



DYOMICS

Colours for Life



Catalogue

Fluorescent Dyes for Bioanalytical
and Hightech Applications

7th Edition – autumn 2011

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Mission Statement

Development and growth in the use of organic dyes in biotechnology, biomedical research and clinical diagnostics is continuing. Powerful new markers and/or probes which have been constructed using knowledge of both organic chemistry and molecular biology, are combined with automated imaging workstations to define the content, activity, and dynamics of different materials in life science. These innovative applications have led to a new dye chemistry. At present, the design of functional chromophores and fluorophores is the most significant and enticing field in dye chemistry.

With more than three decades of experience in synthesis of functional fluorescent dyes, the staff of Dyomics have an excellent scientific foundation in the search for labels for bioanalytics and diagnostics. Dyomics accepts the challenge to develop new high quality "tailor made" dyes for bioanalysis and related fields of high technology. This is, however, not the only goal of our company.

As an established company, we are open-minded and flexible in nearly every aspect of colour chemistry. Our product portfolio ranges from different labels and stains for optical sensing which includes highly hydrophilic probes and dyes to new MegaStokes dyes with enhanced Stokes shift.

In cooperation with business partners in the area of molecular biology and pharmaceutical research we continue the development of new tailored products. Customers' requests for special modifications and/or synthesis are always welcome and belong to our day-to-day business.

Our greatest wish is that our products and services will satisfy you as the customer so that we may establish a long lasting relationship with each one of you.

Products

The production of new chromophores has extensively changed the possibilities for detection of biomolecules and their interaction in recent years and has enabled previously impossible insights to be made.

The advantages of optical detection methods are in the speed of the measuring process and the high sensitivity (up to single molecule detection). In principle, the collection of data over time as well as parallel or multiple readings are feasible. Hence it is possible to answer an array of questions in biotechnology or medical diagnostics in a cost-saving and reliable manner including the application of miniaturized analysers.

In this area Dyomics offers technical know how for diverse users and hardware developers in the choice and design of suitable chromophores optimized for particular excitation and emission wavelengths.

Currently, there are more than 60 fluorophores and quenchers of the DY family available as NHS- and TFP-esters, maleimides and other derivatives for the labeling of biomolecules.

In particular, our company has great experience in the development of fluorophores which can be excited with low-cost diode lasers and show emission in the red and near infrared areas of the spectrum, as well as with custom made FRET systems.

According to the demands of the customer, the activated DY dyes can be coupled covalently to amino acids, antibodies, hormones, sugars, dNTPs, oligonucleotides or other affinity tags to corresponding probes or markers.

The proprietary DY dyes are especially useful in the analysis of enzyme kinetics and DNA sequences, in the evaluation of chromosomes by Fluorescence-In-Situ-Hybridisation (FISH), in Fluorescence Correlation Spectroscopy (FCS), in FRET assays, in microscopy and in DNA and protein chips although new applications are always being discovered. Current developments, e.g., include cell organelle specific stains and probes.

Quality Statement

Dyomics is committed to a high standard of quality. Since the second year after the foundation of the company in 1999, Dyomics has worked in accordance to the quality management system DIN EN 9001:2000 (www.tuevclub.com).

The continuous monitoring and documentation of all processes, the exclusive use of certified reagents and the final control of the products using ^1H - and ^{13}C -NMR, LC/MS, HPLC, UV-VIS and fluorescence spectroscopy guarantee a high conformity and quality of every single product, a guarantee for reproducible results.

Quality management relates not only to the different products, but also to the rapid processing of both routine and R&D contracts. Characteristic therefore is the obliging and customer orientated service and the efficient implementation of synthesis. Ultimately the key to the fulfilment of our quality claims is our qualified and experienced work force.

Sincerely yours,



*Dr. Peter Czerney
(Managing Director)*

Table with spectral characteristics of DY-Labels

label	absorption maximum* [nm]	molar absorbance* [$l \cdot mol^{-1} \cdot cm^{-1}$]	emission maximum* [nm]	molecular weight# [g · mol ⁻¹]	order number	page
DY-350	353	19,000	432	777.03	350-	7
DY-405	400	32,000	423	695.59	405-	8
DY-415	418	34,000	463	476.48	415-	9
DY-490	491	73,000	515	769.66	490-	10
DY-495	494	70,000	521	525.94	495-	11
DY-505	507	80,000	528	523.97	505-	12
DY-530	533	100,000	554	695.75	530-	13
DY-547	558	150,000	573	638.74	547-	14
DY-548	558	150,000	572	754.81	548-	15
DY-549	562	150,000	577	884.91	549-	16
DY-549P1	563	150,000	578	942.99	549P1-	17
DY-550	562	120,000	577	667.76	550-	18
DY-554	544	100,000	570	580.08	554-	19
DY-555	547	100,000	573	636.19	555-	20
DY-556	548	100,000	574	715.80	556-	21
DY-560	560	120,000	578	671.84	560-	22
DY-590	581	120,000	600	719.88	590-	23
DY-591	581	120,000	598	765.91	591-	24
DY-594	594	92,000	615	937.06	594-	25
DY-605	600	110,000	624	992.18	605-	26
DY-610	610	80,000	632	667.76	610-	27
DY-615	623	200,000	643	578.73	615-	28
DY-630	638	200,000	658	634.84	630-	29
DY-631	637	200,000	657	736.88	631-	30
DY-632	636	200,000	658	852.96	632-	31
DY-633	638	200,000	658	750.91	633-	32
DY-634	636	200,000	657	969.03	634-	33
DY-635	648	200,000	670	658.86	635-	34
DY-636	647	200,000	670	760.91	636-	35
DY-647	653	250,000	673	664.78	647-	36
DY-648	655	250,000	676	780.85	648-	37
DY-649	656	250,000	670	910.95	649-	38
DY-649P1	654	250,000	672	969.03	649P1-	39
DY-650	656	220,000	676	686.92	650-	40
DY-651	655	220,000	677	788.96	651-	41
DY-652	653	220,000	676	905.03	652-	42
DY-654	653	220,000	677	1007.08	654-	43
DY-675	675	180,000	699	706.91	675-	44
DY-676	675	180,000	699	808.95	676-	45
DY-677	674	180,000	698	925.02	677-	46
DY-678	674	180,000	694	1027.07	678-	47
DY-679	679	200,000	698	1099.09	679-	48
DY-679P1	679	200,000	697	1143.14	679P1-	49
DY-680	691	140,000	709	634.84	680-	50
DY-681	692	140,000	709	736.88	681-	51
DY-682	692	140,000	709	852.96	682-	52

Table with spectral characteristics of DY-Labels

label	absorption maximum* [nm]	molar absorbance* [$l \cdot mol^{-1} \cdot cm^{-1}$]	emission maximum* [nm]	molecular weight# [g · mol ⁻¹]	order number	page
DY-700	707	140,000	728	668.86	700-	54
DY-701	709	140,000	730	770.90	701-	55
DY-703	705	140,000	721	800.93	703-	56
DY-704	706	140,000	721	902.97	704-	57
DY-730	734	240,000	755	660.88	730-	58
DY-731	736	240,000	755	762.92	731-	59
DY-732	735	240,000	756	879.00	732-	60
DY-734	733	240,000	755	995.07	734-	61
DY-749	759	240,000	780	936.99	749-	62
DY-750	751	270,000	774	712.96	750-	63
DY-751	752	270,000	772	814.99	751-	64
DY-752	750	270,000	771	931.07	752-	65
DY-754	748	270,000	771	1033.12	754-	66
DY-776	772	240,000	787	834.98	776-	67
DY-777	770	240,000	788	951.06	777-	68
DY-778	767	240,000	787	1053.11	778-	69
DY-780	783	170,000	799	660.88	780-	70
DY-781	784	170,000	796	762.92	781-	71
DY-782	785	170,000	794	879.00	782-	72
DY-800	777	280,000	791	953.08	800-	73
DY-831	844	220,000	875	788.96	831-	74
DY-480XL	504	50,000	631	514.60	480XL-	76
DY-481XL	519	50,000	648	630.67	481XL-	77
DY-485XL	488	50,000	559	502.59	485XL-	78
DY-510XL	509	50,000	590	554.67	510XL-	79
DY-520XL	522	50,000	662	514.60	520XL-	80
DY-521XL	526	50,000	666	630.67	521XL-	81
DYQ-425	425	24,500	—	784.37	DYQ-425-	86
DYQ-505	504	39,000	—	713.75	DYQ-505-	86
DYQ-1	543	48,000	—	788.77	DYQ1-	82
DYQ-2	641	90,000	—	745.81	DYQ2-	83
DYQ-660	660	140,000	—	668.86	660Q-	88
DYQ-661	662	140,000	—	770.90	661Q-	89
DYQ-3	683	80,000	—	802.86	DYQ3-	84
DYQ-700	696	48,500	—	808.92	DYQ-700-	87
DYQ-4	766	180,000	—	796.94	DYQ4-	85

Chemical structures in the supplement.

* The given values can change depending on the environment of the label (nature of conjugate and solvent).

These are the molecular weights of the carboxylic acids.

Aminoreactive Dyes

Primary amino groups are natural components of peptides and proteins which consist of amino acids. At least one amino group is positioned at the N-terminus of the peptide chain. Other biomolecules can be modified in such way that an amino group is incorporated without changing its activity or functionality. The most common examples are 5'-aminomodified DNA oligomers and cDNA bearing aminoallyl-dU-units.

There are several functional groups which react with primary amino groups forming a covalent bond between the dye and the amino containing (bio-)molecule under mild conditions not harmful to biologically active compounds. The most common reactive group being the NHS-ester (N-Hydroxysuccinimidyl-ester). At a pH between 8 and 9, the NHS-ester reacts specifically with primary amino groups. Depending on the purity (content of active NHS-ester), quantitative yields can be achieved.



Dyomics offers further aminoreactive groups too, e.g. TFP-ester (2,3,5,6-Tetrafluorophenyl-ester) or STP-ester (4-Sulfo-2,3,5,6-Tetrafluorophenyl-ester).

At Dyomics, special attention is paid to the activity of our labels. We carefully store and ship our DY-NHS-esters to supply our customers with high quality reagents. If the activity of the label suffered during shipment we replace the label for free.

Maleimides

As well as primary amino groups, thiol groups are preferred targets for fluorescent labels. There are various chemical moieties which form stable chemical bonds with thiol groups, including maleimides and iodoacetamides. The low influence of a neutral thiol-reactive label on the isoelectric point of a protein is an advantage over the labeling of amino groups in proteins and makes differential 2D-gel electrophoresis possible. Maleimides couple to thiol groups under very mild conditions at pH around 7.



Aminomodified labels

Labels with amino groups can be used for cross coupling to carboxy groups via a mild in-situ activation. Aminomodified labels can also be incorporated into a matrix during polymerization. Coupling reactions should be performed at pH 8.5 or higher since the dyes are delivered as ammonium salts.

Biotinylated labels and others

These modifications can be used to bind to streptavidin or avidin under physiological conditions, making it quite easy to detect the above proteins which are widely used in bioanalytics.

Dyomics offers a wide range of further reactive groups e.g. Hydrazide, Azide, Acetylene, Cadaverine. If you have need for a special modified label, please ask us.

DY-350

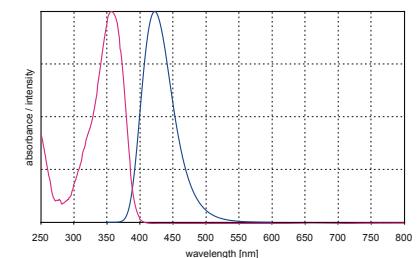
Absorption/emission max.:
353 nm / 432 nm (in PBS)

Molar absorbance:
19,000 M⁻¹cm⁻¹

Comments:
- soluble in water, methanol,
ethanol, DMF, DMSO



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	777.03	C ₁₉ H ₂₀ N ₂ O ₁₁ S ₂ * 2 C ₈ H ₂₀ N	350-00
NHS-ester	874.10	C ₂₃ H ₂₃ N ₃ O ₁₃ S ₂ * 2 C ₈ H ₂₀ N	350-01
Amino-derivative	689.86	C ₂₁ H ₂₇ N ₄ O ₁₀ S ₂ * C ₈ H ₂₀ N	350-02
Maleimide	899.15	C ₂₅ H ₂₆ N ₄ O ₁₂ S ₂ * 2 C ₈ H ₂₀ N	350-03



Structure on request

Label

Blue-green excitation

DY-405



Absorption/emission max.:
400 nm / 423 nm (in PBS)

Molar absorbance:
 $32,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in water, methanol, ethanol, DMF, DMSO

Absorption/emission max.:
418 nm / 463 nm (in Ethanol)

Molar absorbance:
 $34,000 \text{ M}^{-1}\text{cm}^{-1}$

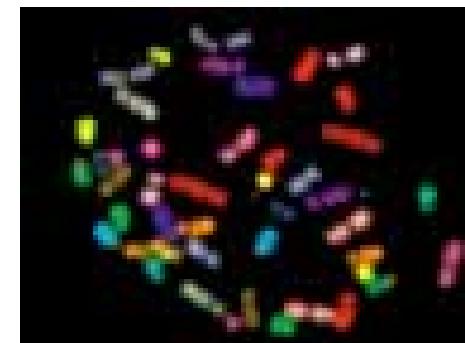
Comments:

- soluble in water, methanol, ethanol, DMF, DMSO

Label

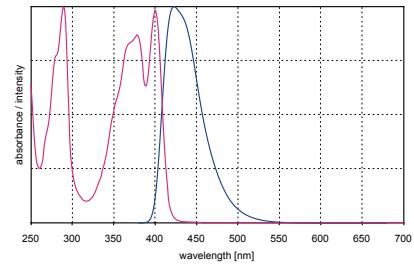
Blue-green excitation

DY-415

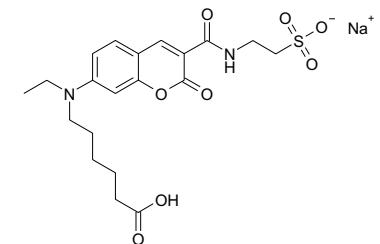
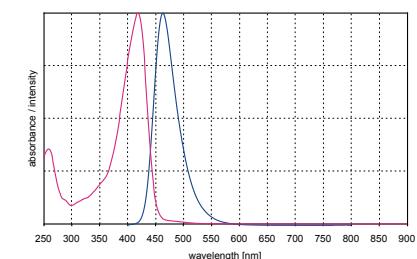


Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	695.59	C ₂₆ H ₂₀ N ₁₀ O ₁₃ S ₃ * 3 Na	405-00
NHS-ester	792.66	C ₂₆ H ₂₃ N ₂ O ₁₅ S ₃ * 3 Na	405-01
Amino-derivative	715.69	C ₂₆ H ₂₇ N ₃ O ₁₂ S ₃ * 2 Na	405-02
Maleimide	817.72	C ₃₀ H ₂₆ N ₃ O ₁₄ S ₃ * 3 Na	405-03

Structure on request



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	476.48	C ₂₀ H ₂₅ N ₂ O ₈ S * Na	415-00
NHS-ester	573.56	C ₂₄ H ₂₈ N ₃ O ₁₀ S * Na	415-01
Amino-derivative	496.59	C ₂₂ H ₃₂ N ₄ O ₇ S	415-02
Maleimide	598.61	C ₂₆ H ₃₁ N ₄ O ₉ S * Na	415-03



DY-490



Absorption/emission max.:
491 nm / 515 nm (in PBS)

