

FT- AYPMB0

Monofunctional Pegylation agents

(mPEG-X – MethylOxyPEG-X, Methoxy-PEG-X)

Presentation:

These Monofunctional PEG derivatives can be used to modify peptides and proteins and other materials, to create conjugates or modify surfaces. These Pegylation agents are available with a variety chemical reactivities (one or 2 per molecule) and with lengths of the PEG hydrophilic spacer ranging from 350Da to 40 000Da (40KDa).

- Compared with other spacers, PEG structure increases water-solubility and stability. It also can reduce immunogenicity of conjugates and suppress the non-specific binding of charged molecules to the modified surfaces.
- mPEG agents have a PEG terminus blocked with a methoxy group (or methyloxy).
- The functional groups can react on specific chemical groups to create a covalent bond between the PEG spacer and the target compound, by conventional chemistry.

See information about functional groups and their reactivity

Product name synonymes	Cat.number Qty ^{0-250mg} ,	MW (g/mol ⁻¹)	Structure
synonymes	1-1g	(8/11101)	
• AMINE			
mPEG – Amine	B36BR1	350	O ₁ NH ₂
Methoxy-PEG-Amine, PEG- NH ₂	B36BU1 B36BV1	550 750	of _n var
PG1-AM (L/M)	WT9821	1 000	
[B] [N]	WT9811	2 000	See sets of 3 or 6 MW #AWK8Q-
Soluble in regular aqueous solution as	B36BT1	3 400	See HCl salt #(2K) [B]
well as most organic solvents.	WT9801	5 000	
	WT9831	10 000	
	WT9841	20 000	
	WT9851	30 000	, [D /N]
	JQ5081	40 000	
mPEG – AminoOxy	AWK7U0	550	
Methoxy-PEG-AO	B33UJ1	5 000	See also mPEO _n -AminoOxy (<u>FT-AWK7U0</u>)
PG1-AO (L/M)	B33UM1	30 000	
			See also HomoBifunctional Amine-PEG _x -Amine (<u>FT-WU0101</u>)
• CARBOXYL			
mPEG - COOH	B36C21 B36C31	350 550	0
Methoxy-PEG-COOH, PEG-	B36C41	750	0/ 0 10
Proprionic acid, PEG-(CH ₂) ₂ -COOH	WT9881	1 000	/ O/ OH
PG1- CA (L) [B] [N]	WT9871	2 000	, ""
Soluble in regular aqueous solution as	JO4951	3 400	
well as most organic solvents.	WT9861	5 000	C 1 mPFO Conhamil#
	WT9891	10 000	See also mPEO _n -Carboxyl #
	WT9901	20 000	
	WT9911	30 000	+
	B36C51	40 000	0
Ask also for:			O CONTRACTOR OH
mPEG-Acetic Acid (mPEG-AA):			mPEG-Succinamide Acid (mPEG-SAA):
CH ₃ O—(CH ₂ CH ₂ O)n OH			O (O) OH
Ö			mPEG-Glutaramide Acid (mPEG-GAA):
mPEG-Succinic Acid (mPEG-SA):			a/
о (о) п о			mPEG-Myristic acid (mPEG-MTA):
mPEG-Glutaric Acid (mPEG-GA):			
			See HomoBifunctional Carboxyl-PEG _x -Carboxyl (FT-WU0101)
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Product name	Cat.number	MW	Structure
synonymes	Qty 0-250mg,	(g/mol ⁻¹)	
	1-1g		
• NHS (Succinimide Ester)			_
mPEGx-NHSuccinumide Ester	DZ3511	1 000	Ö Ö
mPEG-SCM, PG1-NS (M)	DZ3521	2 000	
	DZ3531	5 000	CH ₃ O-(CH ₂ CH ₂ O) _n OH
	DZ3541	10 000	, , , , , , , , , , , , , , , , , , , ,
	DZ3551	20 000	See also mPEO _n -NHSuccinimide #RPX58G(PEG16)
	DZ3561	30 000	#8X543/C (mPEG ₅ -t-Butyl ester)
	HQ3581	40 000	#B4K6L (mPEG ₃ -S-PEG1-tButyl ester)
			#IO508/N (mPEG-Succinimidyl Glutarate Est MW 1000Da)
• MALEIMIDE			The second of th
mPEGx-Maleimide	DZ5781	1 000	
mPEG-MAL, PG1-ML(M)	DZ5761	2 000	
IIII EG-WIAE, I GI-WIE	DZ5781	10 000	
	DZ5781 DZ579	20 000	
	DZ579	30 000	See also mPEO_n-Maleimide #A1LYG2 (PEG ₄), RJ2021(PEG ₂₄),
	GV7341	40 000	
• AZIDE	0 1 / 341	40 000	
mPEG – Azide	AWJLI1	350	0 / \(\tau \) \(\tau \)
	Inquire	550	0/n \ 13
Methoxy-PEG-Azide, PEG-Azide,	Inquire	750	· ''' .
mPEG-AZ, PG1-AZ (M) [B][N]	WU0001	1 000	
	WT9991	2 000	
Soluble in regular aqueous solution as	WT9981	3 400	
well as most organic solvents.	WT9971	5 000	See also mPEO _n -Azide #6W540
	WU0011	10 000	
	WU0021	20 000	
	WU0031	30 000	
	B3AB21	40 000	+ [<u>D</u> N]
mPEG-DBCO	B4YBD-	350	^ Ā ^
(K,M) PG1-DB	B4YBB-	550	
C William William C 10YO	B36C81	1 000	
Appearance : White crystalline (for 10K) Soluble in Water, DMSO, DMF, chloroform	1Q7031	2 000	N V
BMF001048N	IOT811	5 000	$\frac{1}{2}$
	IOT821	10 000	
*1-100mg 2-1g	IOT831	20 000	Ö
2-1g	IOT841	30 000	See FT-B36C81 with other PEG DBCO reagents.
	101011	20 000	See HomoBifunctional Azide-PEG _x -Azide (<u>FT-WU0101</u>)
• HYDRAZIDE			
mPEG – Hydrazide	Inquire	350 to 750Da	
Methoxy-PEG- oxo -hydrazide,	AYPM71	1 000	
mPEG-HZ, mPEG-HAZ	AYPM81	2 000	NHNH ₂
PG1-HAZ (M)	AYPM91	3 400	CH ₃ O—(CH ₂ CH ₂ O)n
[B]	AYPMA1	4 000	
Soluble in regular aqueous solution as	AYPMB1	5 000	0
well as most organic solvents.	AYPMC1	10 000	
	AYPMD1	20 000	See also mPEO _n -Hydrazide #AYQU01(PEO ₂) to AYQUB0(PEO ₃
	AYPME1	30 000	
	AYPME1	40 000	$+\frac{ \mathbf{D}/\mathbf{N} }{ \mathbf{D} }$
	AITWILI	40 000	See HomoBifunctional Hydrazide-PEG _x -Hydrazide FT-WU0101
• ALDEHVDE			See Homodiffunctional Hydrazide-LEOx-Hydrazide F1-W00101
• ALDEHYDE			



