chemistry

Potassium Acyltrifluoroborates (KATs): A New Approach to Bioconjugation

As a unique class of functional groups, KATS:

- Offer innovative approaches to protein conjugation and hydrogel formation
- Undergo rapid, chemoselective amide-forming reactions under aqueous conditions without the need for activating agents or protecting groups

Late Stage Conjugation: Native Chemical Tagging of Bioactive Small Molecules

We offer Baran Diversinate crosslinkers with a metal sulfinate on one end to attach directly to otherwise inert C-H bonds and an azide on the other end for attaching to alkynes. We offer various alkyne and strained alkyne crosslinkers for attachment to the azide crosslinker.

QBD10501
Azido-dPEG®$_4$-NHS ester

17758
11-Azido-3,6,9-trioxaundecan-1-amine
technical, ≥90% (GC)

908401
1-Methyl-7-nitroisatoic anhydride

802409
7-Azido-4-Methylcoumarin

97%

QBD10200
NHS-dPEG®$_4$-biotin

QBD10503
Azido-dPEG®$_8$-NHS ester

736678
Tetra(ethylene glycol) dithiol
97%

A-PLSC010
AnteoBind™ Biosensor

QBD10211
m-dPEG®$_4$-NHS ester

QBD10109
Acid-dPEG®$_5$-NHS ester

900606
Fmoc-Trp-BODIPY
≥95%

QBD10502
Azido-dPEG®$_4$-acid

767751
2-{2-[2-(2-Mercaptoethoxy)ethoxy]ethoxy}ethanol
97%

92893
3,6,9-Trioxaundecanedioic acid
technical, ≥70% (T)

76318
***O*-(2-Aminoethyl)-*O'*-(2-azidoethyl)heptaethylene glycol**
≥90% (oligomer purity)

900572
Biotinylated isoxazole
95%

905283
(6-Aminohexyl)triphenylphosphonium bromide hydrobromide
≥95%

QBD10175
m-dPEG®$_4$-amine

QBD10249
Amino-dPEG®$_4$-OH

QBD11815
DBCO-TFP ester

672688
***O*-(2-Carboxyethyl)-*O'*-(2-mercaptoethyl)heptaethylene glycol**
≥95% (oligomer purity)

A-VMPAKMP
Universal Coupling Kit

QBD11217
Bromoacetamido-dPEG ®$_3$-azide

905267
(6-Bromohexyl)triphenylphosphonium bromide
≥95%

QBD11812
DBCO-dPEG ®$_{12}$-carboxyfluorescein

QBD10067
Amino-dPEG®$_6$-acid

07969
***O*-(2-Aminoethyl)polyethylene glycol 3,000**
M_p 3,000

689440
O-(2-Azidoethyl)heptaethylene glycol
≥95% (oligomer purity)

QBD10567
Azido-dPEG ®$_4$-TFP ester

QBD11100
Biotin-dPEG®$_3$-oxyamine HCl

QBD10314
MAL-dPEG®$_{24}$-NHS ester

QBD11028
MAL-dPEG ®$_{12}$-DSPE

QBD10278
m-dPEG®$_8$-amine

QBD10505
Azido-dPEG®$_{12}$-NHS ester

QBD11221
Bromoacetamido-dPEG ®$_4$-amido-DBCO

QBD10127
Acid-dPEG®$_{13}$-NHS ester

QBD11811
Biotin-dPEG ®$_{12}$-DBCO

QBD10569
Azido-dPEG ®$_{12}$-TFP ester
>90%

QBD10198
NHS-dPEG®$_{12}$-biotin

689777

O-[*N*-(3-Maleimidopropionyl)aminoethyl]-*O'*-[3-(*N*-succinimidyloxy)-3-oxopropyl]heptacosaethylene glycol
≥90% (oligomer purity)

76172
***O*-(2-Aminoethyl)-*O'*-(2-azidoethyl)pentaethylene glycol**
≥90% (oligomer purity)

QBD10774
NHS-dPEG®$_{24}$-biotin

QBD10219
Biotin-dPEG®$_4$-hydrazide

QBD10277
Amino-dPEG®$_8$-acid

QBD11167
DOTA-tris(acid)-amido-dPEG ®$_{11}$-Maleimide

712515
***O*-(3-Carboxypropyl)-*O'*-[2-(3-mercaptopropionylamino)ethyl]-polyethylene glycol**
M_w 3000

A-LMPAKMM
Activation Kit For Multiplex Microspheres

QBD10201
Biotin-dPEG®$_3$-MAL

QBD10593
DBCO-dPEG ®$_{12}$-MAL

QBD10523
Azido-dPEG®$_7$-amine

QBD10374
SPDP-dPEG®$_4$-NHS ester

QBD10784
Biotin-dPEG®$_{11}$-azide

689882
***O*-(2-Aminoethyl)-*O'*-[2-(biotinylamino)ethyl]octaethylene glycol**
≥95% (oligomer uniformity)

71613
O-(2-Azidoethyl)-O-[2-(diglycolyl-amino)ethyl]heptaethylene glycol
≥90% (oligomer purity)

JKA12021
Propionic acid-PEG4-Propionic acid

83060
***tert*-Butyl 12-amino-4,7,10-trioxadodecanoate**
technical, ≥80% (T)

QBD10317
Amino-dPEG®$_{24}$-acid

QBD11362
DBCO-dPEG ®$_4$-TFP ester

QBD10196
Biotin-dPEG®$_{11}$-NH2

QBD10551
***MAL*-dPEG®$_4$-TFP ester**

70023
***O*-(2-Aminoethyl)-*O'*-[2-(Boc-amino)ethyl]hexaethylene glycol**
≥90% (oligomer purity)