Molar absorbance:
73,000 M⁻¹cm⁻¹

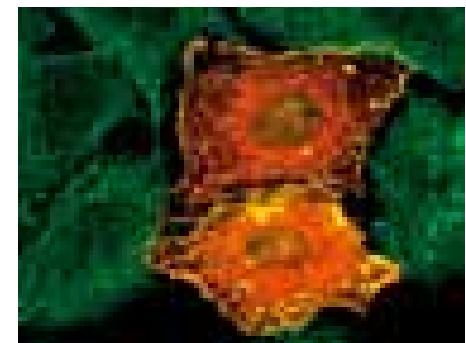
Comments:
– alternative to AlexaFluor™ 488
– soluble in water, methanol
– bright green, pH-stable emission
– suitable for flow cytometry, FISH microscopy, gel electrophoresis

Absorption/emission max.:
494 nm / 521 nm (in pH9-buffer)

Molar absorbance:
70,000 M⁻¹cm⁻¹

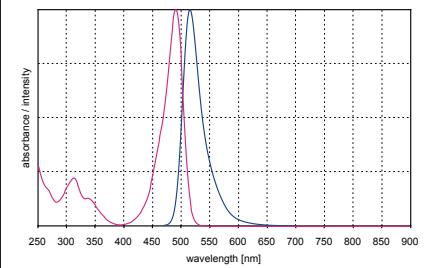
Comments:
– soluble in methanol, ethanol, DMF, DMSO
– bright solid state emission
– suitable for microarray experiments, FISH microscopy, gel electrophoresis

DY-495

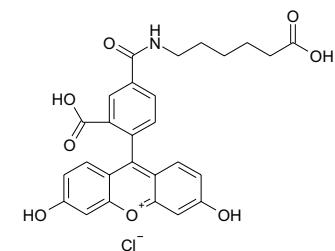
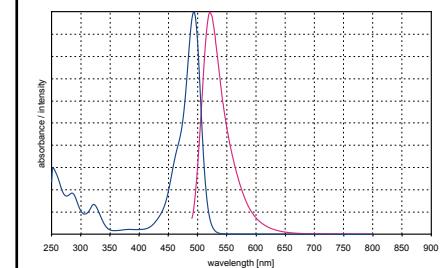


Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	769.66	C ₂₆ H ₂₂ N ₃ O ₁₃ S ₃ * CF ₃ COO	490-00
NHS-ester	1011.20	C ₂₈ H ₂₂ N ₄ O ₁₅ S ₃ * 2 C ₈ H ₂₀ N	490-01
Amino-derivative	925.78	C ₂₆ H ₂₉ N ₅ O ₁₂ S ₃ * 2 CF ₃ COO	490-02
Maleimide	799.74	C ₃₀ H ₂₆ N ₅ O ₁₄ S ₃ * Na	490-03

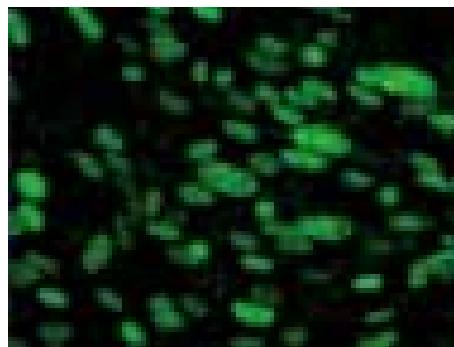
Structure on request



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	525.94	C ₂₇ H ₂₄ NO ₈ * Cl	495-00
NHS-ester	623.02	C ₃₁ H ₂₇ N ₂ O ₁₀ * Cl	495-01
Amino-derivative	604.49	C ₂₉ H ₃₁ N ₃ O ₇ * 2 Cl	495-02
Maleimide	647.07	C ₃₃ H ₃₀ N ₃ O ₉ * Cl	495-03



DY-505



Absorption/emission max.:
507 nm / 528 nm (in Ethanol)

Molar absorbance:
 $80,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments,
FISH microscopy, gel electrophoresis

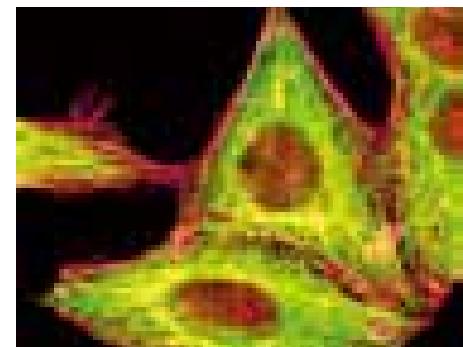
Absorption/emission max.:
533 nm / 554 nm (in PBS)

Molar absorbance:
 $100,000 \text{ M}^{-1}\text{cm}^{-1}$

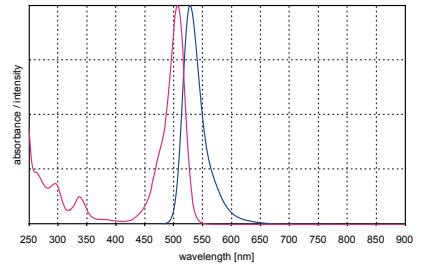
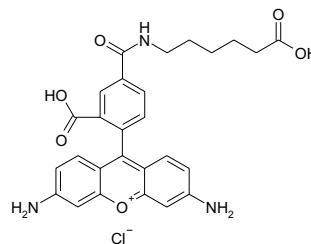
Comments:

- soluble in water, methanol

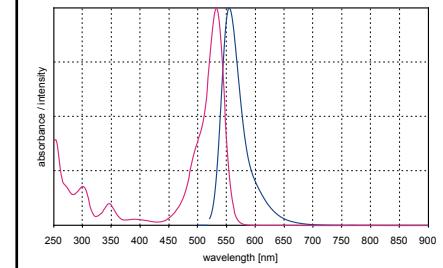
DY-530



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	523.97	C ₂₇ H ₂₆ N ₃ O ₆ * Cl	505-00
NHS-ester	621.05	C ₃₁ H ₂₉ N ₄ O ₈ * Cl	505-01
Amino-derivative	602.53	C ₂₉ H ₃₃ N ₅ O ₅ * 2 Cl	505-02
Maleimide	723.67	C ₃₃ H ₃₂ N ₅ O ₇ * CF ₃ CO ₂	505-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	695.75	C ₃₁ H ₃₄ N ₃ O ₁₀ S ₂ * Na	530-00
NHS-ester	792.82	C ₃₅ H ₃₇ N ₄ O ₁₂ S ₂ * Na	530-01
Amino-derivative	715.85	C ₃₃ H ₄₁ N ₅ O ₉ S ₂	530-02
Maleimide	817.88	C ₃₇ H ₄₀ N ₅ O ₁₁ S ₂ * Na	530-03



Structure on request

DY-547



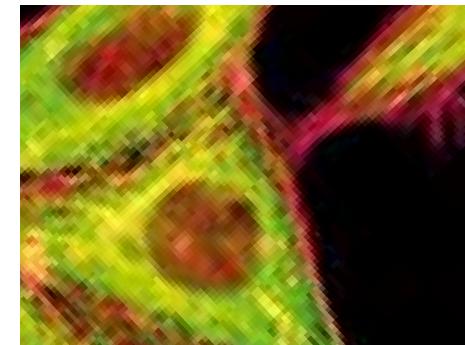
Absorption/emission max.:
558 nm / 573 nm (in Ethanol)

Molar absorbance:
 $150,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

DY-548



Absorption/emission max.:
558 nm / 572 nm (in Ethanol)

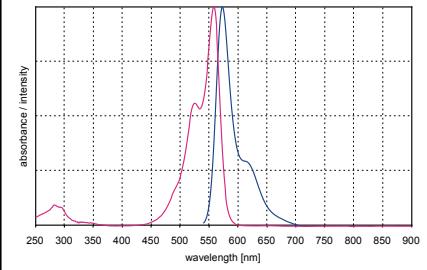
Molar absorbance:
 $150,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in water, methanol, DMF, DMSO

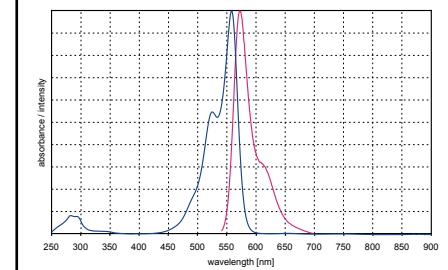
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	638.74	C ₃₀ H ₃₅ N ₂ O ₈ S ₂ * Na	547-00
NHS-ester	735.81	C ₃₄ H ₃₈ N ₃ O ₁₀ S ₂ * Na	547-01
Amino-derivative	658.84	C ₃₂ H ₄₂ N ₄ O ₇ S ₂	547-02
Maleimide	760.87	C ₃₆ H ₄₁ N ₄ O ₉ S ₂ * Na	547-03

Structure on request



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	754.81	C ₃₃ H ₃₆ N ₂ O ₁₁ S ₂ * 2 Na	548-00
NHS-ester	851.89	C ₃₅ H ₃₉ N ₃ O ₁₃ S ₃ * 2 Na	548-01
Amino-derivative	774.91	C ₃₃ H ₄₃ N ₄ O ₁₀ S ₃ * Na	548-02
Maleimide	876.94	C ₃₇ H ₄₂ N ₄ O ₁₂ S ₃ * 2 Na	548-03

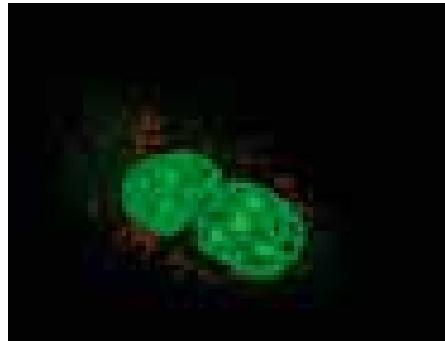
Structure on request



Label

Green excitation

DY-549



Absorption/emission max.:
562 nm / 577 nm (in Ethanol)

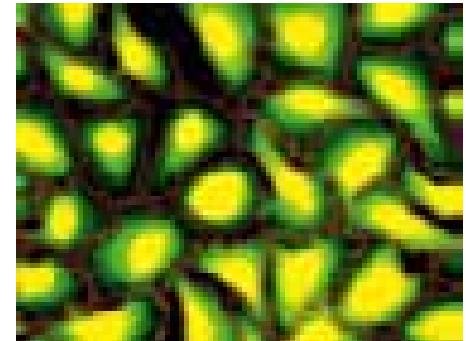
Molar absorbance:
 $150,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– soluble in water, methanol

Absorption/emission max.:
563 nm / 578 nm (in Ethanol)

Molar absorbance:
 $150,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– soluble in water, methanol

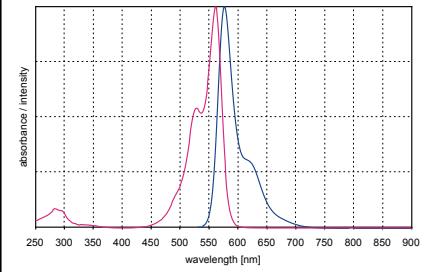


Green excitation

DY-549P1

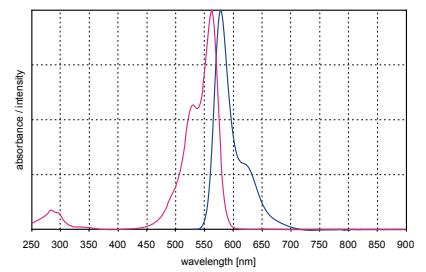
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	884.91	C ₃₃ H ₃₉ N ₂ O ₁₄ S ₄ * 3 Na	549-00
NHS-ester	981.98	C ₃₇ H ₄₂ N ₃ O ₁₆ S ₄ * 3 Na	549-01
Amino-derivative	905.01	C ₃₅ H ₄₆ N ₄ O ₁₃ S ₄ * 2 Na	549-02
Maleimide	1007.04	C ₃₉ H ₄₅ N ₄ O ₁₅ S ₄ * 3 Na	549-03

Structure on request

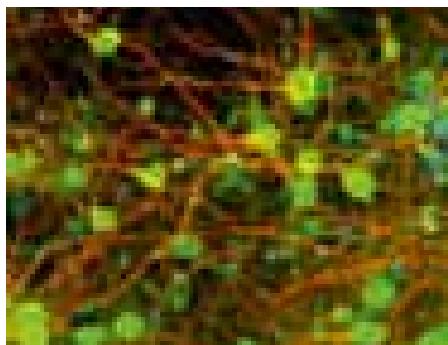


Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	942.99	C ₃₆ H ₄₅ N ₂ O ₁₅ S ₄ * 3 Na	549P1-00
NHS-ester	1040.06	C ₄₀ H ₄₈ N ₃ O ₁₇ S ₄ * 3 Na	549P1-01
Amino-derivative	963.09	C ₃₈ H ₅₂ N ₄ O ₁₄ S ₄ * 2 Na	549P1-02
Maleimide	1065.12	C ₄₂ H ₅₁ N ₄ O ₁₆ S ₄ * 3 Na	549P1-03

Structure on request



DY-550



Absorption/emission max.:
562 nm / 577 nm (in Ethanol)

Molar absorbance:
120,000 M⁻¹cm⁻¹

Comments:

- soluble in water, methanol, DMF, DMSO

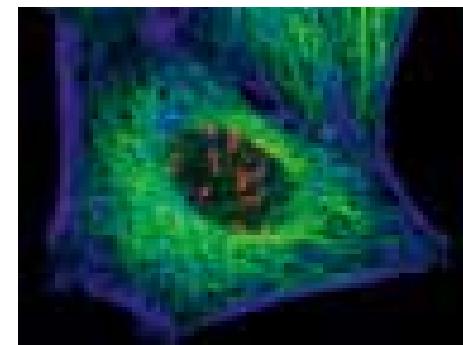
Absorption/emission max.:
544 nm / 570 nm (in Ethanol)

Molar absorbance:
100,000 M⁻¹cm⁻¹

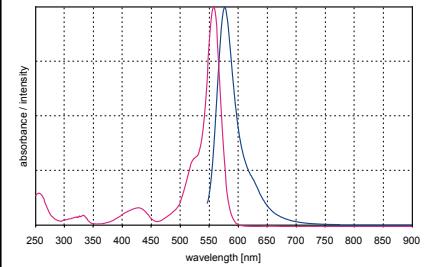
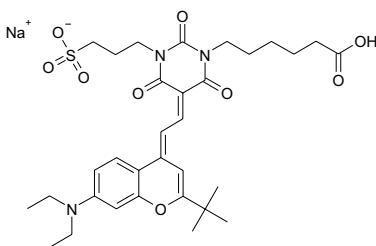
Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments,
FISH microscopy, gel electrophoresis

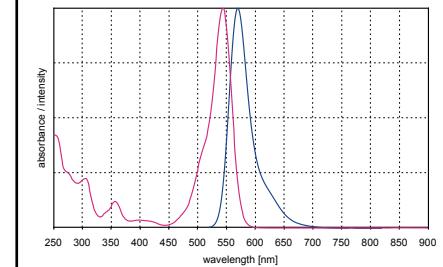
DY-554



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	667.76	C ₃₂ H ₄₂ N ₃ O ₉ S * Na	550-00
NHS-ester	764.83	C ₃₆ H ₄₅ N ₄ O ₁₁ S * Na	550-01
Amino-derivative	687.86	C ₃₄ H ₄₉ N ₅ O ₈ S	550-02
Maleimide	789.89	C ₃₈ H ₄₈ N ₅ O ₁₀ S * Na	550-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	580.08	C ₃₁ H ₃₄ N ₃ O ₆ * Cl	554-00
NHS-ester	677.16	C ₃₅ H ₃₇ N ₄ O ₈ * Cl	554-01
Amino-derivative	622.17	C ₃₃ H ₄₀ N ₅ O ₅ * Cl	554-02
Maleimide	779.78	C ₃₇ H ₄₀ N ₅ O ₇ * CF ₃ CO ₂	554-03