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MPEG - Aldehyde Methoxy-PEG-Aldehyde, mPEG-CHO, mPEG-CH2-CHO MT8931 1 000 WT8911 5 000 WT8911	FT- AYPMB0			
1g	Product name	Cat.number	MW	Structure
MPEG - Aldehyde Methoxy-PEG-Aldehyde, mPEG-CH2-CHO MT8931 1 000 WT8911 5 000 WT8911 5 000 WT8911 5 000 WT8911 20 000 AYRIL1 5 000 AYRIL1 5 000 AYRIL1 5 000 AYRIL1 20 000 B36BP. WT8911 WT8911 20 000 B36BP. WT8911 20 000 B36BP. WT8911 WT8911 20 000 B36BP. WT8911	synonymes		(g/mol ⁻¹)	
Methoxy-PEG-Aldehyde, mPEG-CHO, mPEG-CH2-CHO PG1-AL M WT8911 Soluble in regular aqueous solution as well as most organic solvents. ■ METHOD Soluble in regular aqueous solution as well as most organic solvents. ■ METHOD METH		¹⁻ 1g		
CHO, mPEG-CH2-CHO WT8911 Soluble in regular aqueous solution as well as most organic solvents. WT8951 20 000 WT8961		AWK7G-		B36BL0(350); B36BM0(550); B36BN0(750)
PG1-AL (M) WT891 10 000 WT8951 10 000 WT8951	Methoxy-PEG-Aldehyde, mPEG-	WT8931	1 000	
B Soluble in regular aqueous solution as well as most organic solvents.	CHO, mPEG-CH ₂ -CHO	WT8921	2 000	~0< ~~~ H
Soluble in regular aqueous solution as well as most organic solvents. WT8951 WT9964 W	PG1-AL (M)	WT8911	5 000	r con T
Soluble in regular aqueous solution as well as most organic solvents. WT8951 WT8961 KV8511 WT8961 KV8511 WT8961 KV8511 WT8961 KV8511 WT8961 KV8511 WT8961 KV8511 WT8961 AV8181 AV8181 AV8181 AV8181 AV8181 AV8181 AV8181 AV8181 AV8181 BINI Soluble in regular aqueous solution as well as most organic solvents. WT8951 WT8961 AV877 AV8781 AV800 AV8781 AV8781 AV8781 AV800 AV8781 AV800 AV8781 AV8781 AV800 AV8781 AV800 A	L J	WT8941	10 000	Ö
WT8961 XV8511 X			20 000	
KV8511	well as most organic solvents.			DN Ask also proprional debyde (ALD) MW-250 20000De (ET PS/GV1) and
mPEG - Amide-Acetaldehyde Methoxy-PEG-NH-CHO, PEG-AAD PG1-ALA (M) Soluble in regular aqueous solution as well as most organic solvents. MEG-Propionaldehyde Methoxy-PEG-Propionaldehyde Mayrini Birini Soluble in regular aqueous solution as well as most organic solvents. MEGEANOUS MPEG - Thiol Methoxy-PEG-Thiol, PEG-SH PG1-SH (M) C B MP: 60-64 °C Soluble in regular aqueous solution as well as most organic solvents. MEGEANOUS MP: 60-64 °C Soluble in regular aqueous solution as well as most organic solvents. MEGEANOUS MP: 60-64 °C Soluble in regular aqueous solution as well as most organic solvents. Size: 1-1g 2-5g 3-10g MEHOXY-PEG-Hydroxyl Methoxy-PEG-Hydroxyl, PEG-OH, methyloxy-PEG, Polyethylene Glycol MEHOXY-PEG-Hydroxyl, PEG-OH, methyloxy-PEG, Polyethylene Glycol MEHOXY-PEG, Polyethylene Glycol Mayrini AYRINI AYR				butyraldehyde (ALR) but also amide-propionaldehyde urethane-propionaldehyde
mPEG - Amide-Acetaldehyde Methoxy-PEG-NH-CHO, PEG-AAD PG1-ALA (M) B N Soluble in regular aqueous solution as well as most organic solvents. mPEG-Propionaldehyde mPEG-Propionaldehyde Inquire MISCEANOUS mPEG - Thiol Methoxy-PEG-Thiol, PEG-SH PG1-SH (M) C B MP : 60-64 °C Soluble in regular aqueous solution as well as most organic solvents. Size: 1⁻¹¹g 2⁻⁵g 3⁻¹0g mPEG - Hydroxyl Methoxy-PEG-Hydroxyl Methoxy-PEG-Hydroxyl, PEG-OH, methyloxy-PEG, Polyethylene Glycol mPEG - Hydroxyl Methoxy-PEG, Polyethylene Glycol mPEG - Hydroxyl Methoxy-PEG, Polyethylene Glycol mPEG-Hydroxyl, PEG-OH, methyloxy-PEG, Polyethylene Glycol methyloxy-PEG, Polyethylene Glycol See also mPEOn-Propionaldehyde (FT-0B6601) AWKL71 350 to 1KDa AYR1K1 2 000 AYR1K1 2 000 AYR1K1 2 000 AYR1K1 3 0000 CH, SOUND AYR1M1 10 000 B36E51 750 B36E51 750 B36E71 10 000 B36E71 10		11 / 00 11	.0 000	amide-butyraldehyde, urethane-butyraldehyde (see NT-PEGYLs). E.g. mPEG