QBD10512
Azido-dPEG®$_8$-acid

QBD10262
m-dPEG®$_{12}$-NHS ester

QBD11171
DOTA-tris(acid)-amido-dPEG ®$_{23}$-Maleimide

QBD11155
DOTA-tris(TBE)-amido-dPEG ®$_{4}$-TFP ester

77787
***O*-(2-Aminoethyl)-*O'*-(2-azidoethyl)nonaethylene glycol**
≥90% (oligomer purity)

QBD10319
m-dPEG®$_{24}$-MAL

712523
***O*-(3-Carboxypropyl)-*O'*-[2-(3-mercaptopropionylamino)ethyl]-polyethylene glycol**
M_w 5000

QBD10525
Azido-dPEG®$_{23}$-amine

QBD11160
DOTA-tris(acid)-amido-dPEG ®$_{4}$-TFP ester

QBD10800
m-dPEG®$_{8}$-Lipoamide

900751
Biotinylated-D-lysine TFA salt

QBD10260
m-dPEG®$_{8}$-NHS ester

QBD10009
Biotin-dPEG®$_{4}$-TFP ester

QBD10318
m-dPEG®$_{24}$-amine

QBD11372
Bis-dPEG ®$_{11}$-DBCO

QBD11102
Biotin-dPEG®$_{11}$-oxyamine HCl

QBD10119
Acid-dPEG®$_{9}$-NHS ester

QBD10524
Azido-dPEG®$_{11}$-amine

QBD10194
NHS-S-S-dPEG®$_{4}$-biotin (cleavable)

QBD11150
DOTA-tris(acid)-amido-dPEG ®$_{3}$-bromoacetamide

QBD10289
m-dPEG®$_{12}$-MAL

QBD10826
Biotin-dPEG®$_{7}$-NH2

QBD11223
Bromoacetamido-dPEG ®$_{12}$-amido-DBCO

QBD10304
m-dPEG®$_{24}$-NHS ester

QBD11166
DOTA-tris(TBE)-amido dPEG ®$_{11}$-Maleimide

QBD11370
DBCO-dPEG ®$_{24}$-TFP ester

QBD10785
Biotin-dPEG®$_{23}$-MAL

QBD11371
Tetra(-dPEG ®$_{11}$-DBCO)PE

77090
***O*-(2-Aminoethyl)-*O'*-[2-(Boc-amino)ethyl]decaethylene glycol**
≥90% (oligomer purity)

QBD10215
Bis-MAL-dPEG®$_3$

QBD10240
Amino-dPEG®$_8$-OH

QBD10819
MAL-dPEG®$_{11}$-Lipoamide

QBD10181
dPEG®$_4$-SATA (S-acetyl-dPEG®$_4$-NHS ester)

QBD10786
Biotin-dPEG®$_{23}$-NH2

QBD10954
Bis-dPEG®$_{13}$-NHS ester

QBD10822
Biotin-dPEG®$_{11}$-Lipoamide

QBD10825
Biotin-dPEG®$_7$-azide

QBD10306
m-dPEG®$_{12}$-TFP ester

LIGHTOX14
LightOx™14
≥98% (HPLC)

QBD10184
dPEG®$_8$-SATA (S-acetyl-dPEG®$_8$-NHS ester)

QBD10406
MAL-dPEG®$_4$-(m-dPEG®$_{12}$)$_3$

QBD10745
m-dPEG®$_4$-MAL

QBD10807
Lipoamido-dPEG®$_8$-acid

689998
***O*-[2-(Biotinylamino)ethyl]-*O'*-(2-carboxyethyl)undecaethylene glycol**
≥95% (oligomer purity)

671592
***O*-(2-Aminoethyl)-*O'*-(2-carboxyethyl)polyethylene glycol 5,000 hydrochloride**
M_p 5,000

671487
***O*-(2-Aminoethyl)-*O'*-(2-carboxyethyl)polyethylene glycol hydrochloride**
M_p 3,000

QBD10852
dPEG®$_{12}$-SATA (S-acetyl-dPEG®$_{12}$-NHS ester)

QBD10361
Diamido-dPEG®$_{11}$-diamine

QBD11383
DBCO-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

QBD10288
m-dPEG®$_{12}$-amine

QBD10868
Amino-dPEG®$_{24}$-OH

79141
***O*-(2-Aminoethyl)-*O'*-[2-(Boc-amino)ethyl]octaethylene glycol**

≥90% (oligomer purity)

QBD10313
Fmoc-N-amido-dPEG®$_{24}$-acid

QBD11366
DBCO-dPEG ®$_{12}$-TFP ester

QBD10787
Biotin-dPEG®$_{23}$-azide

QBD10195
Biotin-dPEG®$_{11}$-MAL

QBD10594
DBCO-dPEG ®$_{24}$-MAL

QBD10540
m-dPEG ®$_{24}$-Azide (Azido-m-dPEG ®$_{24}$)

QBD10401
NHS-dPEG®$_4$-(m-dPEG®$_{12}$)$_3$-ester

QBD10397
Bis-MAL-dPEG®$_{11}$

QBD10526
Azido-dPEG ®$_{35}$-amine
>90%

QBD10596
m-dPEG ®$_{12}$-DBCO

QBD11200
Bromoacetamido-dPEG ®$_4$-TFP ester

QBD10170
Amino-dPEG®$_{12}$-OH

QBD10514
Azido-dPEG®$_{24}$-acid

QBD10907
Amino-dPEG®$_{36}$-acid

QBD11152
DOTA-tris(acid)-amido-dPEG ®$_{11}$-bromoacetamide

QBD10746
m-dPEG®$_{8}$-MAL

QBD11005
Fmoc-N-amido-dPEG ®$_{8}$-TFP ester

LIGHTOX19
LightOx™19
≥95% (HPLC)

LIGHTOX58
LightOx™58
≥95% (HPLC)