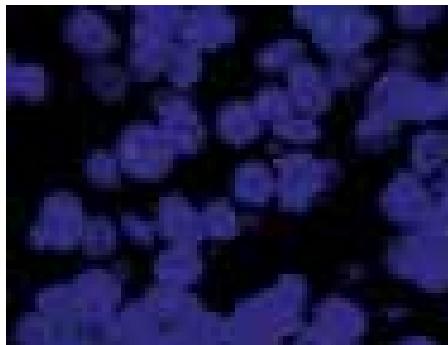


Structure on request

Label

Green excitation

DY-555



Absorption/emission max.:
547 nm / 573 nm (in Ethanol)

Molar absorbance:
 $100,000 \text{ M}^{-1}\text{cm}^{-1}$

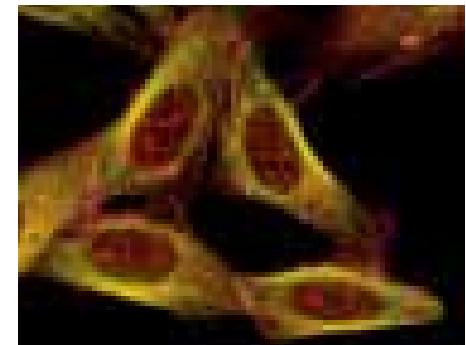
Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments, FISH microscopy, gel electrophoresis

Label

Green excitation

DY-556



Absorption/emission max.:
548 nm / 574 nm (in Ethanol)

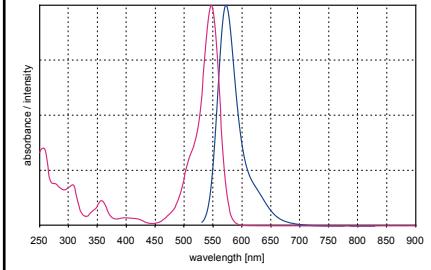
Molar absorbance:
 $100,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments, FISH microscopy, gel electrophoresis

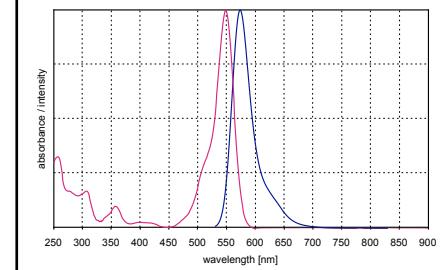
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	636.19	C ₃₅ H ₄₂ N ₃ O ₆ * Cl	555-00
NHS-ester	733.27	C ₃₉ H ₄₅ N ₄ O ₈ * Cl	555-01
Amino-derivative	678.28	C ₃₇ H ₄₈ N ₅ O ₅ * Cl	555-02
Maleimide	721.86	C ₄₁ H ₄₇ N ₅ O ₇	555-03

Structure on request



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	715.80	C ₃₆ H ₄₂ N ₃ O ₉ S * Na	556-00
NHS-ester	812.88	C ₄₀ H ₄₅ N ₄ O ₁₁ S * Na	556-01
Amino-derivative	735.91	C ₃₈ H ₄₉ N ₅ O ₈ S	556-02
Maleimide	837.93	C ₄₂ H ₄₈ N ₅ O ₁₀ S * Na	556-03

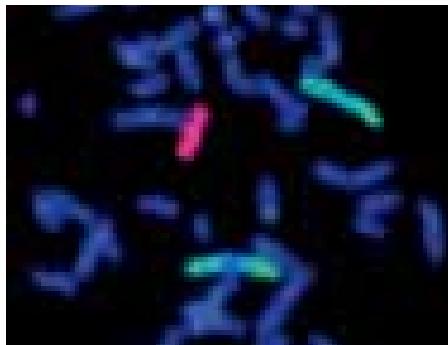
Structure on request



Label

Green excitation

DY-560



Absorption/emission max.:
560 nm / 578 nm (in Ethanol)

Molar absorbance:
 $120,000 \text{ M}^{-1}\text{cm}^{-1}$

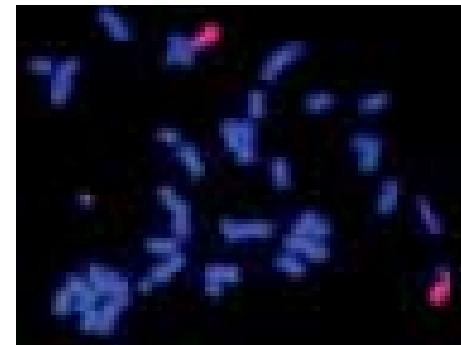
Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

Label

Orange excitation

DY-590



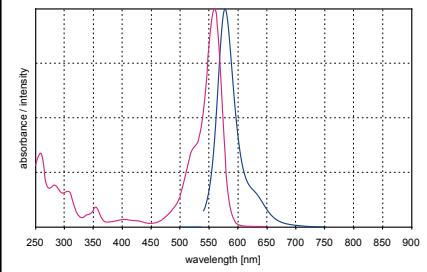
Absorption/emission max.:
581 nm / 600 nm (in Ethanol)

Molar absorbance:
 $120,000 \text{ M}^{-1}\text{cm}^{-1}$

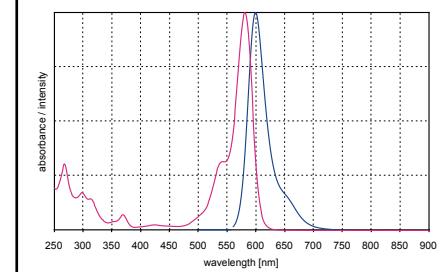
Comments:

- soluble in methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	671.84	C ₃₃ H ₄₁ N ₃ O ₈ S ₂	560-00
NHS-ester	768.91	C ₃₇ H ₄₄ N ₄ O ₁₀ S ₂	560-01
Amino-derivative	750.38	C ₃₅ H ₄₈ N ₅ O ₇ S ₂ * Cl	560-02
Maleimide	793.97	C ₃₉ H ₄₇ N ₅ O ₉ S ₂	560-03



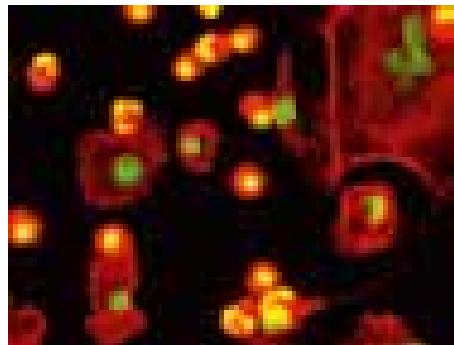
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	719.88	C ₃₇ H ₄₁ N ₃ O ₈ S ₂	590-00
NHS-ester	816.96	C ₄₁ H ₄₄ N ₄ O ₁₀ S ₂	590-01
Amino-derivative	798.42	C ₃₉ H ₄₈ N ₅ O ₇ S ₂ * Cl	590-02
Maleimide	842.01	C ₄₃ H ₄₇ N ₅ O ₉ S ₂	590-03



Label

Orange excitation

DY-591



Absorption/emission max.:
581 nm / 598 nm (in Ethanol)

Molar absorbance:
 $120,000 \text{ M}^{-1}\text{cm}^{-1}$

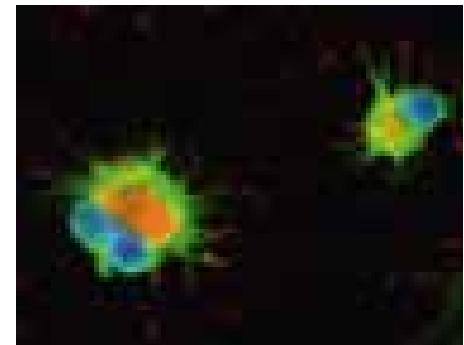
Comments:

- more hydrophilic than DY-590
- similar to Texas Red™

Label

Orange excitation

DY-594



Absorption/emission max.:
594 nm / 615 nm (in Ethanol)

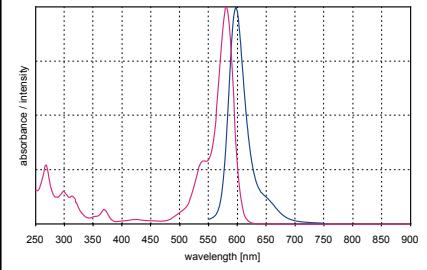
Molar absorbance:
 $92,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- similar to AlexaFluor™ 594

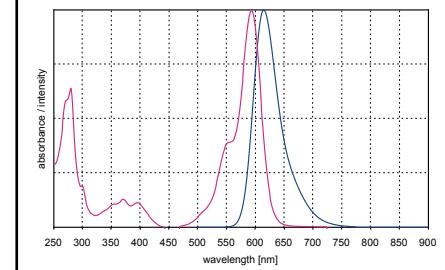
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	765.91	C ₃₈ H ₄₃ N ₃ O ₁₀ S ₂	591-00
NHS-ester	862.98	C ₄₂ H ₄₆ N ₄ O ₁₂ S ₂	591-01
Amino-derivative	844.45	C ₄₀ H ₅₀ N ₅ O ₉ S ₂ * Cl	591-02
Maleimide	888.04	C ₄₄ H ₄₉ N ₅ O ₁₁ S ₂	591-03

Structure on request



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	937.06	C ₃₉ H ₄₄ N ₄ O ₁₅ S ₄	594-00
NHS-ester	1078.10	C ₄₃ H ₄₅ N ₅ O ₁₇ S ₄ * 2 Na	594-01
Amino-derivative	979.14	C ₄₁ H ₅₀ N ₆ O ₁₄ S ₄	594-02
Maleimide	1059.19	C ₄₅ H ₅₀ N ₆ O ₁₆ S ₄	594-03

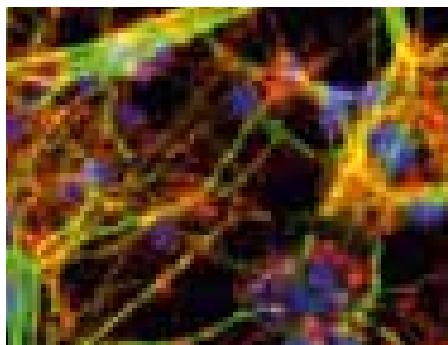
Structure on request



Label

Orange excitation

DY-605



Absorption/emission max.:
600 nm / 624 nm (in Ethanol)

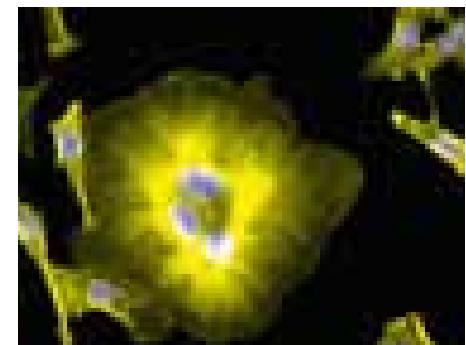
Molar absorbance:
 $110,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– hydrophilic dye

Label

Orange excitation

DY-610

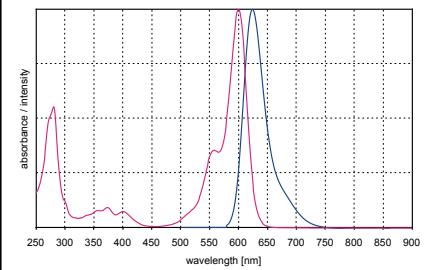


Absorption/emission max.:
610 nm / 632 nm (in Ethanol)

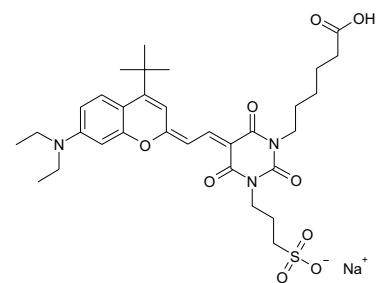
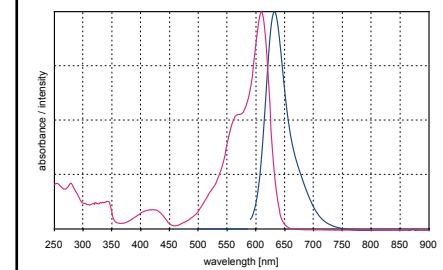
Molar absorbance:
 $80,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– soluble in water, methanol, DMF, DMSO
– bright solid state emission

Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	992.18	C ₄₃ H ₅₃ N ₅ O ₁₄ S ₄	605-00
NHS-ester	1089.26	C ₄₇ H ₅₆ N ₆ O ₁₆ S ₄	605-01
Amino-derivative	1148.30	C ₄₅ H ₆₀ N ₇ O ₁₃ S ₄ * C ₂ F ₃ O ₂	605-02
Maleimide	1114.31	C ₄₉ H ₅₉ N ₇ O ₁₅ S ₄	605-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	667.76	C ₃₂ H ₄₂ N ₃ O ₉ S * Na	610-00
NHS-ester	764.82	C ₃₆ H ₄₅ N ₄ O ₁₁ S * Na	610-01
Amino-derivative	687.86	C ₃₄ H ₄₉ N ₅ O ₈ S	610-02
Maleimide	789.89	C ₃₈ H ₄₈ N ₅ O ₁₀ S * Na	610-03



Label**Orange excitation****DY-615**

Absorption/emission max.:
623 nm / 643 nm (in Ethanol)

Molar absorbance:
 $200,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission

Absorption/emission max.:
638 nm / 658 nm (in Ethanol)

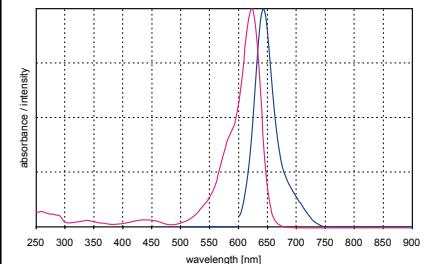
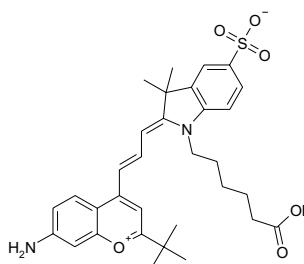
Molar absorbance:
 $200,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

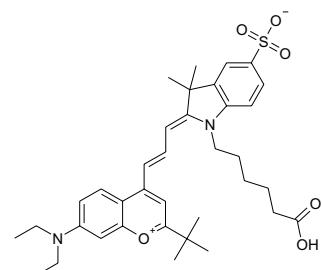
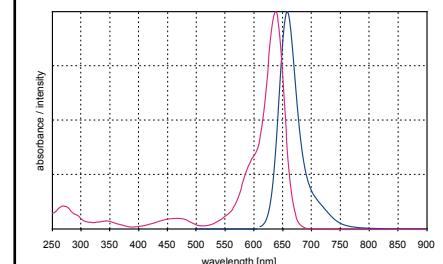
- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments,
FISH microscopy, gel electrophoresis

