



TAMRA and its derivatives

Product Description

A popular orange-red fluorophore used in various bioconjugations for immunochemistry, notably with nucleic acids as well as peptides and proteins. Also an excellent fluorescence acceptor for fluorescein derivatives in FRET-based assays. Available derivatized with various chemical reactive groups.

Chemical structure of TAMRA:



(FP-AM381)

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Fluorescent properties:

 $\begin{array}{l} \mbox{Extinction Coeff. : ϵ (540) : 95 000 M^{-1} cm^{-1}$ \\ $\lambda_{exc. max.} \mbox{wem.max.}(nm): 540/565 nm (coupled)$ \\ \mbox{Note: In pH 8 buffer compared to MeOH, absorption and emission (540/565nm) are red-shifted ~8 nm with EC lowered by ~10\%. \\ \end{array}$





Typical absorption and emission spectra of TAMRA in pH7.0 buffer (coupled).

Product name	Formula	MW	Soluble in	Reactivity	Comments
cat.number	[CAS]	(g·mol ⁻¹)			
5(6)-TAMRA	C25H22N2O5.HCl	466.92	DMSO, DMF,	COOH	Mixed isomer of 5-(and 6) used to modify
(K) <u>FP-46644A</u> , 100mg			MeOH, or pH >6.5		amino and hydroxy groups using EDAC-
<u>F</u> P-JQ4610, 1mg			buffer		mediated couplings when there are
					difficulties in using 5-(and 6)-TAMRA-SE
5-TAMRA	$C_{25}H_{22}N_2O_5$	430.5	DMSO, DMF,	COOH	Single isomer used when they are
(K) FP-05867A , 10mg			MeOH, or pH >6.5		difficulties with the mixed isomer,
			buffer		typically used in protein labeling
6-TAMRA	$C_{25}H_{22}N_2O_5.$	430.5	DMSO, DMF,	COOH	Single isomer typically used in nucleotide
(K) <u>FP-M1306A</u> , 10mg	[91809-66-4]		MeOH, or pH >6.5		labeling.
			buffer		C C
5-TAMRA-PEO3-amine	$C_{37}H_{42}F_6N_4O_{11}\\$	832.74	DMSO, DMF, or	NH2	The amine derivate, with an hydrophilic
(K) <u>FP-BV3780</u> , 5mg			H ₂ O		spacer
5(6)-TAMRA-SE	$C_{29}H_{25}N_3O_7.$	527.5	DMF or DMSO	NHS/Amines	Contains an reactivity toward amines in
(K') <u>FP-52498A</u> , 25mg	[150408-83-6]				physiological conditions
5-TAMRA-SE	$C_{29}H_{25}N_3O_7$	527.5	DMF or DMSO	NHS/Amines	Is preferred for peptides and proteins.
(K') <u>FP-67480A</u> , 5mg	[150810-69-7]				



FT-52498	FT-52498					
Product name cat.number	Formula [CAS]	MW (g·mol ⁻¹)	Soluble in	Reactivity	Comments	
6- TAMRA-SE (K') <u>FP-84634A,</u> 5mg FP-CD6831, 100µl at 90mg/mL	C ₂₉ H ₂₅ N ₃ O ₇ ; [150810-69-8]	527.5	DMF or DMSO	NHS/Amines	Is preferred for labeling nucleotides and DNA sequencing.	
6-TAMRA-X-SE (K') <u>FP-33406A</u> , 10 mg		640.69	DMF or DMSO	NHS/Amines	Contains an extended spacer to reduces self quenching upon high level conjugation.	
5-TAMRA-PEO₈-SE (K') <u>FP-CD6821</u> , 1mg	$C_{48}H_{62}N_4O_{16}$	951.02	DMF or DMSO	NHS/Amines	Contains an hydrophilic spacer to improve conjugates conjugation, fluorescence and stability	
5-TAMRA-PEO ₁₂ -SE (K') <u>FP-CD6811</u> , 1mg	$C_{56}H_{78}N_4O_{20}$	1127.23	DMF or DMSO	NHS/Amines	The extreme extension version.	
TAMRA-PEG8-SE (K [°]) FP-B3A8E0, 5mg	$C_{48}H_{62}N_4O_{16}$	951.04	DMF or DMSO	NHS/Amines	CAS 2067281-44-9. $\mathbb{P}_{\mathcal{A}} = (\mathbb{P}_{\mathcal{A}} = \mathbb{P}_{\mathcal{A}} = \mathbb{P}_{$	
5(6)-TAMRA-C₅-maleimide (K') <u>FP-AM381A</u> , 5mg	$C_{34}H_{34}N_4O_6$	594.66	DMF or DMSO	Mal/Thiols	Contains an reactivity toward thiols in physiological conditions and an extended spacer	
5(6)-TAMRA-C ₆ -maleimide (K') FP-LV8390, 5mg	C ₃₅ H ₃₆ N ₄ O ₆	608.70	DMF or DMSO	Mal/Thiols	Contains an reactivity toward thiols in physiological conditions and an extended spacer	
5-TAMRA-maleimide (K') FP-21873A, 5mg	$C_{28}H_{23}N_3O_5$	481,51	DMF or DMSO	Mal/Thiols	Contains an reactivity toward thiols in physiological conditions – single isomer	
6-TAMRA-maleimide (K') FP-862231, 1mg	$C_{28}H_{23}N_3O_5$	481,51	DMF or DMSO	Mal/Thiols	Contains an reactivity toward thiols in physiological conditions – single isomer	
5(6)-TAMRA-maleimide (K') FP-IT2471, 5mg	C ₂₈ H ₂₃ N ₃ O ₅	481,51	DMF or DMSO	Mal/Thiols	Contains an reactivity toward thiols in physiological conditions	
TAMRA-MTS (K') <u>FP-60222A</u> , 5 mg	$C_{28}H_{29}N_6O_6S_2$	567.69	DMF or DMSO	MTS/Thiols	Superior specificity to maleimide.	