QBD10994
Fmoc-N-amido-dPEG ®$_{4}$-NHS ester

QBD10492
Aminooxy-dPEG®$_{12}$-amido-dPEG®$_{12}$-(m-dPEG®$_{11}$)$_3$

QBD10808
Lipoamido-dPEG®$_{12}$-acid

QBD11303
MAL-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-TFP ester

LIGHTOX72
LightOx™72
≥98% (HPLC)

QBD10008
Biotin-dPEG®$_{12}$-TFP ester

QBD11204
Bromoacetamido-dPEG ®$_{11}$-azide

QBD10814
Lipoamido-dPEG®$_{12}$-TFP ester

QBD11205
Bromoacetamido-dPEG ®$_{23}$-azide

QBD11224
Bromoacetamido-dPEG ®$_{24}$-amido-DBCO

QBD10376
SPDP-dPEG®$_{8}$-NHS ester

QBD10513
Azido-dPEG®$_{12}$-acid

LIGHTOX17
LightOx™17
≥98% (HPLC)

QBD10573
Azido-dPEG ®$_{36}$-TFP ester

QBD10075
4-formyl-benzamido-dPEG®$_{12}$-EDA-MAL

QBD10552
MAL-dPEG®$_{8}$-TFP ester

QBD10801
m-dPEG®$_{12}$-Lipoamide

QBD11414
Tetra(-amido-dPEG®$_{23}$-MAL)pentaerythritol

QBD10203
TFP-dPEG®$_4$-biotinidase resistant biotin

QBD10554
MAL-dPEG®$_{24}$-TFP ester

QBD11135
Phthalimidooxy-dPEG®$_{12}$-NHS ester

QBD10571
Azido-dPEG®$_{24}$-TFP ester

QBD11157
DOTA-tris(TBE)-amido-dPEG®$_{12}$-TFP ester

QBD11162
DOTA-tris(acid)-amido-dPEG®$_{12}$-TFP ester

QBD11384
DOTA-tris(acid)-amido-dPEG®$_{24}$-amido-dPEG®$_{24}$-DSPE

QBD10011
Phthalimidooxy-dPEG®$_4$-NHS ester

QBD10015
Bis-dPEG®$_5$-PFP ester

QBD10081
4-formyl-benzamido-dPEG®$_{12}$-TFP ester

QBD10204
TFP-dPEG®$_{12}$-biotinidase resistant biotin

QBD10007
Biotin-dPEG®$_{24}$-TFP ester

QBD10378

SPDP-dPEG®$_{12}$-NHS ester

QBD11341
Hydroxy-dPEG ®$_4$-TFP ester

QBD10995
Fmoc-N-amido-dPEG ®$_8$-NHS ester

QBD10996
Fmoc-N-amido-dPEG ®$_{12}$-NHS ester

QBD11170
DOTA tris(TBE)-amido-dPEG ®$_{23}$-Maleimide

QBD11435
Tetra(-amido-dPEG ®$_{11}$-MAL)pentaerythritol

QBD10303
m-dPEG®$_{24}$-TFP ester

QBD10287
Amino-dPEG®$_{12}$-acid

LIGHTOX21
LightOx™21
≥95% (HPLC)

LIGHTOX23
LightOx™23
≥95% (HPLC)

QBD11000
Fmoc-N-amido-dPEG ®$_4$-TFP ester

QBD11008
Fmoc-N-amido-dPEG ®$_{36}$-TFP ester

QBD11202
Bromoacetamido-dPEG ®$_{12}$-TFP ester

QBD11305

Amino-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-acid

QBD10553
MAL-dPEG®$_{12}$-TFP ester

QBD10804
m-dPEG®$_{24}$-Lipoamide

QBD10983
Bis-dPEG®$_{9}$-PFP ester

QBD11007
Fmoc-N-amido-dPEG ®$_{24}$-TFP ester

LIGHTOX26
LightOx™26
≥95% (HPLC)

QBD11086
m-dPEG ®$_{13}$-NHS ester

QBD11093
MAL-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

QBD11163
DOTA-tris(acid)-amido-dPEG ®$_{24}$-TFP ester

QBD10631
Bis-MAL-Lysine-dPEG®$_{4}$-TFP ester

QBD11373
Methoxy-dPEG ®$_{25}$-DSPE

QBD10379
SPDP-dPEG®$_{24}$-NHS ester

QBD10642
Lipoamido-dPEG®$_{8}$-TFP ester

LIGHTOX22
LightOx™22
≥95% (HPLC)

QBD10555
MAL-dPEG ®$_{36}$-TFP ester

QBD11006
Fmoc-N-amido-dPEG ®$_{12}$-TFP ester

QBD11087
m-dPEG ®$_{13}$-TFP ester

QBD11158
DOTA-tris(TBE)-amido-dPEG ®$_{24}$-TFP ester

QBD11345
Hydroxy-dPEG ®$_{12}$-TFP ester

QBD11349
Hydroxy-dPEG ®$_{24}$-TFP ester

QBD11385
Carboxyfluorescein-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

QBD11386
Biotin-dPEG ®$_{4}$ amido-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

LIGHTOX25
LightOx™25
≥95% (HPLC)

QBD11095
m-dPEG ®$_{25}$-amido-dPEG ®$_{24}$-DSPE

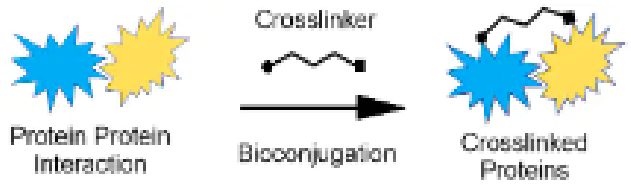
QBD11203
Bromoacetamido-dPEG ®$_{24}$-TFP ester

QBD11388
Mal-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

QBD10984
Bis-dPEG®$_{13}$-PFP ester

QBD11338
Bis-Bromoacetamido-dPEG ®$_{11}$

Crosslinkers



We are proud to offer you a wide variety of linkers and crosslinkers for all your biochemistry needs. Our protein reagents are suitable for stabilizing structures in protein-protein, protein-peptide and peptide/protein-small molecule interactions, in immobilizing proteins onto a solid support for assays or purification, as well as for various peptide-nucleic acid and nucleic acid-nucleic acid conjugations, among many other applications. Explore our extensive portfolio of versatile linkers and crosslinkers to find the most efficient and optimal reagent for your scientific breakthroughs.