MPEG - Amide-Acetaldehyde Methoxy-PEG-NH-CHO, PEG-AAD AYR1K1 2 000 AYR1K1 5 000 AYR1K1 10 000 AYR1K1 AYR1K1 AYR1K1 AYR1K1 AYR1K1 AYR1K1 AYR1K1 AYR1K1 AYR1K1 10 000 AYR1K1				PropionAldehydes
MPEG - Amide-Acetaldehyde Methoxy-PEG-NH-CHO, PEG-AAD AYR1K1 2 000 AYR1K1 5 000 AYR1K1 10 000 AYR1K1 10 000 AYR1K1				
Methoxy-PEG-NH-CHO, PEG-AAD PGI-ALA MO PGI-ALA MO PGI-ALA MO PGI-ALA MO PGI-ALA MO AYRIMI Soluble in regular aqueous solution as well as most organic solvents. AYRIMI 10 0000 AYRIMI 10 0000 AYRIMI 10 0000 B36BP- 40 0000 B36BP- B36				See also mPEO _n -PropionAldehydes (FT- <u>0B6601</u>)
PG1-ALA (M) B				
Soluble in regular aqueous solution as well as most organic solvents.				
Soluble in regular aqueous solution as well as most organic solvents.				\
well as most organic solvents. B36BO- B36BP- 40 000 1		AYR1M1		0
MPEG-Propionaldehyde Inquire 2K, 5K, 10K, 20K(1R3030) See HomoBifunctional Aldehyde-PEG _x -Aldehyde (FT-WU0101)	1	AYR1N1	20 000	
mPEG-Propionaldehyde Inquire 2K, 5K, 10K, 20K(1R3030) • MISCEANOUS See HomoBifunctional Aldehyde-PEG₂-Aldehyde (FT-WU0101) • MISCEANOUS AWKL91 350 mPEG - Thiol AWKL91 350 B36E51 550 Methoxy-PEG-Thiol, PEG-SH PG1-SH PG1-SH PG1-SH PG1-SH PG1-SH PG1-SH PG2-SH PG1-SH PG1-S	well as most organic solvents.	B36BO-	30 000	
MISCEANOUS **mPEG - Thiol** Methoxy-PEG-Thiol, PEG-SH PG1-SH (M) (C)[[B] MP : 60-64 °C Soluble in regular aqueous solution as well as most organic solvents. Size: 1-1g 2-5g 3-10g See HomoBifunctional Aldehyde-PEGx-Aldehyde (FT-WU0101) **MKL91		B36BP-	40 000	
• MISCEANOUS mPEG - Thiol Methoxy-PEG-Thiol, PEG-SH PG1-SH PG1-SH MP: 60-64 °C Soluble in regular aqueous solution as well as most organic solvents. Size: 1-1g 2-5g 3-10g mPEG - Hydroxyl Methoxy-PEG-Hydroxyl, PEG-OH, methyloxy-PEG, Polyethylene Glycol mPEG - Hydroxyl Methoxy-PEG, Polyethylene Glycol AWKL91 B36E51 B36E51 B36E51 B36E51 B36E61 750 B36E71 B36E81 B36E8	mPEG-Propionaldehyde	Inquire		2K, 5K, 10K, 20K(1R3030)
MPEG - Thiol Methoxy-PEG-Thiol, PEG-SH B36E51 B36E61 750 B36E71 1 000 B36E81 2 000 B36E81 10 000 B36E81 40 000 B36E81 40 000 B36AA1 B36AA1 B36AA1 B36AB1 B3AB1 B3AB1 B3AB1 B3AB1 B3AB1 B3AB1 B3AB1 B3AB1 B3AB1	-			See HomoBifunctional Aldehyde-PEG _x -Aldehyde (<u>FT-WU0101</u>)
Methoxy-PEG-Thiol, PEG-SH B36E51 550 CH₃O−(CH₂CH₂O)n−CH₂CH₂−SH PG1-SH (M) B36E71 1 000 D36E81 2 000 MP: 60-64 °C B36E81 2 000 B36E91 3 400 Soluble in regular aqueous solution as well as most organic solvents. B36E81 5 000 B36EB1 10 000 B36EB1 10 000 B36ED1 30 000 B36ED1 30 000 B36EB1 40 000 B36AB1 350 B36AB1 550 Methoxy-PEG-Hydroxyl, PEG-OH, methyloxy-PEG, Polyethylene Glycol B36AB1 550 B36AC1 750	• MISCEANOUS			