Label**Red excitation****DY-630**

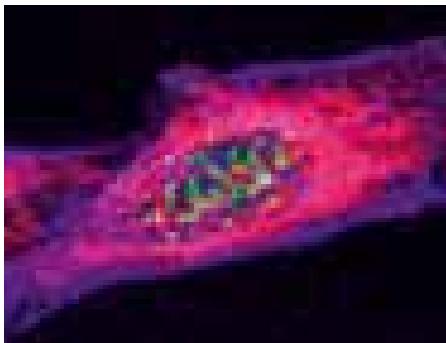
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	578.73	C ₃₂ H ₃₈ N ₂ O ₆ S	615-00
NHS-ester	675.80	C ₃₆ H ₄₁ N ₃ O ₈ S	615-01
Amino-derivative	657.28	C ₃₄ H ₄₅ N ₄ O ₅ S * Cl	615-02
Maleimide	700.86	C ₃₈ H ₄₄ N ₄ O ₇ S	615-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	634.84	C ₃₆ H ₄₆ N ₂ O ₆ S	630-00
NHS-ester	731.92	C ₄₀ H ₄₉ N ₃ O ₈ S	630-01
Amino-derivative	713.38	C ₃₈ H ₅₃ N ₄ O ₅ S * Cl	630-02
Maleimide	756.97	C ₄₂ H ₅₂ N ₄ O ₇ S	630-03



DY-631



Absorption/emission max.:
637 nm / 657 nm (in Ethanol)

Molar absorbance:
200,000 M⁻¹cm⁻¹

Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

Absorption/emission max.:
636 nm / 658 nm (in Ethanol)

Molar absorbance:
200,000 M⁻¹cm⁻¹

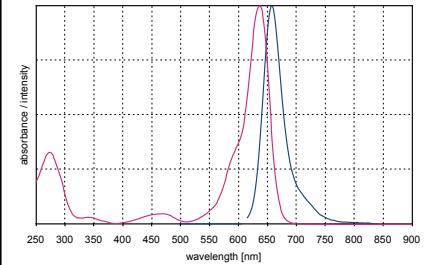
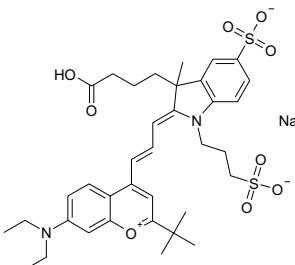
Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments, FISH microscopy, gel electrophoresis

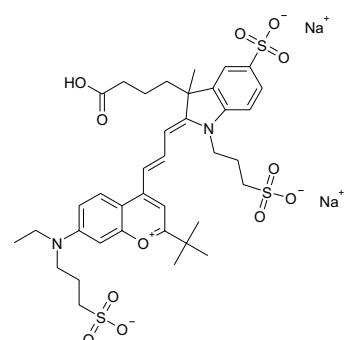
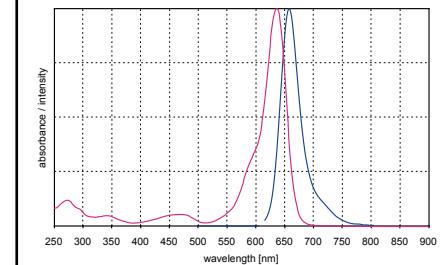
DY-632



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	736.88	C ₃₆ H ₄₅ N ₂ O ₉ S ₂ * Na	631-00
NHS-ester	833.95	C ₄₀ H ₄₈ N ₃ O ₁₁ S ₂ * Na	631-01
Amino-derivative	756.99	C ₃₈ H ₅₂ N ₄ O ₈ S ₂	631-02
Maleimide	859.01	C ₄₂ H ₅₁ N ₄ O ₁₀ S ₂ * Na	631-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	852.96	C ₃₇ H ₄₆ N ₂ O ₁₂ S ₂ * 2 Na	632-00
NHS-ester	950.03	C ₄₁ H ₄₉ N ₃ O ₁₄ S ₃ * 2 Na	632-01
Amino-derivative	873.05	C ₃₉ H ₅₃ N ₄ O ₁₁ S ₃ * Na	632-02
Maleimide	975.08	C ₄₃ H ₅₂ N ₄ O ₁₃ S ₃ * 2 Na	632-03



Label

Red excitation

DY-633



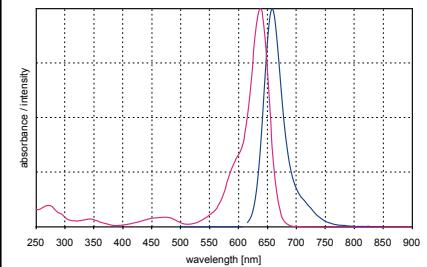
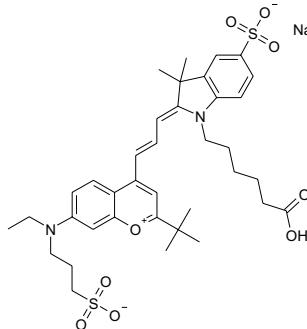
Absorption/emission max.:
638 nm / 658 nm (in Ethanol)

Molar absorbance:
200,000 M⁻¹cm⁻¹

Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	750.91	C ₃₇ H ₄₇ N ₂ O ₉ S ₂ * Na	633-00
NHS-ester	847.98	C ₄₁ H ₅₀ N ₃ O ₁₁ S ₂ * Na	633-01
Amino-derivative	771.02	C ₃₉ H ₅₄ N ₄ O ₈ S ₂	633-02
Maleimide	873.04	C ₄₃ H ₅₃ N ₄ O ₁₀ S ₂ * Na	633-03

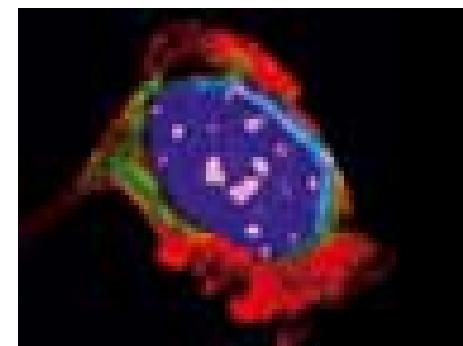


Absorption/emission max.:
636 nm / 657 nm (in Ethanol)

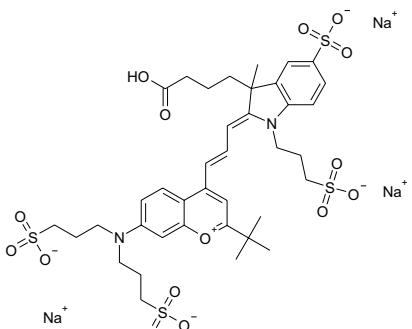
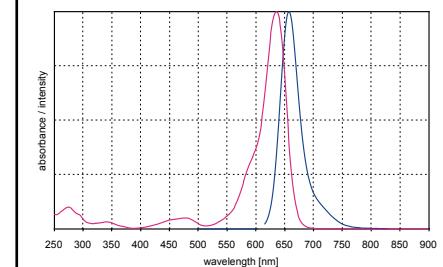
Molar absorbance:
200,000 M⁻¹cm⁻¹

Comments:

- soluble in water, methanol, DMF, DMSO



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	969.03	C ₃₈ H ₄₇ N ₂ O ₁₅ S ₂ * 3 Na	634-00
NHS-ester	1066.10	C ₄₂ H ₅₀ N ₃ O ₁₇ S ₄ * 3 Na	634-01
Amino-derivative	989.12	C ₄₀ H ₅₄ N ₄ O ₁₄ S ₄ * 2 Na	634-02
Maleimide	1091.16	C ₄₄ H ₅₃ N ₄ O ₁₆ S ₄ * 3 Na	634-03

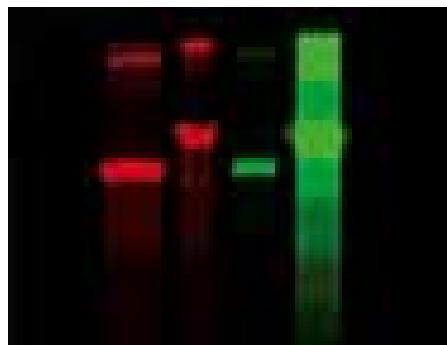


Label

Red excitation

DY-634

DY-635



Absorption/emission max.:
648 nm / 670 nm (in Ethanol)

Molar absorbance:
 $200,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments, FISH microscopy, gel electrophoresis

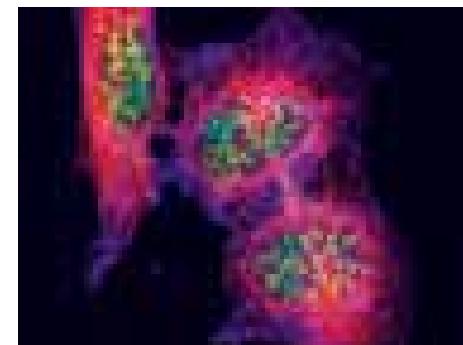
Absorption/emission max.:
647 nm / 670 nm (in Ethanol)

Molar absorbance:
 $200,000 \text{ M}^{-1}\text{cm}^{-1}$

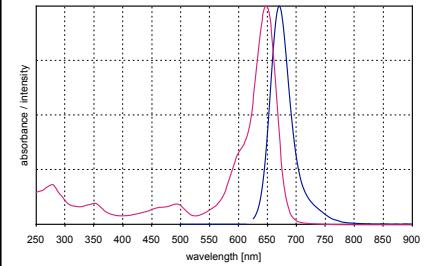
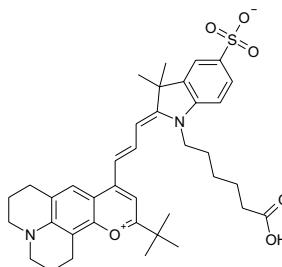
Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

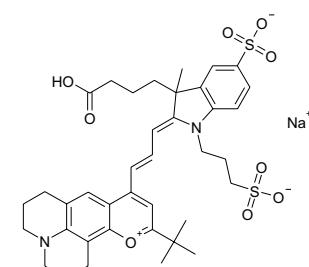
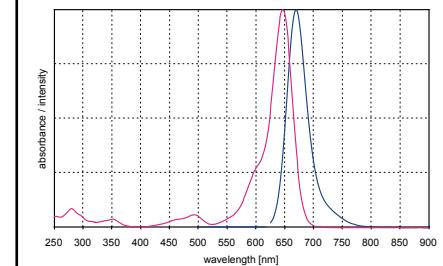
DY-636



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	658.86	C ₃₈ H ₄₆ N ₂ O ₆ S	635-00
NHS-ester	755.93	C ₄₂ H ₄₉ N ₃ O ₈ S	635-01
Amino-derivative	737.41	C ₄₀ H ₅₃ N ₄ O ₅ S * Cl	635-02
Maleimide	780.99	C ₄₄ H ₅₂ N ₄ O ₇ S	635-03



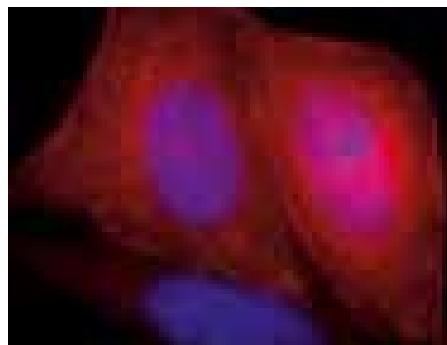
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	760.91	C ₃₈ H ₄₅ N ₂ O ₉ S ₂ * Na	636-00
NHS-ester	857.98	C ₄₂ H ₄₈ N ₃ O ₁₁ S ₂ * Na	636-01
Amino-derivative	781.01	C ₄₀ H ₅₂ N ₄ O ₈ S ₂	636-02
Maleimide	883.04	C ₄₄ H ₅₁ N ₄ O ₁₀ S ₂ * Na	636-03



Label

Red excitation

DY-647



Absorption/emission max.:
653 nm / 673 nm (in Ethanol)

Molar absorbance:
 $250,000 \text{ M}^{-1}\text{cm}^{-1}$

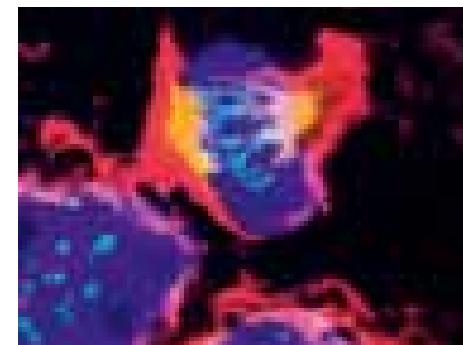
Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

Label

Red excitation

DY-648



Absorption/emission max.:
655 nm / 676 nm (in Ethanol)

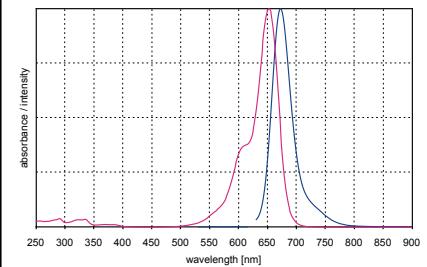
Molar absorbance:
 $250,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in water, methanol, DMF, DMSO

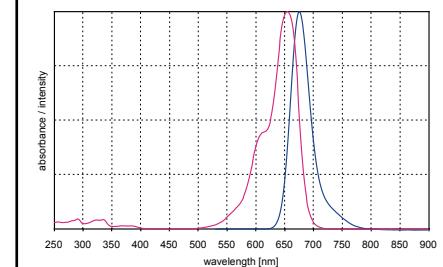
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	664.78	C ₃₂ H ₃₇ N ₂ O ₈ S ₂ * Na	647-00
NHS-ester	761.85	C ₃₆ H ₄₀ N ₃ O ₁₀ S ₂ * Na	647-01
Amino-derivative	684.88	C ₃₄ H ₄₄ N ₄ O ₇ S ₂	647-02
Maleimide	786.91	C ₃₈ H ₄₃ N ₄ O ₉ S ₂ * Na	647-03

Structure on request



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	780.85	C ₃₃ H ₃₈ N ₂ O ₁₁ S ₂ * 2 Na	648-00
NHS-ester	877.92	C ₃₇ H ₄₁ N ₃ O ₁₃ S ₃ * 2 Na	648-01
Amino-derivative	800.95	C ₃₅ H ₄₅ N ₄ O ₁₀ S ₃ * Na	648-02
Maleimide	902.98	C ₃₉ H ₄₄ N ₄ O ₁₂ S ₃ * 2 Na	648-03

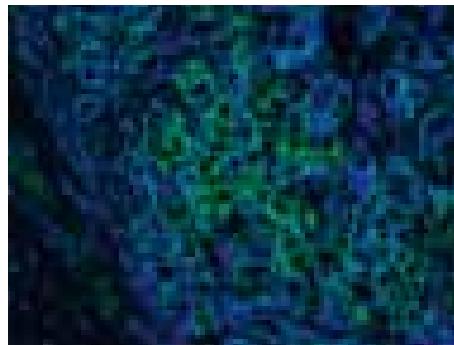
Structure on request



Label

Red excitation

DY-649



Absorption/emission max.:
656 nm / 670 nm (in Ethanol)

Molar absorbance:
 $250,000 \text{ M}^{-1}\text{cm}^{-1}$

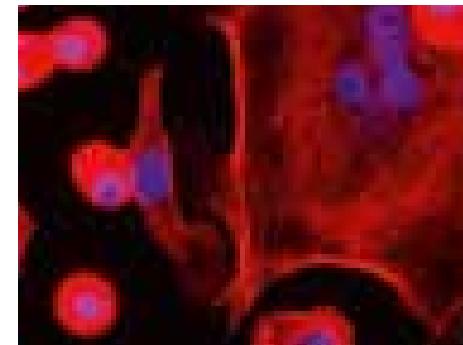
Comments:

- soluble in water, methanol
- very hydrophilic
- three negative charges
- usually combined with DY-549

Label

Red excitation

DY-649P1



Absorption/emission max.:
654 nm / 672 nm (in Ethanol)

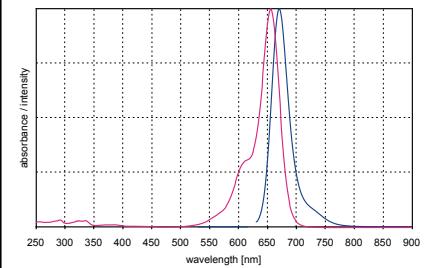
Molar absorbance:
 $250,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- replacement / alternative to DY-649

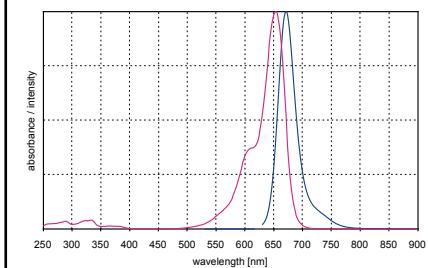
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	910.95	C ₃₅ H ₄₁ N ₂ O ₁₄ S ₄ * 3 Na	649-00
NHS-ester	1008.02	C ₃₉ H ₄₄ N ₃ O ₁₆ S ₄ * 3 Na	649-01
Amino-derivative	931.05	C ₃₇ H ₄₈ N ₄ O ₁₃ S ₄ * 2 Na	649-02
Maleimide	1033.08	C ₄₁ H ₄₇ N ₄ O ₁₅ S ₄ * 3 Na	649-03

Structure on request

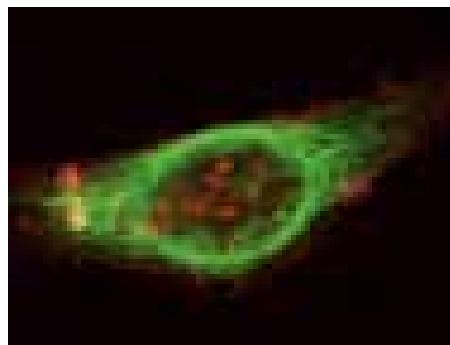


Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	969.03	C ₃₈ H ₄₇ N ₂ O ₁₅ S ₄ * 3 Na	649P1-00
NHS-ester	1066.10	C ₄₂ H ₅₀ N ₃ O ₁₇ S ₄ * 3 Na	649P1-01
Amino-derivative	989.13	C ₄₀ H ₅₄ N ₄ O ₁₄ S ₄ * 2 Na	649P1-02
Maleimide	1091.16	C ₄₄ H ₅₃ N ₄ O ₁₆ S ₄ * 3 Na	649P1-03

Structure on request



DY-650



Absorption/emission max.:
656 nm / 676 nm (in Ethanol)

Molar absorbance:
 $220,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments, FISH microscopy, gel electrophoresis

Absorption/emission max.:
655 nm / 677 nm (in Ethanol)

Molar absorbance:
 $220,000 \text{ M}^{-1}\text{cm}^{-1}$

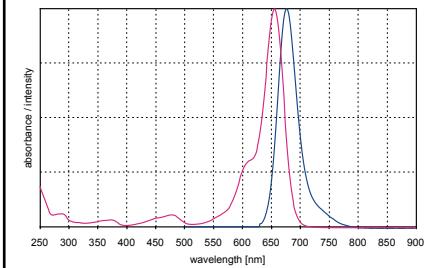
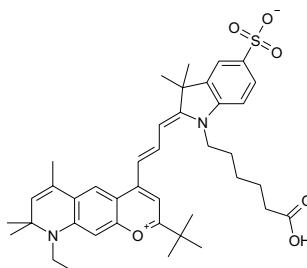
Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

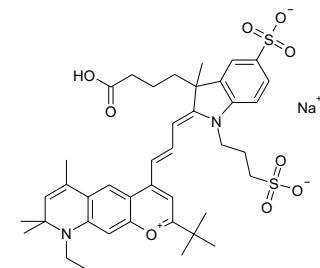
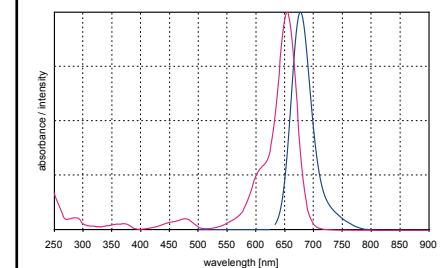
DY-651

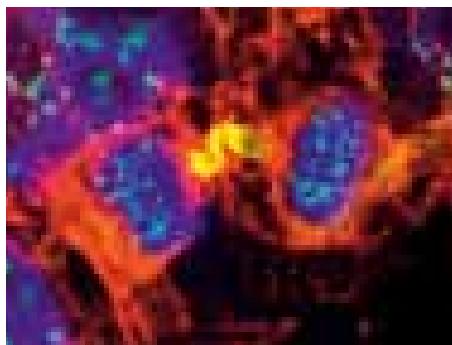


Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	686.92	C ₄₀ H ₅₀ N ₂ O ₆ S	650-00
NHS-ester	783.99	C ₄₄ H ₅₃ N ₃ O ₈ S	650-01
Amino-derivative	765.46	C ₄₂ H ₅₇ N ₄ O ₅ S * Cl	650-02
Maleimide	809.05	C ₄₆ H ₅₆ N ₄ O ₇ S	650-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	788.96	C ₄₀ H ₄₉ N ₂ O ₉ S ₂ * Na	651-00
NHS-ester	886.04	C ₄₄ H ₅₂ N ₃ O ₁₁ S ₂ * Na	651-01
Amino-derivative	809.06	C ₄₂ H ₅₆ N ₄ O ₈ S ₂	651-02
Maleimide	911.09	C ₄₆ H ₅₅ N ₄ O ₁₀ S ₂ * Na	651-03



Label**Red excitation****DY-652**

Absorption/emission max.:
653 nm / 676 nm (in Ethanol)

Molar absorbance:
 $220,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

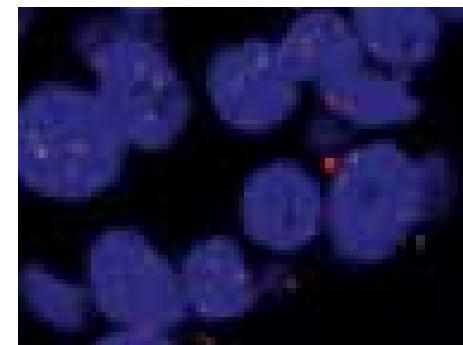
- soluble in methanol, ethanol, DMF, DMSO

Absorption/emission max.:
653 nm / 677 nm (in Ethanol)

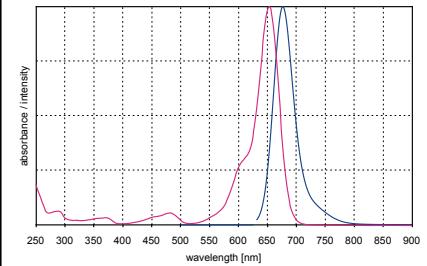
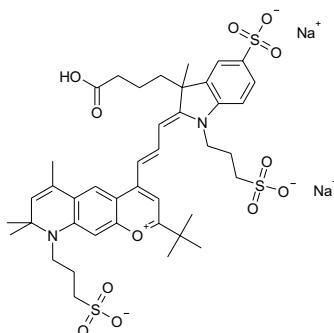
Molar absorbance:
 $220,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

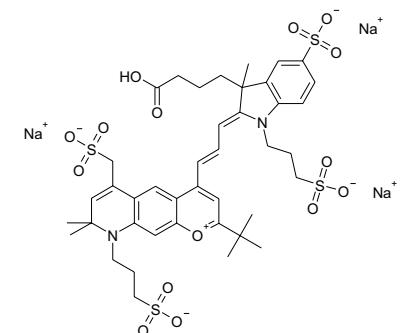
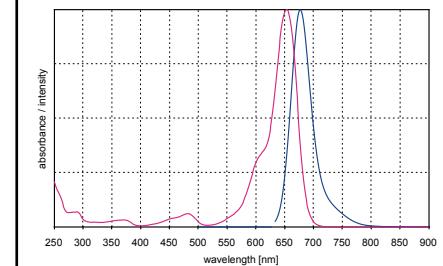
- soluble in water, methanol, DMF

Label**Red excitation****DY-654**

Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	905.03	C ₄₃ H ₅₀ N ₂ O ₁₂ S ₃ * 2 Na	652-00
NHS-ester	1002.11	C ₄₅ H ₅₃ N ₃ O ₁₄ S ₃ * 2 Na	652-01
Amino-derivative	925.13	C ₄₃ H ₅₇ N ₄ O ₁₁ S ₃ * Na	652-02
Maleimide	1027.16	C ₄₇ H ₅₆ N ₄ O ₁₃ S ₃ * 2 Na	652-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	1007.08	C ₄₃ H ₄₉ N ₂ O ₁₅ S ₄ * 3 Na	654-00
NHS-ester	1102.14	C ₄₅ H ₅₀ N ₃ O ₁₇ S ₄ * 3 Na	654-01
Amino-derivative	1027.18	C ₄₃ H ₅₆ N ₄ O ₁₄ S ₄ * 2 Na	654-02
Maleimide	1129.21	C ₄₇ H ₅₅ N ₄ O ₁₆ S ₄ * 3 Na	654-03



Label

Far red excitation

DY-675



Absorption/emission max.:
675 nm / 699 nm (in Ethanol)

Molar absorbance:
 $180,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments, FISH microscopy, gel electrophoresis

Absorption/emission max.:
675 nm / 699 nm (in Ethanol)

Molar absorbance:
 $180,000 \text{ M}^{-1}\text{cm}^{-1}$

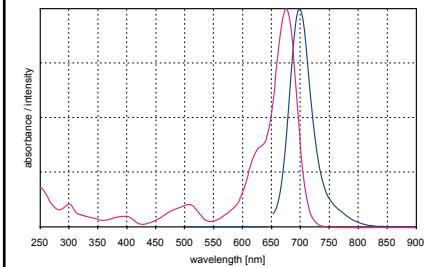
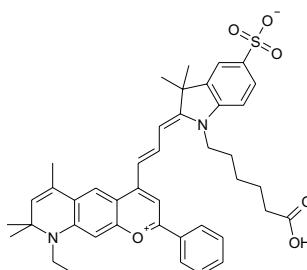
Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

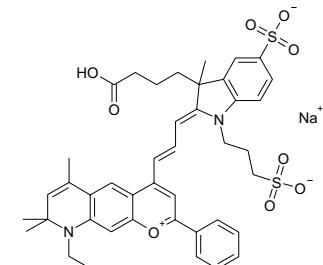
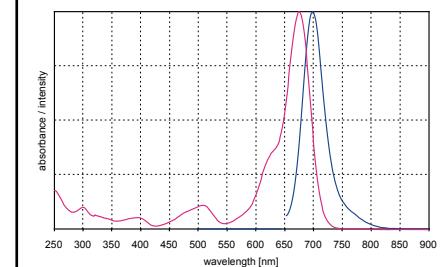


DY-676

Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	706.91	C ₄₂ H ₄₆ N ₂ O ₆ S	675-00
NHS-ester	803.98	C ₄₆ H ₄₉ N ₃ O ₈ S	675-01
Amino-derivative	785.45	C ₄₄ H ₅₃ N ₄ O ₅ S * Cl	675-02
Maleimide	829.04	C ₄₈ H ₅₂ N ₄ O ₇ S	675-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	808.95	C ₄₂ H ₄₅ N ₂ O ₉ S ₂ * Na	676-00
NHS-ester	906.02	C ₄₆ H ₄₈ N ₃ O ₁₁ S ₂ * Na	676-01
Amino-derivative	829.06	C ₄₄ H ₅₂ N ₄ O ₈ S ₂	676-02
Maleimide	931.08	C ₄₈ H ₅₁ N ₄ O ₁₀ S ₂ * Na	676-03



Label**Far red excitation****DY-677**

Absorption/emission max.:
674 nm / 698 nm (in Ethanol)

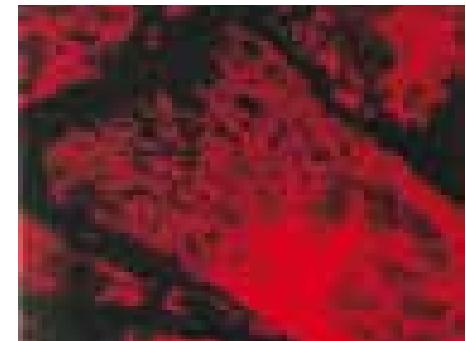
Molar absorbance:
 $180,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– soluble in methanol, DMF, DMSO

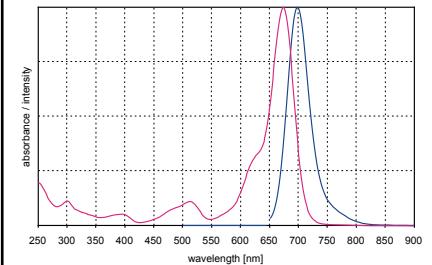
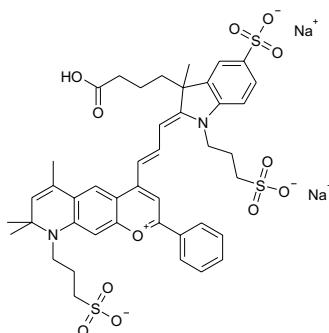
Absorption/emission max.:
674 nm / 694 nm (in Ethanol)

Molar absorbance:
 $180,000 \text{ M}^{-1}\text{cm}^{-1}$

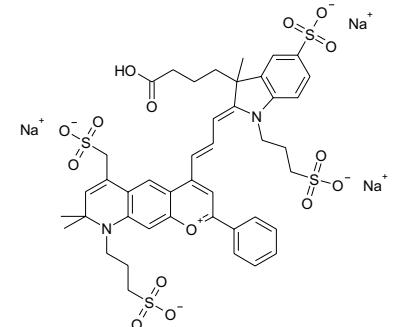
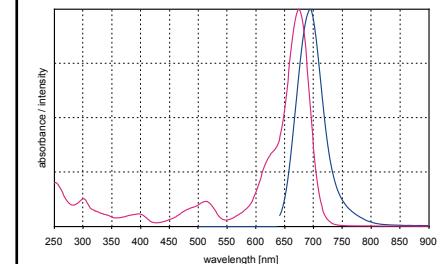
Comments:
– very hydrophilic

Label**Far red excitation****DY-678**

Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	925.02	C ₄₃ H ₄₆ N ₂ O ₁₂ S ₃ * 2 Na	677-00
NHS-ester	1022.10	C ₄₇ H ₄₉ N ₃ O ₁₄ S ₃ * 2 Na	677-01
Amino-derivative	945.12	C ₄₅ H ₅₁ N ₄ O ₁₁ S ₃ * Na	677-02
Maleimide	1047.15	C ₄₉ H ₅₂ N ₄ O ₁₃ S ₃ * 2 Na	677-03