HOMOBIFUNCTIONAL AND HETEROFUNCTIONAL LINKERS

Our homobifunctional linkers and heterofunctional linkers contain diverse functional groups, such as primary amines, sulfhydryls, acids, alcohols and bromides. Many of our linkers are functionalized with maleimide (sulfhydral reactive) and succinimidyl ester (NHS) or isothiocyanate (ITC) groups that react with amines. We also offer a broad selection of mono-protected (Boc, Fmoc, and Cbz) linkers.

CROSSLINKING REAGENTS

Crosslinking reagents are molecules containing two or more reactive ends that are 'activated' to attach themselves to certain functional groups (e.g., amines and sulfhydryls) via a covalent bond. The usefulness of crosslinking chemistry is realized in applications such as:

- protein structure and/or function determination
- protein or other biomolecule immobilization
- general biomolecule-biomolecule conjugation
- antibody-drug conjugates

Some of the most common crosslinkers contain maleimide, sulfhydryl reactive groups, or succinimidyl esters (often referred to as NHS esters), which react with amines. Our portfolio comprises all functional groups with a variety of linker lengths and solubilities. Sulfosuccinimidyl esters allow for more water-soluble crosslinkers which can be useful when working with large biomolecules that are not amenable to organic solvents. Our crosslinkers with cleavable linkers (e.g., disulfides) are optimal for applications where a permanent linkage is not desired.

SELECTING CROSSLINKERS

When selecting a crosslinker for your application it is important to consider several factors, such as the reagent's chemical and physical properties, the functional groups it targets for coupling, its length, its molecular size, its water solubility, and its cleavability:

Chemical Specificity: One of the most fundamental aspects of crosslinker design is whether the reagent is homobifunctional or heterobifunctional. Homobifunctional compounds will react at both ends with the same target functional group, thus forming a covalent crosslink between two molecules using the same type of bond. They are used in single-step reactions for polymerization of like functional groups, in the creation of intramolecular crosslinks, and in the evaluation of protein

interactions. Heterobifunctional reagents have two different end groups, allowing for each end to react with a different functional group. They are used for controlled two-step reactions to avoid undesirable cross-reactions and polymerization.

Functional Groups Targeted: Amines, thiols, and hydroxyls are the main nucleophilic groups, and under the right conditions, they react directly with the electrophilic reactive groups present on many bioconjugation reagents. In contrast, functional groups consisting of carboxylates, aldehydes, organic phosphates, and reactive hydrogen sites require special activation agents or secondary coupling agents before they form covalent bonds with another functional group.

Crosslinker Length: The dimensions or overall linear length of the target molecule before and after conjugation should be considered when choosing a crosslinker. The spacer arm or cross-bridge of the reagent mainly determines the molecular length of the resulting compound. This length can be determined by use of certain molecular modeling software programs.

Cleavable vs. Noncleavable Crosslinkers: If interacting biomolecules that have been captured by crosslinking subsequently need to be isolated and analyzed, it is important for the spacer arm of the crosslinker to be cleavable. Additionally, cleavable linkers are used as transfer reagents for the study of interacting proteins. Our disulfides are some of the most common cleavable crosslinkers. Our portfolio also offers cleavable esters and sulfones.

Hydrophobic vs Hydrophilic Crosslinkers: In some applications, reagent hydrophobicity can be an advantage, especially when an application involves penetration of cell membranes. Hydrophobic reagents without strongly polar groups can quickly pass through membranes and crosslink or label internal cell proteins. However, hydrophobic compounds that contain one or more negatively charged sulfo-NHS groups will be restricted to reacting with the proteins on the outer membrane surface of cells due to their negative charge. The ability to switch between cell surface labeling and internal cellular labeling by choosing charged or uncharged reagents is one benefit of using hydrophobic crosslinkers.

744867
(1*R*,8*S*,9*s*)-Bicyclo[6.1.0]non-4-yn-9-ylmethyl *N*-succinimidyl carbonate
for Copper-free Click Chemistry

40421
1,11-Diamino-3,6,9-trioxaundecane
≥98.0% (GC)

765953
1,11-Diazido-3,6,9-trioxaundecane

17758
11-Azido-3,6,9-trioxaundecan-1-amine
technical, ≥90% (GC)

30953
15-(Boc-amino)-4,7,10,13-tetraoxapentadecanoic acid
purum, ≥97.0% (TLC)

767751
2-{2-[2-(2-Mercaptoethoxy)ethoxy]ethoxy}ethanol
97%

458910
2-(Boc-amino)ethanethiol
97%

17354
2-(Boc-amino)ethyl bromide
≥97.0% (GC)

445185
2-(Fmoc-amino)ethanol
97%

74291
2-(Fmoc-amino)ethyl bromide
≥95.0% (AT)

41314
2-Maleimidoethyl mesylate
technical, ≥90.0% (HPLC)

P3415
3-(2-Pyridyldithio)propionic acid *N*-hydroxysuccinimide ester
≥95%, powder

906298
3-(4-Bromophenyl)-3-(trifluoromethyl)-3*H*-diazirine

416444
3-(Boc-amino)-1-propanol
97%

17356
3-(Boc-amino)propyl bromide
≥96.0% (GC)