B36E61 750 B36E71 1 000 B36E81 2 000 B36E81 3 400 B36E81 3 000 B36E81 3 000 B36E81 3 000 B36E81 3 000 B36E81 40 000 B36A81 B3	mPEG - Thiol			
PG1-SH (M) C [B] B36E71 1 000 MP : 60-64 °C B36E81 2 000 Soluble in regular aqueous solution as well as most organic solvents. B36E81 10 000 Size: 1 ^{-1g} 2 ^{-5g} 3 ^{-10g} B36E01 30 000 B36E1 40 000 MPEG - Hydroxyl Methoxy-PEG-Hydroxyl, PEG-OH, methyloxy-PEG, Polyethylene Glycol B36AC1 750 Methoxy-PEG, Polyethylene Glycol B36AC1 750 Methoxy-PEG, Polyethylene Glycol B36AC1 750 CH30-(CH2CH2O)H-CH2CH2-SH 1 000 B36E81 2 000 B36E81 2 000 B36E81 10 000 B36EB1 10 000 B36EB1 10 000 B36EB1 3 400 B36EB1 10 000 B36EB1 10 000 B36EB1 3 400 B36EB1 3 400	Methoxy-PEG-Thiol, PEG-SH			011 0 (011 011 0)- 011 011 or
B36E71				CH ₃ O-(CH ₂ CH ₂ O)n-CH ₂ CH ₂ -SH
Soluble in regular aqueous solution as well as most organic solvents. B36E91 B36EA1 B36EB1 B36EB1 B36EB1 B36EB1 B36EB1 B36EB1 B36EC1 B36EC1 B36ED1 B36ED1 B36ED1 B36ED1 B36ED1 B36EB1 B36ED1 B36ED1 B36ED1 B36AB1 B36BB1 B36AB1 B36BB1	[C][B]			
B36EA1 5 000 B36EB1 10 000	MP: 60-64 °C			
well as most organic solvents. B36EB1	Soluble in regular aqueous solution as			, [D]
Size: 1 ^{-1g} 2 ^{-5g} 3 ^{-10g} B36EC1 20 000 B36ED1 30 000 B36EE1 40 000 mPEG – Hydroxyl Methoxy-PEG-Hydroxyl, PEG-OH, methyloxy-PEG, Polyethylene Glycol B36A21 B36AB1 B36BB1 B36	well as most organic solvents.			+·
B36ED1 30 000 B36EE1 40 000 mPEG – Hydroxyl Methoxy-PEG-Hydroxyl, PEG-OH, methyloxy-PEG, Polyethylene Glycol B36AB1 B36B1	Ū			
B36EE1 40 000 MPEG - Hydroxyl B36A91 160 S36AA1 S50 B36AB1 S50 B36AC1 T50 S36AC1 T5	Size: 1 ^{-1g} 2 ^{-5g} 3 ^{-10g}			
mPEG – Hydroxyl Methoxy-PEG-Hydroxyl, PEG-OH, methyloxy-PEG, Polyethylene Glycol B36A91 B36A91 B36A91 B36A91 B36AA1 B36AB1 B36AB				
Methoxy-PEG-Hydroxyl, PEG-OH, methyloxy-PEG, Polyethylene Glycol B36AC1 750 O/n				
methyloxy-PEG, Polyethylene Glycol B36AB1 B36AC1 T50				O (OH OH
methyloxy-PEG, Polyethylene Glycol B36AC1 750	Methoxy-PEG-Hydroxyl, PEG-OH,			\ On .
	Monomethyl Ether	AWK7M1	1 000	Appearance: Viscous liquid or white solid depends on MW.
CAS: 25322-68-3. IO5011 2 000 Moisture Content: <0.5%.	CAS: 25322-68-3.			
PG1-OH (ZL) 105021 3 400 MW: +-15% around its mean MW.	PG1-OH (ZL)			MW: +-15% around its mean MW.
[C][B] [-] IO5031 5 000				
MP: 60-64 °C 105041 10 000				
Soluble in regular aqueous solution as 105051 20 000	Soluble in regular aqueous solution as			
well as most organic solvents. See also nydrolyzed PEGs (PEG-OH, 10-20KDa)	well as most organic solvents.			See also hydrolyzed PEGs (PEG-OH, 10-26KDa) [B36US0]
103001 30 000 [Hydrolyzed Poly(VinylAlcool)]	Ū			
Size: 1 ^{-10g} 2 ^{-50g} 3 ^{-100g} See also mPEO_n-Hydroxyl #BH9112(PEO ₁₁) to A1LYO2(PEO	Size: 1 ^{-10g} 2 ^{-50g} 3 ^{-100g}	1034/1	+0 000	See also mPEO_n-Hydroxyl #BH9112(PEO ₁₁) to A1LYO2(PEO ₄₈)
mPEG – Acrylate B36BA1 350 Da O	mPEG – Acrylate	B36BA1	350 Da	.0./ ^ \
mPEG-ACR mPEG-MAR B36BF1 550 Da O/n	•	B36BF1		$\sim \sim $
CAS:26015 72 0 B36BHI 750 Da				
DC1 A CD (ML)				0
B30BD1 2 000				
Soluble in regular aqueous solution as B36BG1 5 000				LINI
well as most organic solvents. B36BA1 10 000 1851			10 000	†[N]
B36BC1 20 000	and the second seconds.			
CV8251 30 000		CV8251	30 000	