Storage:

Free acid: at +4°C $_{(K)}$ or -20°C for long term $_{(K)}$. Protect from light

SE, Maleimide and MTS derivatives: at -20° C (M) Protect from light and moisture. Storage can also be done at $+4^{\circ}$ C, but for short-term for SE, Maleimide, MTS derivates.

Technical and scientific information

• **Carboxy tetramethylrhodamine** (**TAMRA**) is one of the most popular orange-red fluorophore used in various bioconjugations for immunochemistry, notably with nucleic acids as well as peptides and proteins. It is an excellent fluorescence acceptor for fluorescein derivatives in FRET-based assays.

The absorption and emission spectra of TAMRA-labeled proteins may vary depending on the labeling location and on the degree of substitution. One can adress such limitations or when a fluorescence quenching is a serious problem, using alternative dye: try an extended spacer version (TAMRA-X-SE, TAMRA-PEOx-SE), or our excellent FluoProbes®547 (#FP-AK773 for the SE form) that has brighter and more stable fluorescence. It is available with various chemical reactive groups, as mixed **isomers** of carboxytetramethylrhodamine, and some derivatives as single isomer. Pure isomers are are preferred for specific biological applications where reproducibility is more critical than material cost since the minor positional difference between 5-TAMRA and 6-TAMRA might significantly affect some biological properties of the underlying conjugates.

• The **carboxyl free acid** derivate (TAMRA) can be coupled to primary or secondary amines via standard peptide chemistry, and to hydroxy groups. EDAC-mediated conjugation is useful when there are difficulties in using SE derivates.





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- The **Amine derivative** can be coupled via standard chemistry to activated carboxyl groups (i.e. via EDC mediated amidation) or sulfonyl chlorides. It is available as standard spacer, and with a long water-soluble PEG spacer separating the dye and the amine functional group (5-TAMRA-PEO₃-amine).
- The **succinimidyl ester** (SE, NHS) reacts readily with primary or secondary amines under mild conditions. It is available with a standard linker, and an improved hydrophilic and long spacer: 5-TAMRA-PEO₁₂ SE has a long water-soluble PEG spacer between tetramethylrhodamine and the amine reactive SE group. This results in improved conjugation of proteins (do not require organic solvents, allows higher ratio of coupling) and higher fluorescence properties as well conjugate stability (reduce aggregation risk).
- The **maleimide derivative** (Mal) reacts readily and specifically with sulfhydryls. It is available with a C₅ linker molecule enhances the fluorescence after the dye is conjugated to proteins.
- The MTS derivative reacts highly and specifically with sulfhydryls.

Please refer to technical notices and the scientific literature for more information on chemical reactivity and protocols, or ask FluoProbes.