M2786
3-Maleimidobenzoic acid *N*-hydroxysuccinimide ester
crystalline

D3669
3,3′-Dithiodipropionic acid di(N-hydroxysuccinimide ester)
powder

55665
3,4-Dihydro-2*H*-pyran-2-methanol
≥98.5% (GC)

92893
3,6,9-Trioxaundecanedioic acid
technical, ≥70% (T)

55262
4-[4-(1-Hydroxyethyl)-2-methoxy-5-nitrophenoxy]butyric acid
≥98.0% (HPLC)

771058
4-(4′-Hydroxyphenylazo)benzoic acid
97%

900602
4-(4-(Prop-2-yn-1-yloxy)benzoyl)benzoic acid
≥95%

15302
4-(Boc-amino)-1-butanol
≥98.0% (GC)

90303
4-(Boc-amino)butyl bromide
technical, ≥90% (AT)

43081
4-(Diphenylhydroxymethyl)benzoic acid
≥98.0% (HPLC)

382639
4-(Hydroxymethyl)benzoic acid
99%

95887
4-(Z-Amino)-1-butanol
≥98.0% (HPLC)

A44150
4-Aminobutyraldehyde diethyl acetal
90%, technical grade

QBD10075
4-formyl-benzamido-dPEG®$_{12}$-EDA-MAL

QBD10081
4-formyl-benzamido-dPEG®$_{12}$-TFP ester

C11804
4-Sulfamoylbenzoic acid
97%

671630
4,7,10,13,16,19,22,25,32,35,38,41,44,47,50,53-Hexadecaoxa-28,29-dithiahexapentacontanedioic acid di-*N*-suc cinimidyl ester
≥94% (oligomer purity)

47299
5-(Fmoc-amino)-1-pentanol
≥98.0% (HPLC)

15304
6-(Boc-amino)-1-hexanol
≥98.0% (GC)

89171
6-(Boc-amino)hexyl bromide
≥97.0% (GC)

568600
6-Maleimidohexanoic acid *N*-hydroxysuccinimide ester
98%

QBD10127
Acid-dPEG®$_{13}$-NHS ester

QBD10109
Acid-dPEG®$_{5}$-NHS ester

QBD10119
Acid-dPEG®$_{9}$-NHS ester

764205
Alkyne-PEG4-maleimide
90% (HPLC)

764167
Alkyne-PEG5-acid

764191
Alkyne-PEG5-*N*-hydroxysuccinimidyl ester

QBD10287
Amino-dPEG®$_{12}$-acid

QBD10170
Amino-dPEG®$_{12}$-OH

QBD10317
Amino-dPEG®$_{24}$-acid

QBD11305
Amino-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-acid

QBD10868
Amino-dPEG®$_{24}$-OH

QBD10907
Amino-dPEG®$_{36}$-acid

QBD10249
Amino-dPEG®$_{4}$-OH

QBD10067
Amino-dPEG®$_{6}$-acid

QBD10277
Amino-dPEG®$_{8}$-acid

QBD10240
Amino-dPEG®$_{8}$-OH

764248
Amino-PEG4-alkyne

902535
Amino-PEG4-t-butyl ester

902659
Amino-PEG6-t-butyl ester

QBD10492
Aminooxy-dPEG®$_{12}$-amido-dPEG®$_{12}$-(m-dPEG®$_{11}$)$_3$

806595
APN-Azide
95%

807745
APN-COCl

QBD10524
Azido-dPEG®$_{11}$-amine

QBD10513
Azido-dPEG®$_{12}$-acid

QBD10505
Azido-dPEG®$_{12}$-NHS ester

QBD10569
Azido-dPEG ®$_{12}$-TFP ester
>90%

QBD10525
Azido-dPEG®$_{23}$-amine

QBD10514
Azido-dPEG®$_{24}$-acid

QBD10571
Azido-dPEG ®$_{24}$-TFP ester

QBD10526
Azido-dPEG ®$_{35}$-amine
>90%

QBD10573
Azido-dPEG ®$_{36}$-TFP ester

QBD10502
Azido-dPEG®$_{4}$-acid

QBD10501
Azido-dPEG®$_{4}$-NHS ester

QBD10567
Azido-dPEG ®$_{4}$-TFP ester

QBD10523
Azido-dPEG®$_{7}$-amine

QBD10512
Azido-dPEG®$_8$-acid

QBD10503
Azido-dPEG®$_8$-NHS ester

478709
Benzyl *N*-(3-hydroxypropyl)carbamate
97%

793329
Biotin-benzyl-tetrazine
95%

QBD10784
Biotin-dPEG®$_{11}$-azide

QBD10822
Biotin-dPEG®$_{11}$-Lipoamide

QBD10195
Biotin-dPEG®$_{11}$-MAL

QBD10196
Biotin-dPEG®$_{11}$-NH2

QBD11102
Biotin-dPEG®$_{11}$-oxyamine HCl

QBD11811
Biotin-dPEG ®$_{12}$-DBCO

QBD10008
Biotin-dPEG®$_{12}$-TFP ester

QBD10787
Biotin-dPEG®$_{23}$-azide

QBD10785
Biotin-dPEG®$_{23}$-MAL

QBD10786
Biotin-dPEG®$_{23}$-NH2

QBD10007
Biotin-dPEG®$_{24}$-TFP ester

QBD10201
Biotin-dPEG®$_{3}$-MAL

QBD11100
Biotin-dPEG®$_{3}$-oxyamine HCl

QBD11386
Biotin-dPEG ®$_{4}$ amido-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