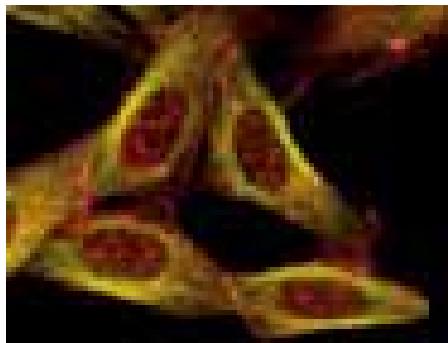
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	1027.07	C ₄₃ H ₄₄ N ₂ O ₁₅ S ₄ * 3 Na	678-00
NHS-ester	1124.14	C ₄₇ H ₄₈ N ₃ O ₁₇ S ₄ * 3 Na	678-01
Amino-derivative	1047.17	C ₄₅ H ₅₁ N ₄ O ₁₄ S ₄ * 2 Na	678-02
Maleimide	1149.30	C ₄₉ H ₅₀ N ₄ O ₁₆ S ₄ * 3 Na	678-03



Label

Far red excitation

DY-679



Absorption/emission max.:
679 nm / 698 nm (in PBS)

Molar absorbance:
 $200,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– soluble in water, methanol,
DMF-water-mixtures

Absorption/emission max.:
679 nm / 697 nm (in PBS)

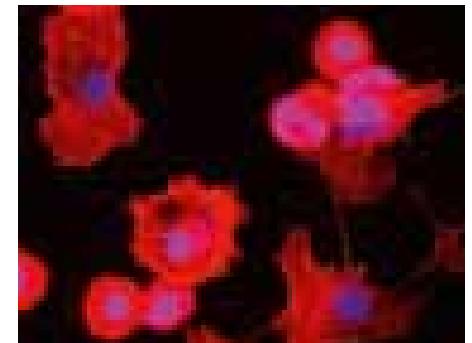
Molar absorbance:
 $200,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– alternative to DY-679

Label

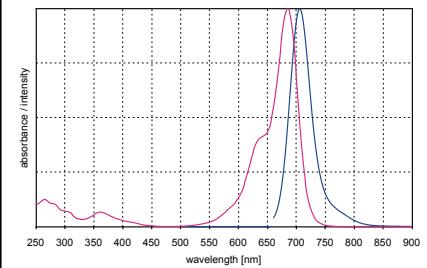
Far red excitation

DY-679P1



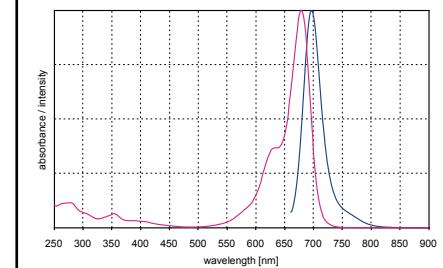
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	1099.09	C ₄₂ H ₄₂ N ₂ O ₁₇ S ₅ * 4 Na	679-00
NHS-ester	1196.16	C ₄₆ H ₄₅ N ₃ O ₁₉ S ₅ * 4 Na	679-01
Amino-derivative	1119.19	C ₄₄ H ₄₉ N ₄ O ₁₆ S ₅ * 3 Na	679-02
Maleimide	1221.21	C ₄₈ H ₄₈ N ₄ O ₁₈ S ₅ * 4 Na	679-03

Structure on request



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	1143.14	C ₄₄ H ₄₆ N ₂ O ₁₈ S ₅ * 4 Na	679P1-00
NHS-ester	1240.21	C ₄₈ H ₄₉ N ₃ O ₂₀ S ₅ * 4 Na	679P1-01
Amino-derivative	1163.24	C ₄₆ H ₅₃ N ₄ O ₁₇ S ₅ * 3 Na	679P1-02
Maleimide	1265.27	C ₅₀ H ₅₂ N ₄ O ₁₉ S ₅ * 4 Na	679P1-03

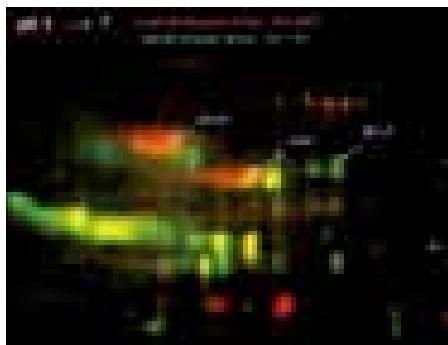
Structure on request



Label

Far red excitation

DY-680



Absorption/emission max.:
691 nm / 709 nm (in Ethanol)

Molar absorbance:
 $140,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments, FISH microscopy, gel electrophoresis



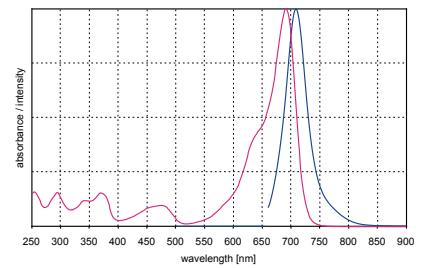
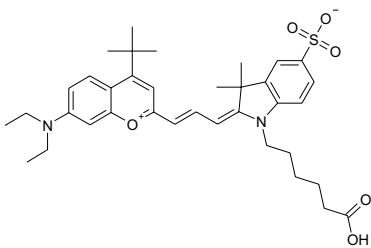
Absorption/emission max.:
692 nm / 709 nm (in Ethanol)

Molar absorbance:
 $140,000 \text{ M}^{-1}\text{cm}^{-1}$

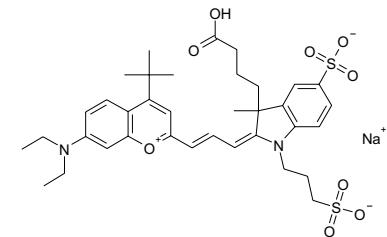
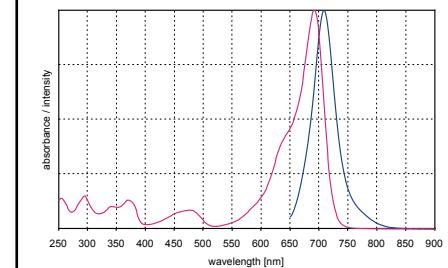
Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	634.84	C ₃₆ H ₄₆ N ₂ O ₆ S	680-00
NHS-ester	731.92	C ₄₀ H ₄₉ N ₃ O ₈ S	680-01
Amino-derivative	713.38	C ₃₈ H ₅₃ N ₄ O ₅ S * Cl	680-02
Maleimide	756.97	C ₄₂ H ₅₂ N ₄ O ₇ S	680-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	736.88	C ₃₆ H ₄₅ N ₂ O ₉ S ₂ * Na	681-00
NHS-ester	833.95	C ₄₀ H ₄₈ N ₃ O ₁₁ S ₂ * Na	681-01
Amino-derivative	756.99	C ₃₈ H ₅₂ N ₄ O ₈ S ₂	681-02
Maleimide	859.01	C ₄₂ H ₅₁ N ₄ O ₁₀ S ₂ * Na	681-03



Label

Far red excitation

DY-681

Far red excitation

DY-682

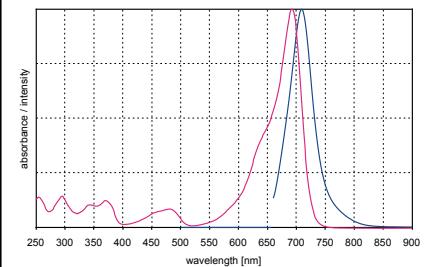
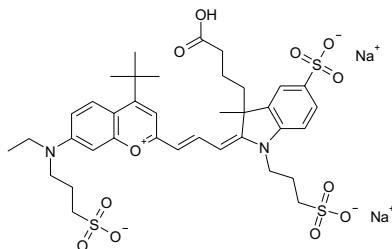


Absorption/emission max.:
692 nm / 709 nm (in Ethanol)

Molar absorbance:
 $140,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– soluble in water, methanol, DMF, DMSO

Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	852.96	C ₃₇ H ₄₆ N ₂ O ₁₂ S ₃ * 2 Na	682-00
NHS-ester	950.03	C ₄₁ H ₄₉ N ₃ O ₁₄ S ₂ * 2 Na	682-01
Amino-derivative	873.06	C ₃₉ H ₅₃ N ₄ O ₁₁ S ₃ * Na	682-02
Maleimide	975.08	C ₄₃ H ₅₂ N ₄ O ₁₃ S ₃ * 2 Na	682-03



DY-700



Absorption/emission max.:
707 nm / 728 nm (in Ethanol)

Molar absorbance:
 $140,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments, FISH microscopy, gel electrophoresis

Absorption/emission max.:
709 nm / 730 nm (in Ethanol)

Molar absorbance:
 $140,000 \text{ M}^{-1}\text{cm}^{-1}$

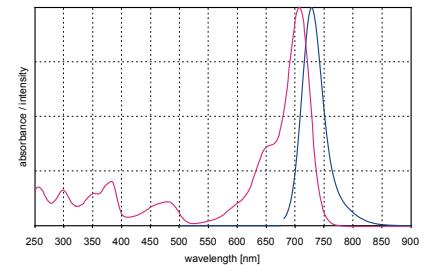
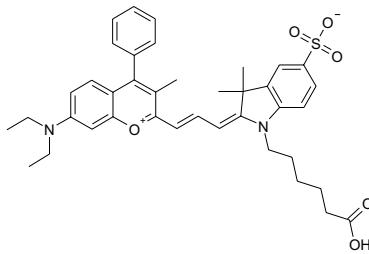
Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

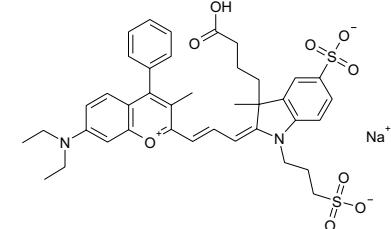
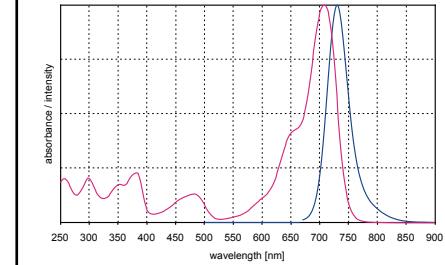
DY-701



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	668.86	C ₃₉ H ₄₄ N ₂ O ₆ S	700-00
NHS-ester	765.93	C ₄₃ H ₄₇ N ₃ O ₈ S	700-01
Amino-derivative	747.40	C ₄₁ H ₅₁ N ₄ O ₅ S * Cl	700-02
Maleimide	790.99	C ₄₅ H ₅₀ N ₄ O ₇ S	700-03



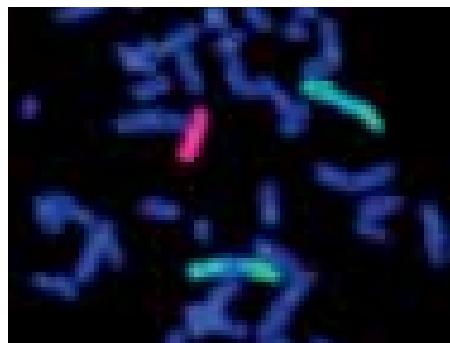
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	770.90	C ₃₉ H ₄₃ N ₂ O ₉ S ₂ * Na	701-00
NHS-ester	867.97	C ₄₃ H ₄₆ N ₃ O ₁₁ S ₂ * Na	701-01
Amino-derivative	791.01	C ₄₁ H ₅₀ N ₄ O ₈ S ₂	701-02
Maleimide	893.03	C ₄₅ H ₄₉ N ₄ O ₁₀ S ₂ * Na	701-03



Label

NIR-excitation

DY-703



Absorption/emission max.:
705 nm / 721 nm (in Ethanol)

Molar absorbance:
 $140,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– soluble in water, methanol,
ethanol, DMF, DMSO

Label

NIR-excitation

DY-704

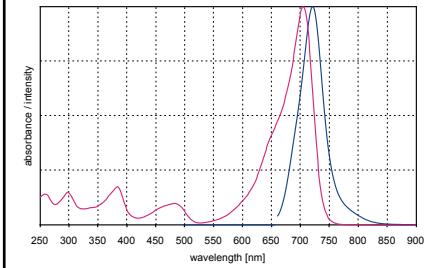
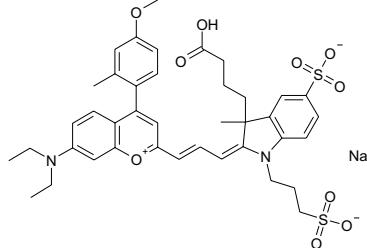


Absorption/emission max.:
706 nm / 721 nm (in Ethanol)

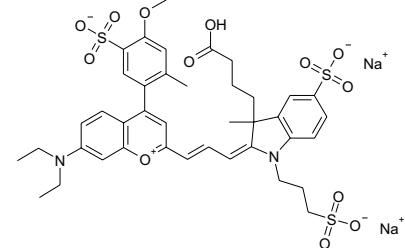
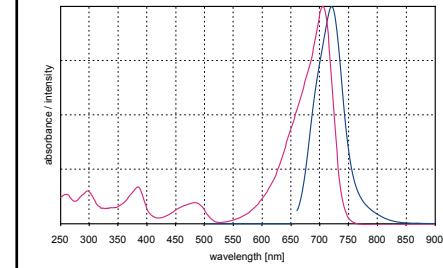
Molar absorbance:
 $140,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– very hydrophilic

Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	800.93	C ₄₀ H ₄₅ N ₂ O ₁₀ S ₂ * Na	703-00
NHS-ester	898.00	C ₄₄ H ₄₈ N ₃ O ₁₂ S ₂ * Na	703-01
Amino-derivative	821.03	C ₄₂ H ₅₂ N ₄ O ₉ S ₂	703-02
Maleimide	923.06	C ₄₆ H ₅₁ N ₄ O ₁₁ S ₂ * Na	703-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	902.97	C ₄₀ H ₄₄ N ₂ O ₁₃ S ₃ * 2 Na	704-00
NHS-ester	1000.05	C ₄₄ H ₄₇ N ₃ O ₁₅ S ₃ * 2 Na	704-01
Amino-derivative	923.07	C ₄₂ H ₅₁ N ₄ O ₁₂ S ₃ * Na	704-02
Maleimide	1025.10	C ₄₆ H ₅₀ N ₄ O ₁₄ S ₃ * 2 Na	704-03



DY-730



Absorption/emission max.:
734 nm / 755 nm (in Ethanol)

Molar absorbance:
 $240,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments, FISH microscopy, gel electrophoresis

Absorption/emission max.:
736 nm / 755 nm (in Ethanol)

Molar absorbance:
 $240,000 \text{ M}^{-1}\text{cm}^{-1}$

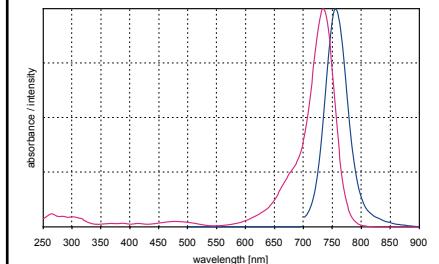
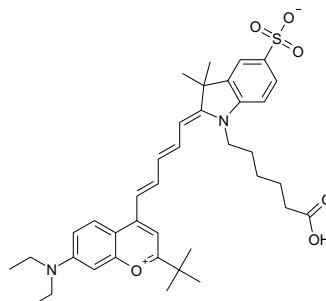
Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