References :

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TAMRA-SE :

- Fuller ME., et al., « Application of a Vital Fluorescent Staining Method for Simultaneous, Near-Real-Time Concentration Monitoring of Two Bacterial Strains in an Atlantic Coastal Plain Aquifer in Oyster, Virginia », *Appl. Envir. Microbiol.*, **70**,1680 (2004) <u>Article</u>
- Byersdorfer C.A., et al., « Visualization of Early APC/T Cell Interactions in the Mouse Lung Following Intranasal Challenge », J. Immunol., 167, 6756(2001) Article

Handling and Storage

All products should from light and moisture.

Storage of amine and carboxylic acid derivates can be stored at +4°C, or at -20°C for long term.

SE, Maleimide and MTS derivates should be stored at -20° C, or for short-term at $+4^{\circ}$ C.

Allow reaching room temperature before opening. Dissolve in suitable solvent or buffer as depicted in table page 1. Typically SE and MTS derivatives are dissolved in anhydrous DMSO, and can be stored frozen some weeks (SE) or months (MAL). Be aware that stability is reduced once in solution, because of light sensitivity, and because of hydrolysis for SE and Maleimide derivatives.





FT-52498

Name : **Catalog Number :** Structure / MW : Solubility:

Absorption / Emission :

5-(and-6)-Carboxytetramethylrhodamine, mixed isomers FP-46644A, 100mg C25H22N2O5.HC1; MW=466.92

5(6)-TAMRA

5-TAMRA

FP-05867A, 10mg

pH>6.5 buffers

6-TAMRA

FP-M1306A, 10mg

or pH>6.5 buffers

especially in solution

 $\lambda_{exc} \lambda_{em}$ (MeOH) = 540/565 nm

Dark red solid soluble in DMSO, DMF, MeOH, or pH >6.5 buffer $\lambda_{\text{exc}} \lambda_{\text{em}} (\text{MeOH}) = 540/565 \text{ nm}$ Store at $+4^{\circ}$ C or -20° C and protect from light (K)

5-Carboxytetramethylrhodamine, single isomer

C25H22N2O5.; CAS: [91809-66-4]; MW=430.5

Dark red solid soluble in DMSO, DMF, MeOH or

Store at $+4^{\circ}$ C or -20° C and protect from light (K)

6-Carboxytetramethylrhodamine, single isomer

C25H22N2O5.; CAS: [91809-66-4]; MW=430.5

Dark red solid soluble in DMSO, DMF, or MeOH,

Store at +4°C or -20°C and protect from light (K)







(CH ₃) ₂ N
C0 2
I CO ₂ H

Name :

Catalog Number : Structure : Molecular Weight : Solubility:

Absorption / Emission : Storage:

Name :

Name :

Catalog Number : Structure : Molecular Weight :

Solubility: **Absorption / Emission :** Storage:

5-TAMRA-PEO3-amine

Catalog Number : Structure : **Molecular Weight :** Solubility: **Absorption / Emission :** Storage:

Tetramethylrhodamine-PEO3-amine, TFA salt

 $\lambda_{exc} \lambda_{em}$ (MeOH) = 540/565 nm

FP-BV3780, 5mg C37H42F6N4O11; MW=832.74

Dark purple solid soluble in DMSO, DMF, or H2O $\lambda_{exc} \lambda_{em}$ (MeOH) = 544/571 nm

Store at +4°C or -20°C and protect from light (K)

5(6)-TAMRA-SE Name : 5-(and-6)-Carboxytetramethylrhodamine succinimidyl ester (mixed isomers) **Catalog Number :** FP-52498A, 25mg Structure / MW : C₂₉H₂₅N₃O₇. CAS: [150408-83-6]; MW=527.5 Solubility: Red solid soluble in DMF or DMSO **Absorption / Emission :** $\lambda_{exc} = (MeOH) = 540/565 \text{ nm}$ $\Box = 95\ 000\ M^{-1}\ cm^{-1}$ **Reactivity :** High and specific amine reactivity Storage : Store at -20°C and protect from light (M)