QBD10219
Biotin-dPEG®$_{4}$-hydrazide

QBD10009
Biotin-dPEG®$_{4}$-TFP ester

QBD10825
Biotin-dPEG®$_{7}$-azide

QBD10826
Biotin-dPEG®$_{7}$-NH2

QBD11338
Bis-Bromoacetamido-dPEG ®$_{11}$

QBD11372
Bis-dPEG ®$_{11}$-DBCO

QBD10954
Bis-dPEG®$_{13}$-NHS ester

QBD10984
Bis-dPEG®$_{13}$-PFP ester

QBD10015
Bis-dPEG®$_{5}$-PFP ester

QBD10983
Bis-dPEG®$_9$-PFP ester

QBD10397
Bis-MAL-dPEG®$_{11}$

QBD10215
Bis-MAL-dPEG®$_3$

QBD10631
Bis-MAL-Lysine-dPEG®$_4$-TFP ester

803537
Bis(NHS)PEG5

902551
BocNH-PEG4-acid

902683
BocNH-PEG5-acid

QBD11204
Bromoacetamido-dPEG ®$_{11}$-azide

QBD11223
Bromoacetamido-dPEG ®$_{12}$-amido-DBCO

QBD11202
Bromoacetamido-dPEG ®$_{12}$-TFP ester

QBD11205
Bromoacetamido-dPEG ®$_{23}$-azide

QBD11224
Bromoacetamido-dPEG ®$_{24}$-amido-DBCO

QBD11203
Bromoacetamido-dPEG ®$_{24}$-TFP ester

QBD11217
Bromoacetamido-dPEG ®$_3$-azide

QBD11221
Bromoacetamido-dPEG ®$_4$-amido-DBCO

QBD11200
Bromoacetamido-dPEG ®$_4$-TFP ester

B8271
Bromoacetic acid N-hydroxysuccinimide ester
≥95%, powder

803464
BS(PEG)9 (PEGylated bis(sulfosuccinimidyl)suberate)

QBD11385
Carboxyfluorescein-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

807664
CBTF

QBD11812
DBCO-dPEG ®$_{12}$-carboxyfluorescein

QBD10593
DBCO-dPEG ®$_{12}$-MAL

QBD11366
DBCO-dPEG ®$_{12}$-TFP ester

QBD11383
DBCO-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

QBD10594
DBCO-dPEG ®$_{24}$-MAL

QBD11370
DBCO-dPEG ®$_{24}$-TFP ester

QBD11362
DBCO-dPEG ®$_4$-TFP ester

QBD11815
DBCO-TFP ester

17353
di-Boc-cystamine
≥98.0% (TLC)

QBD10361
Diamido-dPEG®$_{11}$-diamine

760668
Dibenzocyclooctyne-maleimide
for Copper-free Click Chemistry

761524
Dibenzocyclooctyne-*N*-hydroxysuccinimidyl ester
for Copper-free Click Chemistry

760676
Dibenzocyclooctyne-PEG4-maleimide
for Copper-free Click Chemistry

D2388
Dimethyl 3,3′-dithiopropionimidate dihydrochloride
powder

QBD11170
DOTA tris(TBE)-amido-dPEG ®$_{23}$-Maleimide

QBD11152
DOTA-tris(acid)-amido-dPEG ®$_{11}$-bromoacetamide

QBD11167
DOTA-tris(acid)-amido-dPEG ®$_{11}$-Maleimide

QBD11162
DOTA-tris(acid)-amido-dPEG ®$_{12}$-TFP ester

QBD11171
DOTA-tris(acid)-amido-dPEG ®$_{23}$-Maleimide

QBD11384
DOTA-tris(acid)-amido-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

QBD11163
DOTA-tris(acid)-amido-dPEG ®$_{24}$-TFP ester

QBD11150
DOTA-tris(acid)-amido-dPEG ®$_3$-bromoacetamide

QBD11160
DOTA-tris(acid)-amido-dPEG ®$_4$-TFP ester

QBD11166
DOTA-tris(TBE)-amido dPEG ®$_{11}$-Maleimide

QBD11157
DOTA-tris(TBE)-amido-dPEG ®$_{12}$-TFP ester

QBD11158
DOTA-tris(TBE)-amido-dPEG ®$_{24}$-TFP ester

QBD11155
DOTA-tris(TBE)-amido-dPEG ®$_4$-TFP ester

QBD10852
dPEG®$_{12}$-SATA (S-acetyl-dPEG®$_{12}$-NHS ester)

QBD10181
dPEG®$_4$-SATA (S-acetyl-dPEG®$_4$-NHS ester)

QBD10184
dPEG®$_8$-SATA (S-acetyl-dPEG®$_8$-NHS ester)