FT- AYPMB0			
Product name	Cat.number	MW	Structure
synonymes	Qty 0-250mg,	(g/mol ⁻¹)	Structure
synonymes	1-1g	(g/IIIOI)	
mPEG – MethAcrylate	B36D91	350 Da	0
mPEG-MA, mPEG-MAR	B36DA1	550 Da	
CAS:26915-72-0	B36DC1	800 Da	\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\
PG1-MAR (ML)	B36D61	1 000	\ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \
	B36D81	2 000	I '
[ی][د،]	B36DB1	5 000	
Soluble in regular aqueous solution as	B36D51	10 000	
well as most organic solvents.	B36D71	20 000	
wen as most organic sorvents.	AWK7B1	40 000	
mPEG – NPC	Inquire	350 to 800Da	NO ₂
mPEG-Nitrophenyl Carbonate	WT9101	1 000	
PG1-MPC (ML)	DY6621	2 000	1.0.7 人 人 人 丿
[B][N]	WT9091	5 000	
	DY6631	10 000	
OH and Amine reactive	WT9111	20 000	
Soluble in regular aqueous solution as	WT9121	30 000	
well as most organic solvents.	HV4931	40 000	
mPEG-silane	GV7362	350	\
PG1-SL(L,M)	GV7372	550	_0
	GV7382	750	0/ 0 10 0 10 0 50
	OO7022	1 000	
	EV4982	2 000	\ \n\ \\n\ \\
	WT9712	3 400	0 \
	DY8032	5 000	