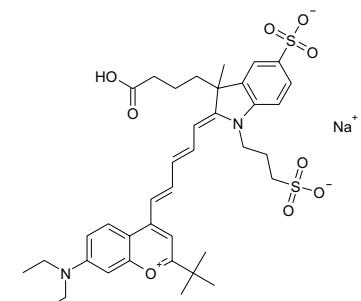
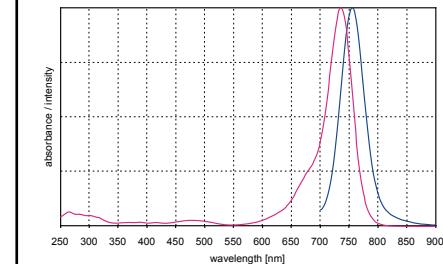
DY-731



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	660.88	C ₃₈ H ₄₈ N ₂ O ₆ S	730-00
NHS-ester	757.96	C ₄₂ H ₅₁ N ₃ O ₈ S	730-01
Amino-derivative	739.42	C ₄₀ H ₅₅ N ₄ O ₅ S * Cl	730-02
Maleimide	783.01	C ₄₄ H ₅₄ N ₄ O ₇ S	730-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	762.92	C ₃₈ H ₄₇ N ₂ O ₉ S ₂ * Na	731-00
NHS-ester	859.99	C ₄₂ H ₅₀ N ₃ O ₁₁ S ₂ * Na	731-01
Amino-derivative	783.03	C ₄₀ H ₅₄ N ₄ O ₈ S ₂	731-02
Maleimide	885.05	C ₄₄ H ₅₃ N ₄ O ₁₀ S ₂ * Na	731-03



DY-732



Absorption/emission max.:
735 nm / 756 nm (in Ethanol)

Molar absorbance:
 $240,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

Absorption/emission max.:
733 nm / 755 nm (in Ethanol)

Molar absorbance:
 $240,000 \text{ M}^{-1}\text{cm}^{-1}$

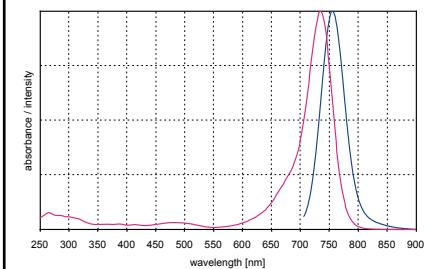
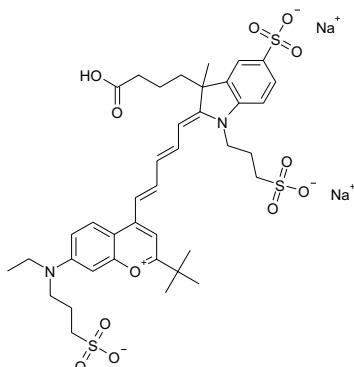
Comments:

- soluble in water, methanol, DMF, DMSO

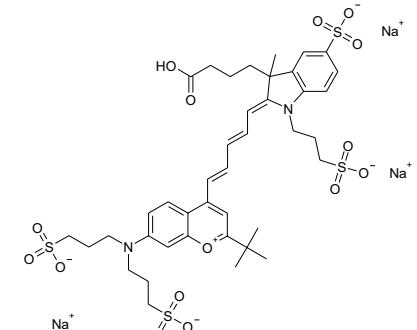
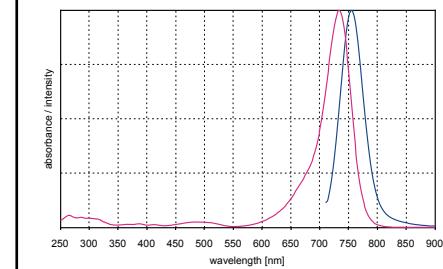
DY-734



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	879.00	C ₃₉ H ₄₈ N ₂ O ₁₂ S ₃ * 2 Na	732-00
NHS-ester	976.07	C ₄₃ H ₅₁ N ₃ O ₁₄ S ₃ * 2 Na	732-01
Amino-derivative	899.10	C ₄₁ H ₅₅ N ₄ O ₁₁ S ₃ * Na	732-02
Maleimide	1001.12	C ₄₅ H ₅₄ N ₄ O ₁₃ S ₃ * 2 Na	732-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	995.07	C ₄₀ H ₄₉ N ₂ O ₁₅ S ₄ * 3 Na	734-00
NHS-ester	1092.14	C ₄₄ H ₅₂ N ₃ O ₁₇ S ₄ * 3 Na	734-01
Amino-derivative	1015.19	C ₄₂ H ₅₆ N ₄ O ₁₄ S ₄ * 2 Na	734-02
Maleimide	1117.19	C ₄₆ H ₅₅ N ₄ O ₁₆ S ₄ * 3 Na	734-03



Label

NIR-excitation

DY-749



Absorption/emission max.:
759 nm / 780 nm (in Ethanol)

Molar absorbance:
 $240,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
- very hydrophilic

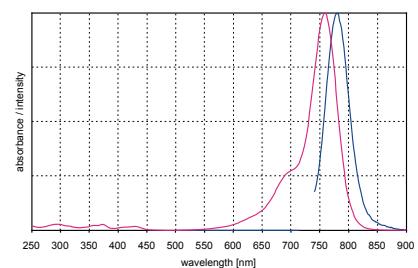
Absorption/emission max.:
751 nm / 774 nm (in Ethanol)

Molar absorbance:
 $270,000 \text{ M}^{-1}\text{cm}^{-1}$

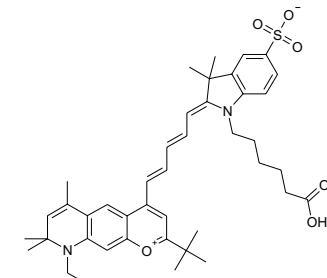
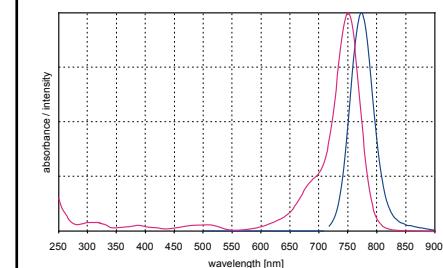
Comments:
- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments,
FISH microscopy, gel electrophoresis



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	936.99	C ₃₇ H ₄₃ N ₂ O ₁₄ S ₄ * 3 Na	749-00
NHS-ester	1034.06	C ₄₁ H ₄₆ N ₃ O ₁₆ S ₄ * 3 Na	749-01
Amino-derivative	957.09	C ₃₉ H ₅₀ N ₄ O ₁₃ S ₄ * 2 Na	749-02
Maleimide	1059.11	C ₄₃ H ₄₉ N ₄ O ₁₅ S ₄ * 3 Na	749-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	712.96	C ₄₂ H ₅₂ N ₂ O ₆ S	750-00
NHS-ester	810.03	C ₄₆ H ₅₅ N ₃ O ₈ S	750-01
Amino-derivative	791.49	C ₄₄ H ₅₉ N ₄ O ₅ S * Cl	750-02
Maleimide	835.08	C ₄₈ H ₅₈ N ₄ O ₇ S	750-03



Label

NIR-excitation

DY-750

DY-751



Absorption/emission max.:
752 nm / 772 nm (in Ethanol)

Molar absorbance:
 $270,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

Absorption/emission max.:
750 nm / 771 nm (in Ethanol)

Molar absorbance:
 $270,000 \text{ M}^{-1}\text{cm}^{-1}$

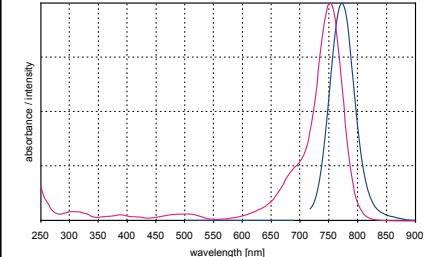
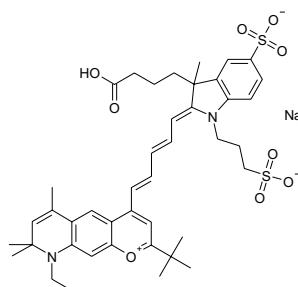
Comments:

- soluble in water, methanol, DMF, DMSO

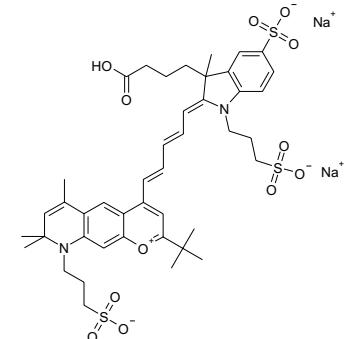
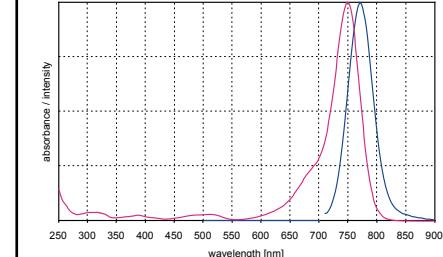
DY-752



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	814.99	C ₄₂ H ₅₁ N ₂ O ₉ S ₂ * Na	751-00
NHS-ester	912.07	C ₄₆ H ₅₄ N ₃ O ₁₁ S ₂ * Na	751-01
Amino-derivative	835.10	C ₄₄ H ₅₈ N ₄ O ₈ S ₂	751-02
Maleimide	937.13	C ₄₈ H ₅₇ N ₄ O ₁₀ S ₂ * Na	751-03



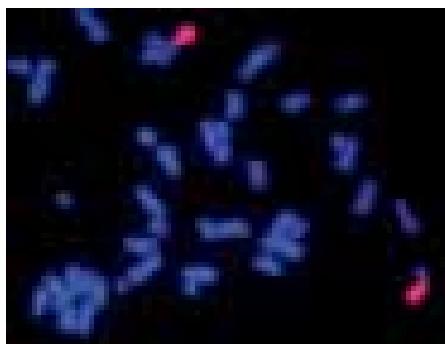
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	931.07	C ₄₃ H ₅₂ N ₂ O ₁₂ S ₂ * 2 Na	752-00
NHS-ester	1028.15	C ₄₇ H ₅₅ N ₃ O ₁₄ S ₃ * 2 Na	752-01
Amino-derivative	951.17	C ₄₅ H ₅₉ N ₄ O ₁₁ S ₃ * Na	752-02
Maleimide	1053.20	C ₄₉ H ₅₈ N ₄ O ₁₃ S ₃ * 2 Na	752-03



NIR-excitation

NIR-excitation

DY-754



Absorption/emission max.:
748 nm / 771 nm (in Ethanol)

Molar absorbance:
 $270,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– soluble in water, methanol, DMF

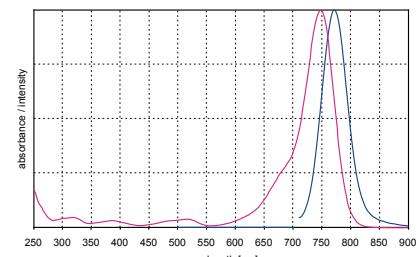
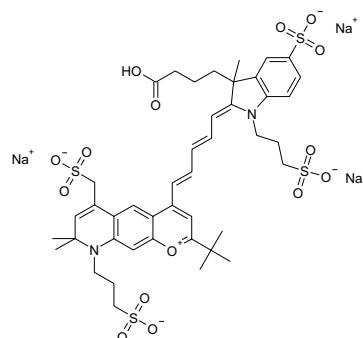
Absorption/emission max.:
772 nm / 787 nm (in Ethanol)

Molar absorbance:
 $240,000 \text{ M}^{-1}\text{cm}^{-1}$

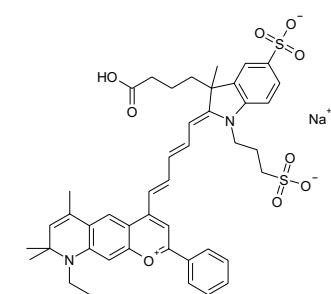
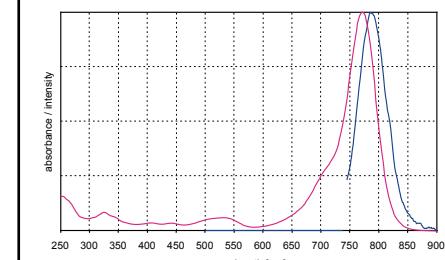
Comments:
– soluble in water, methanol, DMF, DMSO
– bright solid state emission
– suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis



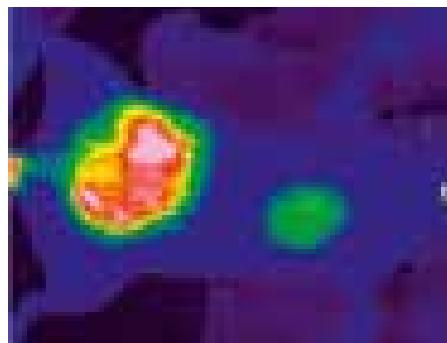
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	1033.12	C ₄₃ H ₅₁ N ₂ O ₁₅ S ₄ * 3 Na	754-00
NHS-ester	1130.19	C ₄₇ H ₅₄ N ₃ O ₁₇ S ₄ * 3 Na	754-01
Amino-derivative	1053.22	C ₄₅ H ₅₈ N ₄ O ₁₄ S ₄ * 2 Na	754-02
Maleimide	1155.25	C ₄₉ H ₅₇ N ₄ O ₁₆ S ₄ * 3 Na	754-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	834.98	C ₄₄ H ₄₇ N ₂ O ₉ S ₂ * Na	776-00
NHS-ester	932.07	C ₄₈ H ₅₀ N ₃ O ₁₁ S ₂ * Na	776-01
Amino-derivative	855.09	C ₄₆ H ₅₄ N ₄ O ₈ S ₂	776-02
Maleimide	957.12	C ₅₀ H ₅₃ N ₄ O ₁₀ S ₂ * Na	776-03



DY-777



Absorption/emission max.:
770 nm / 788 nm (in Ethanol)

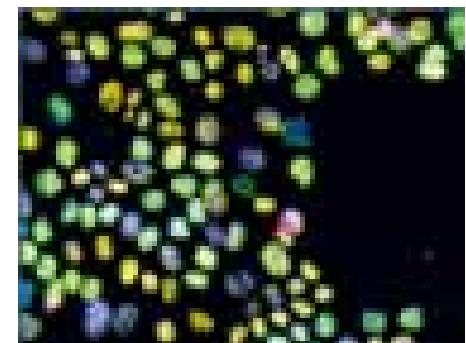
Molar absorbance:
 $240,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
- hydrophilic

Absorption/emission max.:
767 nm / 787 nm (in Ethanol)

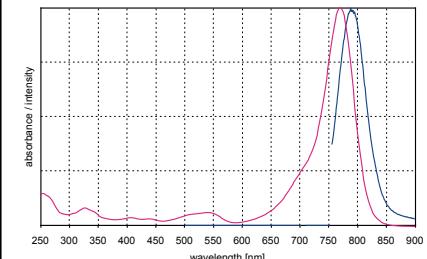
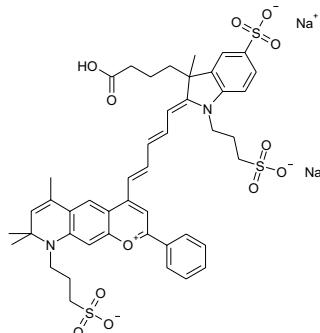
Molar absorbance:
 $240,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
- very hydrophilic