803200
DTSSP (3,3′-dithiobis(sulfosuccinimidyl propionate))

803677
EMCH (N-(ε-maleimidocaproic acid) hydrazide, trifluoroacetic acid salt)

QBD10996

Fmoc-N-amido-dPEG ®12-NHS ester

QBD11006
Fmoc-N-amido-dPEG ®12-TFP ester

QBD10313
Fmoc-N-amido-dPEG®24-acid

QBD11007
Fmoc-N-amido-dPEG ®24-TFP ester

QBD11008
Fmoc-N-amido-dPEG ®36-TFP ester

QBD10994
Fmoc-N-amido-dPEG ®4-NHS ester

QBD11000
Fmoc-N-amido-dPEG ®4-TFP ester

QBD10995
Fmoc-N-amido-dPEG ®8-NHS ester

QBD11005
Fmoc-N-amido-dPEG ®8-TFP ester

QBD11345
Hydroxy-dPEG ®12-TFP ester

QBD11349
Hydroxy-dPEG ®24-TFP ester

QBD11341
Hydroxy-dPEG ®4-TFP ester

902586
Hydroxy-PEG4-t-butyl ester

902578
Hydroxy-PEG6-t-butyl ester
≥95%

I9760
Iodoacetic acid *N*-hydroxysuccinimide ester
powder

803421
LC-SDA (NHS-LC-Diazirine) (succinimidyl 6-(4,4′-azipentanamido)hexanoate)

803383
LC-SMCC (succinimidyl-4-(N-maleimidomethyl)cyclohexane-1-carboxy-(6-amidocaproate))

803642
LC-SPDP (succinimidyl 6-[3(2-pyridyldithio)propionamido]hexanoate)

QBD10808
Lipoamido-dPEG®$_{12}$-acid

QBD10814
Lipoamido-dPEG®$_{12}$-TFP ester

QBD10807
Lipoamido-dPEG®$_{8}$-acid

QBD10642
Lipoamido-dPEG®$_{8}$-TFP ester

QBD10288
m-dPEG®$_{12}$-amine

QBD10596
m-dPEG ®$_{12}$-DBCO

QBD10801
m-dPEG®$_{12}$-Lipoamide

QBD10289
m-dPEG®$_{12}$-MAL

QBD10262
m-dPEG®$_{12}$-NHS ester

QBD10306
m-dPEG®$_{12}$-TFP ester

QBD11086
m-dPEG ®13-NHS ester

QBD11087
m-dPEG ®13-TFP ester

QBD10318
m-dPEG®24-amine

QBD10540
m-dPEG ®24-Azide (Azido-m-dPEG ®24)

QBD10804
m-dPEG®24-Lipoamide

QBD10319
m-dPEG®24-MAL

QBD10304
m-dPEG®24-NHS ester

QBD10303
m-dPEG®24-TFP ester

QBD11095
m-dPEG ®25-amido-dPEG ®24-DSPE

QBD10175
m-dPEG®4-amine

QBD10745
m-dPEG®4-MAL

QBD10211
m-dPEG®4-NHS ester

QBD10278
m-dPEG®8-amine

QBD10800
m-dPEG®$_8$-Lipoamide

QBD10746
m-dPEG®$_8$-MAL

QBD10260
m-dPEG®$_8$-NHS ester

QBD10819
MAL-dPEG®$_{11}$-Lipoamide

QBD11028
MAL-dPEG ®$_{12}$-DSPE

QBD10553
MAL-dPEG®$_{12}$-TFP ester

QBD11388
Mal-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

QBD11093
MAL-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-DSPE

QBD11303
MAL-dPEG ®$_{24}$-amido-dPEG ®$_{24}$-TFP ester

QBD10314
MAL-dPEG®$_{24}$-NHS ester

QBD10554
MAL-dPEG®$_{24}$-TFP ester

QBD10555
MAL-dPEG ®$_{36}$-TFP ester

QBD10406
MAL-dPEG®$_4$-(m-dPEG®$_{12}$)$_3$

QBD10551
MAL-dPEG®$_4$-TFP ester

QBD10552
MAL-dPEG®$_8$-TFP ester

746215
Maleimide-PEG$_{12}$-succinimidyl ester

746223
Maleimide-PEG$_2$-succinimidyl ester
≥95%

746193
Maleimide-PEG$_6$-succinimidyl ester

746207
Maleimide-PEG$_8$-succinimidyl ester

QBD11373
Methoxy-dPEG ®$_{25}$-DSPE

900914
Methyltetrazine-NHS ester
≥95%

56951
***N*-(2-Aminoethyl)maleimide trifluoroacetate salt**
≥95% (HPLC), ≥98% (T)

B66302
***N*-(2-Bromoethyl)phthalimide**
95%

773263
N-(2-Hydroxyethyl)maleimide
97%

440507
***N*-(2-Hydroxyethyl)trifluoroacetamide**
97%

B80003
***N*-(3-Bromopropyl)phthalimide**
98%

100919

***N*-(4-Bromobutyl)phthalimide**
98%

89761
***N*-Boc- 2,2′-(ethylenedioxy)diethylamine**
≥95.0% (NT)

15408
***N*-Boc-1,3-propanediamine**
≥97.0% (GC/NT)

15404
***N*-Boc-1,4-butanediamine**
≥97.0% (GC/NT)

79229
***N*-Boc-1,6-hexanediamine**
≥98.0% (NT)

437018
***N*-Boc-1,6-hexanediamine hydrochloride**
97%

93113
***N*-Boc-4,7,10-trioxa-1,13-tridecanediamine**
≥95.0% (NT)

15406
***N*-Boc-cadaverine**
≥97.0% (NT)

382027
***N*-Boc-ethanolamine**
98%

15369
***N*-Boc-ethylenediamine**
≥98.0% (NT)

53175
***N*-Boc-*m*-phenylenediamine**
≥98.0% (HPLC)

671401
***N*-Boc-*N*′-succinyl-4,7,10-trioxa-1,13-tridecanediamine**
95% (HPLC)

15485
***N*-Boc-*p*-phenylenediamine**
≥97.0% (NT)

47543
***N*-Fmoc-1,6-hexanediamine hydrobromide**
≥98.0% (AT)

47542
***N*-Fmoc-ethylenediamine hydrobromide**
≥98.0% (AT)

726303
N^1,N^4-Bis-Boc-spermidine
≥95.0% (TLC)

QBD10198
NHS-dPEG®$_{12}$-biotin

QBD10774
NHS-dPEG®$_{24}$-biotin

QBD10401
NHS-dPEG®$_4$-(m-dPEG®$_{12}$)$_3$-ester

QBD10200
NHS-dPEG®$_4$-biotin

QBD10194
NHS-S-S-dPEG®$_4$-biotin (cleavable)

689998
***O*-[2-(Biotinylamino)ethyl]-*O'*-(2-carboxyethyl)undecaethylene glycol**
≥95% (oligomer purity)

689777
O-[*N*-(3-Maleimidopropionyl)aminoethyl]-*O'*-[3-(*N*-succinimidyloxy)-3-oxopropyl]heptacosaethylene glycol
≥90% (oligomer purity)