Store at -20°C (M). Keep desiccated. Protect from light. Stable for +12months at -20°C.

GV7352 AWKI52

B2ZUM2

B2ZUN2

Physical Properties:

• Soluble in regular aqueous solution as well as most organic solvents: water, ethanol, chloroform, DMSO, etc

10 000

20 000

30 000

40 000

Handling and Use:

For best use, material should always be kept in low temperature in dry conditions and under inert gaz for best stability. Prepare fresh solution right before use. Avoid frequent thaw and freezing.

Functional groups information (reactivity):

See more information [NT-XLreact]

Other PEG Silanes: see FT-LO5310

See also PEO₆-Silane #B63UT

The **amino** group (-NH2) reacts readily with succinimidyl ester groups (in particular NHS), carboxylic groups (via carbodiimide mediation) and many other amine reactive functional groups either in aqueous buffer or organic solvents.

The **carboxyl** group (-COHH) can be conjugated using severa conventional chemistries: via carbodiimide mediation to amines, with and many other amine reactive functional groups either in aqueous buffer or organic solvents.

The **N-hydroxysuccinimydyl** group (**NHS**) reacts at pH 7.5-9 in aqueous phase on aliphatic primary (– NH2) and secondary amines (=NH) (in fact on its deprotonated form), optimally at neutral pH or higher. A peptidic bond (amide link -CO-NH-) is formed.

NitroPhenyl Carbonate group (**NPC**) can react with both amine and hydroxyl groups. The **Maleimide** group reacts with thiols readily at pH 6.5~7.5 to form stable thioether bonds.

Hydrazine and **hydroxylamine** derivatives also have amine-like reactivity and, in some cases, can be coupled to water-soluble carbodiimide – activated carboxylic acid groups





FT- AYPMB0

The **Hydrazide** group give a variety of reactions: they react with Carbonyls (aldehydes and ketones) much faster than amine does, making them useful for site-specific crosslinking.

Reaction with carboxyls better occur with activation by a carbodiimide.

Reactions with amines are nonetheless useful, but require mediators (NH₄BO₄, via Shiff'base formation)

The **Aldehyde** group react with hydrazides and amines at pH 5-7. The reaction with hydrazides is faster than with amines, making them useful for site-specific crosslinking. It as react react spontaneously with amino groups to form Schiff base intermediates that can be stabilized by reduction with Sodium cyanoborohydride (NaCNB₃).

Related / associated products and documents

See or ask for Heterobifunctional PEG and PEO reagents; 3/4/5/6Arm-PEG reagents

See BioSciences Innovations catalogue and e-search tool.

For any information, please ask: Uptima / Interchim; Hotline: +33(0)470 0373 06

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