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FT-52498		
Name :	5-TAMRA SE	(CH ₃) ₂ N
	5-Carboxytetramethylrhodamine succinimidyl ester (single isomer)	U,
Catalog Number :	rr-0/480A, Sing	
Structure / MW :	$C_{29}H_{25}N_{3}O_{7}$. : CAS: [150810-68-7]; MW=527.5	L.
Solubility:	Dark red solid soluble in DMSO or DMF	Ŭ
Absorption / Emission :	$\lambda_{\text{exc}} \setminus \lambda_{\text{em}} \text{ (MeOH)} = 540/565 \text{ nm}$	L
Reactivity:	High and specific amine reactivity	
Storage:	Store at -20°C and protect from light (M)	
Name :	6-TAMRA SE	(CH 3)2N
	6-Carboxytetramethylrhodamine succinimidyl ester (single	
Catalog Number :	FP-84634A, 5mg	Ŷ
U U	FP-CD6831, 100µl at 90mg/mL (0.17M) in anhydrous DMSO (6-TAMRA SE in DMSO)	Ju-o-
Structure / MW :	$C_{29}H_{25}N_3O_7$; [150810-69-8]; MW=527.5	
Solubility:	Dark red solid soluble in DMSO or DMF	
Absorption / Emission :	$\lambda_{exc} \lambda_{em}$ (MeOH) = 540/565 nm	
Reactivity:	High and specific amine reactivity	
Storage	Store at -20°C and protect from light (M)	
Storage		
Name :	5-TAMRA-X SE	
	6-(Tetramethylrhodamine-5-(and-6)-	-Ń~
Catalog Number :	carboxyamido)hexanoic acid, succinimidyl ester FP-33406A, 1mg	0
Molecular Weight :	MW=640.69	И-о-т-(сн.).
Solubility:	Dark red solid soluble in DMSO or DMF	
Absorption / Emission :	$\lambda_{exc} \lambda_{em}$ (MeOH) = 540/565 nm	
Reactivity:	High and specific amine reactivity	
Storage:	Store at -20°C and protect from light (M)	
~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~		
Name :	5-TAMRA-PEO8 SE	(CH3)2N C CH3)2
	5-Carboxytetramethylrhodamine-PEO8-propionate	J ^{ro} i
Catalog Number :	FP-CD6821, 1mg	~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~~
Structure / MW ·	$C_{10}H_{c0}N_{1}O_{1}C$ MW=951.02	
Schubility	Dark red solid soluble in DMSO or DMF	
Absorption / Emission	$\lambda = (M_{PO}H) = 540/565 \text{ nm}$	
Absorption / Emission :	$\lambda_{\rm exc}$ $\lambda_{\rm em}$ (MeOII) = 540/505 mm	
Reactivity:	Stern at 20% and protect from lights of	
Storage:	Store at -20°C and protect from light(M))	
Name ·	5-TAMRA-PEO12 SE	
Traine .	5-Carboxytetramethylrhodamine-PEO12-propionate	
Catalog Number :	succinimidyl ester FP-CD6811. 1mg	
Structure / MW	$C_{ee}H_{70}N_{4}O_{20}$ · MW=1127 23	
Suluciure / IVI W :	$C_{2011/81}$ $(40_{20}, 101)$ $(1127.25)$	
Solubility:	Dark red solid solution in DIVISO of DIVIF $\lambda = \lambda^2$ (MaOII) = 540/565 mm	
Absorption / Emission :	$\Lambda_{\text{exc}} \Lambda_{\text{em}} (\text{MeOH}) = 540/505 \text{ nm}$	
	Fign and specific amine reactivity	
	solution (M)	











## FT-52498

Name :	5(6)-TAMRA-C5-maleimide
Catalog Number :	5(6)-Carboxytetramethylrhodamine-C5-maleimide FP-AM381A, 5mg
Structure / MW :	$C_{34}H_{34}N_4O_6$ ; MW=594.66
Solubility:	Red solid soluble in DMF or DMSO (>10mg/ml).
Absorption / Emission :	$\lambda_{exc} \geq 40/565 \text{ nm}$
	EC: 61 000
Reactivity:	High and specific thiol reactivity
Storage:	Store at -20°C and protect from light (M)
Name :	TAMRA-MIS Methods This Selfs and a Cash and Take Method Dhadansing
Catalog Number :	FP-60222A, 1mg
Structure / MW :	$C_{28}H_{29}N_6O_6S_2$ ; MW=567.69
Solubility:	Soluble in DMF or DMSO
Absorption / Emission :	$\lambda_{exc} \setminus \lambda_{em} (MeOH) = 540/565 \text{ nm}$
Reactivity:	High and specific thiol reactivity
Storage:	Store at -20°C and protect from light(M)





# **Ordering information**

## **Related products**

- Desalting tools: Gelfiltration columns UP84874, UptiSpin Concentrators, CelluSep dialysis tubings
- FluoProbes®547-SE (#FP-AK773), a great alternative to TAMRA that has brighter and more stable fluorescence λexc./λem.: 557/574nm

Catalog size quantities and prices may be found at <u>http://www.fluoprobes.com</u> Please inquire for higher quantities (availability, shipment conditions).

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