712582
O-[*N*-(3-Maleimidopropionyl)aminoethyl]-*O'*-[3-(*N*-succinimidyloxy)-3-oxopropyl]triethylene glycol
≥90% (NMR)

689882
***O*-(2-Aminoethyl)-*O'*-[2-(biotinylamino)ethyl]octaethylene glycol**

≥95% (oligomer uniformity)

77090
***O*-(2-Aminoethyl)-*O'*-[2-(Boc-amino)ethyl]decaethylene glycol**
≥90% (oligomer purity)

70023
***O*-(2-Aminoethyl)-*O'*-[2-(Boc-amino)ethyl]hexaethylene glycol**
≥90% (oligomer purity)

79141
***O*-(2-Aminoethyl)-*O'*-[2-(Boc-amino)ethyl]octaethylene glycol**
≥90% (oligomer purity)

76318
***O*-(2-Aminoethyl)-*O'*-(2-azidoethyl)heptaethylene glycol**
≥90% (oligomer purity)

77787
***O*-(2-Aminoethyl)-*O'*-(2-azidoethyl)nonaethylene glycol**
≥90% (oligomer purity)

76172
***O*-(2-Aminoethyl)-*O'*-(2-azidoethyl)pentaethylene glycol**
≥90% (oligomer purity)

671592
***O*-(2-Aminoethyl)-*O'*-(2-carboxyethyl)polyethylene glycol 5,000 hydrochloride**
M_p 5,000

671487
***O*-(2-Aminoethyl)-*O'*-(2-carboxyethyl)polyethylene glycol hydrochloride**
M_p 3,000

07969
***O*-(2-Aminoethyl)polyethylene glycol 3,000**
M_p 3,000

71613
O-(2-Azidoethyl)-O-[2-(diglycolyl-amino)ethyl]heptaethylene glycol
≥90% (oligomer purity)

689440

O-(2-Azidoethyl)heptaethylene glycol
≥95% (oligomer purity)

672688
***O*-(2-Carboxyethyl)-*O'*-(2-mercaptoethyl)heptaethylene glycol**
≥95% (oligomer purity)

712523
***O*-(3-Carboxypropyl)-*O'*-[2-(3-mercaptopropionylamino)ethyl]-polyethylene glycol**
M_w 5000

712515
***O*-(3-Carboxypropyl)-*O'*-[2-(3-mercaptopropionylamino)ethyl]-polyethylene glycol**
M_w 3000

713791
***O*,*O'*-Bis[2-(*N*-Succinimidyl-succinylamino)ethyl]polyethylene glycol**
10,000

713783
***O*,*O'*-Bis[2-(*N*-Succinimidyl-succinylamino)ethyl]polyethylene glycol**
2,000

258555
***p*-Phenylene diisothiocyanate**
98%

901767
PC biotin-PEG3-NHS ester

803480
PDPH (3-(2-pyridyldithio)propionyl hydrazide)

803499
PEG4-SPDP (PEGylated, long-chain SPDP crosslinker)

QBD11135
Phthalimidooxy-dPEG®$_{12}$-NHS ester

QBD10011
Phthalimidooxy-dPEG®$_4$-NHS ester

764221
Propargyl-*N*-hydroxysuccinimidyl ester

902667

Propargyl-PEG6-acid

ALD00344
PTAD-Azide
95%

803650
SBAP (succinimidyl 3-(bromoacetamido)propionate)

803413
SDA (NHS-Diazirine) (succinimidyl 4,4′-azipentanoate)

803391
SM(PEG)24 (PEGylated, long-chain SMCC crosslinker)

803626
SMPH (succinimidyl-6-((b-maleimidopropionamido)hexanoate)

803472
SMPT (4-succinimidyloxycarbonyl-alpha-methyl-alpha(2-pyridyldithio)toluene)

QBD10378
SPDP-dPEG®$_{12}$-NHS ester

QBD10379
SPDP-dPEG®$_{24}$-NHS ester

QBD10374
SPDP-dPEG®$_{4}$-NHS ester

QBD10376
SPDP-dPEG®$_{8}$-NHS ester

S5799
Suberic acid bis(3-sulfo-N-hydroxysuccinimide ester) sodium salt
≥95% (H-NMR), powder

S1885
Suberic acid bis(N-hydroxysuccinimide ester)
≥95%, powder

760706
Sulfo-dibenzocyclooctyne-biotin conjugate
for Copper-free Click Chemistry

803219
Sulfo-EGS (ethylene glycol bis(sulfosuccinimidyl succinate))

803359
Sulfo-LC-SDA (Sulfo-NHS-LC-Diazirine) (sulfosuccinimidyl 6-(4,4′-azipentanamido)hexanoate)

803316
Sulfo-LC-SPDP (sulfosuccinimidyl 6-[3′-(2-pyridyldithio)propionamido]hexanoate)

803332
Sulfo-SANPAH (sulfosuccinimidyl 6-(4′-azido-2′-nitrophenylamino)hexanoate)

803340
Sulfo-SDA (Sulfo-NHS-Diazirine) (sulfosuccinimidyl 4,4′-azipentanoate)

803243
Sulfo-SMPB (sulfosuccinimidyl 4-(N-maleimidophenyl)butyrate)

900927
TCO-PNB ester
≥95%

83060
***tert*-Butyl 12-amino-4,7,10-trioxadodecanoate**
technical, ≥80% (T)

QBD11435
Tetra(-amido-dPEG ®$_{11}$-MAL)pentaerythritol

QBD11414
Tetra(-amido-dPEG ®$_{23}$-MAL)pentaerythritol

QBD11371
Tetra(-dPEG ®$_{11}$-DBCO)PE

QBD10204
TFP-dPEG®$_{12}$-biotinidase resistant biotin

QBD10203
TFP-dPEG®$_4$-biotinidase resistant biotin

CF0041
Urea crosslinker – C$_4$-arm, NHS ester (DSBU, BuUrBU)

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