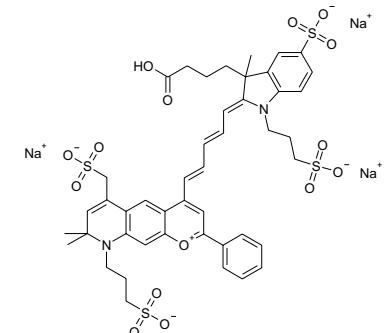
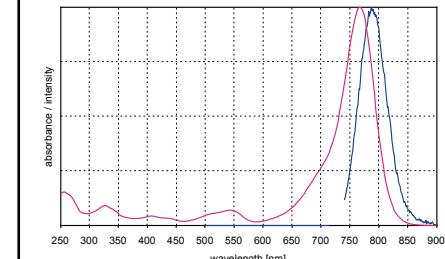


DY-778

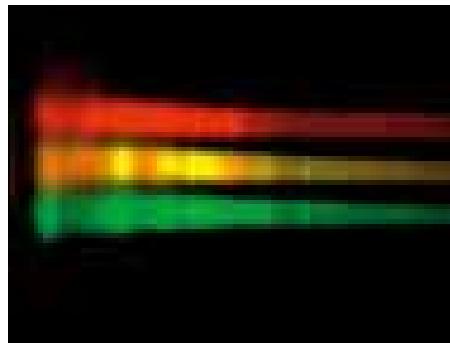
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	951.06	C ₄₅ H ₄₈ N ₂ O ₁₂ S ₃ * 2 Na	777-00
NHS-ester	1048.14	C ₄₉ H ₅₁ N ₃ O ₁₄ S ₃ * 2 Na	777-01
Amino-derivative	971.16	C ₄₇ H ₅₅ N ₄ O ₁₁ S ₃ * Na	777-02
Maleimide	1073.19	C ₅₁ H ₅₄ N ₄ O ₁₃ S ₃ * 2 Na	777-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	1053.11	C ₄₅ H ₄₇ N ₂ O ₁₅ S ₃ * 3 Na	778-00
NHS-ester	1150.18	C ₄₉ H ₅₀ N ₃ O ₁₇ S ₄ * 3 Na	778-01
Amino-derivative	1073.21	C ₄₇ H ₅₄ N ₄ O ₁₄ S ₄ * 2 Na	778-02
Maleimide	1175.24	C ₅₁ H ₅₃ N ₄ O ₁₆ S ₄ * 3 Na	778-03



DY-780



Absorption/emission max.:
783 nm / 799 nm (in Ethanol)

Molar absorbance:
170,000 M⁻¹cm⁻¹

Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments, FISH microscopy, gel electrophoresis

Absorption/emission max.:
784 nm / 796 nm (in Ethanol)

Molar absorbance:
170,000 M⁻¹cm⁻¹

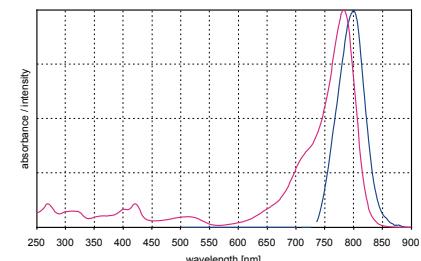
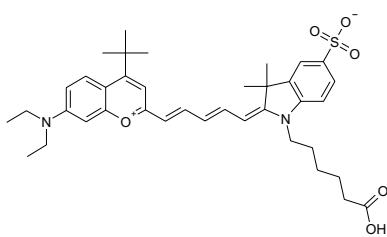
Comments:

- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

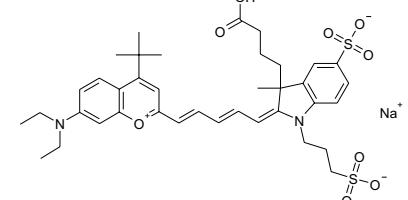
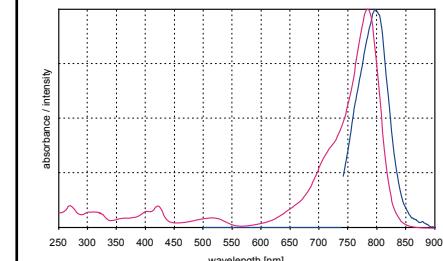


DY-781

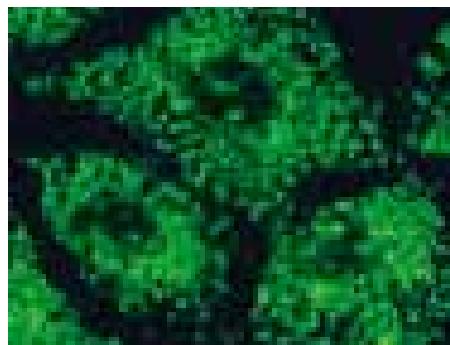
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	660.88	C ₃₈ H ₄₈ N ₂ O ₆ S	780-00
NHS-ester	757.95	C ₄₂ H ₅₁ N ₃ O ₈ S	780-01
Amino-derivative	739.42	C ₄₀ H ₅₅ N ₄ O ₅ S * Cl	780-02
Maleimide	783.01	C ₄₄ H ₅₄ N ₄ O ₇ S	780-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	762.92	C ₃₈ H ₄₇ N ₂ O ₉ S ₂ * Na	781-00
NHS-ester	860.00	C ₄₂ H ₅₀ N ₃ O ₁₁ S ₂ * Na	781-01
Amino-derivative	783.03	C ₄₀ H ₅₄ N ₄ O ₈ S ₂	781-02
Maleimide	885.05	C ₄₄ H ₅₃ N ₄ O ₁₀ S ₂ * Na	781-03



DY-782



Absorption/emission max.:
785 nm / 794 nm (in Ethanol)

Molar absorbance:
170,000 M⁻¹cm⁻¹

Comments:

- soluble in water, methanol
- two negative charges
- enhanced water solubility and polarity

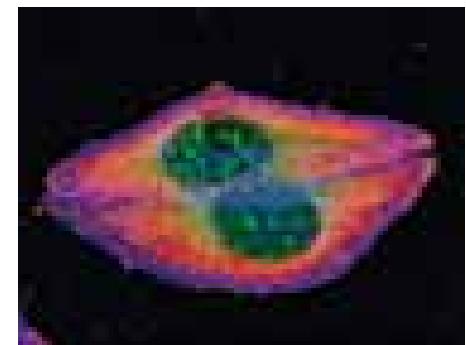
Absorption/emission max.:
777 nm / 791 nm (in Ethanol)

Molar absorbance:
280,000 M⁻¹cm⁻¹

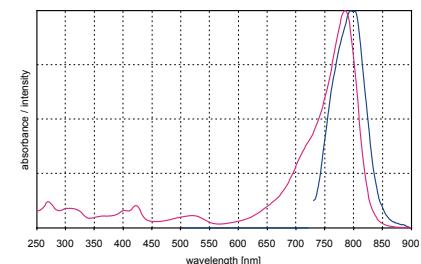
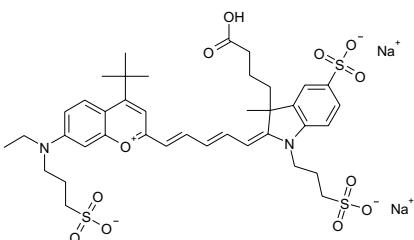
Comments:

- soluble in water, methanol, DMF
- hydrophilic

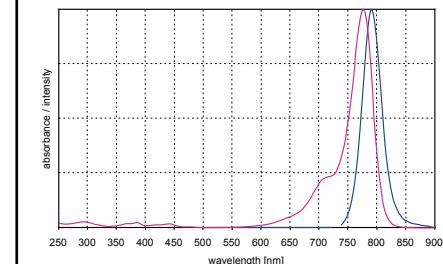
DY-800



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	879.00	C ₃₉ H ₄₈ N ₂ O ₁₂ S ₃ * 2 Na	782-00
NHS-ester	976.07	C ₄₃ H ₅₁ N ₃ O ₁₄ S ₃ * 2 Na	782-01
Amino-derivative	899.10	C ₄₁ H ₅₅ N ₄ O ₁₁ S ₃ * Na	782-02
Maleimide	1001.12	C ₄₅ H ₅₄ N ₄ O ₁₃ S ₃ * 2 Na	782-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	953.08	C ₄₅ H ₅₀ N ₂ O ₁₂ S ₃ * 2 Na	800-00
NHS-ester	1050.15	C ₄₉ H ₅₃ N ₃ O ₁₄ S ₃ * 2 Na	800-01
Amino-derivative	973.18	C ₄₇ H ₅₇ N ₄ O ₁₁ S ₃ * Na	800-02
Maleimide	1075.21	C ₅₁ H ₅₆ N ₄ O ₁₃ S ₃ * 2 Na	800-03



Structure on request

NIR-excitation

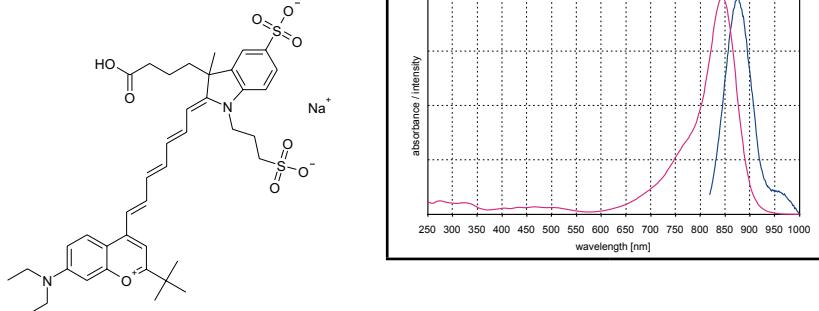
DY-831

Absorption/emission max.:
844 nm / 875 nm (in Ethanol)

Molar absorbance:
220,000 M⁻¹cm⁻¹

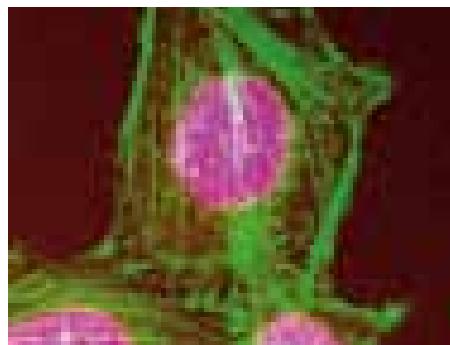
Comments:
– soluble in water, methanol, DMF, DMSO

Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	788.96	C ₄₀ H ₄₉ N ₂ O ₉ S ₂ * Na	831-00
NHS-ester	886.04	C ₄₄ H ₅₂ N ₃ O ₁₁ S ₂ * Na	831-01
Amino-derivative	809.06	C ₄₂ H ₅₆ N ₄ O ₈ S ₂	831-02
Maleimide	911.09	C ₄₆ H ₅₅ N ₄ O ₁₀ S ₂ * Na	831-03



MegaStokes-Dyes for multi-colour-detection

DY-480XL



Absorption/emission max.:
504 nm / 631 nm (in Ethanol)

Molar absorbance:
50,000 M⁻¹cm⁻¹

Comments:

- soluble in methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

Absorption/emission max.:
519 nm / 648 nm (in Ethanol)

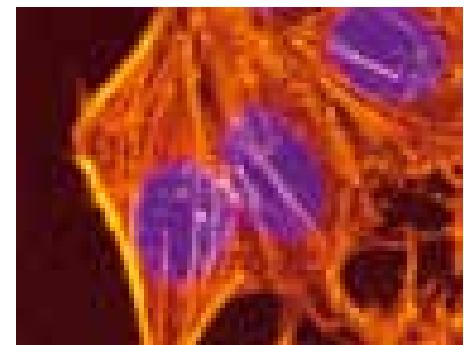
Molar absorbance:
50,000 M⁻¹cm⁻¹

Comments:

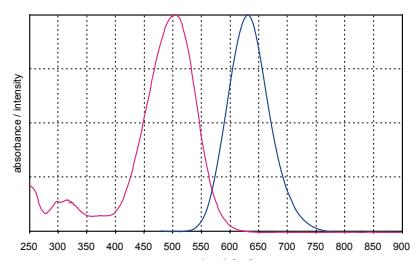
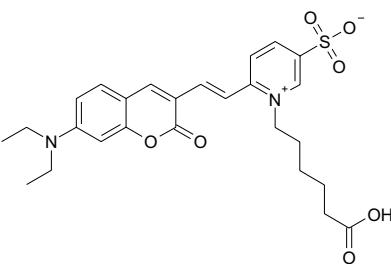
- soluble in water, methanol, ethanol, DMF, DMSO

MegaStokes-Dyes for multi-colour-detection

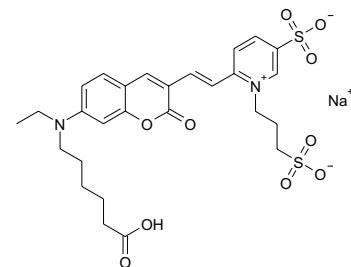
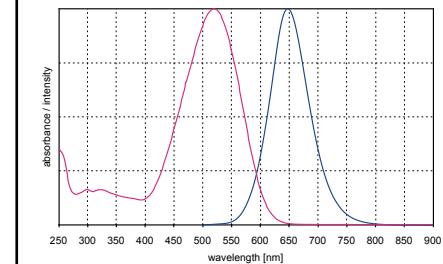
DY-481XL



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	514.60	C ₂₆ H ₃₀ N ₂ O ₇ S	480XL-00
NHS-ester	611.68	C ₃₀ H ₃₃ N ₃ O ₉ S	480XL-01
Amino-derivative	593.14	C ₂₈ H ₃₇ N ₄ O ₆ S * Cl	480XL-02
Maleimide	636.73	C ₃₂ H ₃₆ N ₄ O ₈ S	480XL-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	630.67	C ₂₇ H ₃₁ N ₂ O ₁₀ S ₂ * Na	481XL-00
NHS-ester	727.75	C ₃₁ H ₃₄ N ₃ O ₁₂ S ₂ * Na	481XL-01
Amino-derivative	650.78	C ₂₉ H ₃₈ N ₄ O ₉ S ₂	481XL-02
Maleimide	752.80	C ₃₃ H ₃₇ N ₄ O ₁₁ S ₂ * Na	481XL-03



MegaStokes-Dyes for multi-colour-detection

DY-485XL



Absorption/emission max.:
488 nm / 559 nm (in Ethanol)

Molar absorbance:
 $50,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

Absorption/emission max.:
509 nm / 590 nm (in Ethanol)

Molar absorbance:
 $50,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

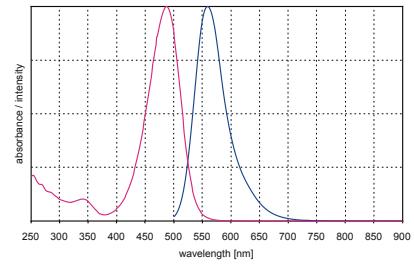
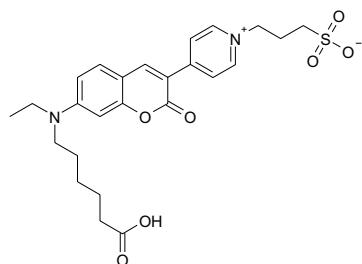
- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

MegaStokes-Dyes for multi-colour-detection

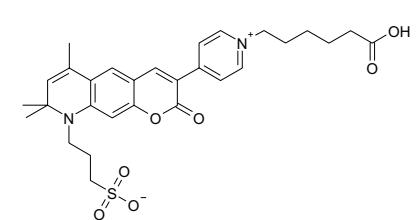
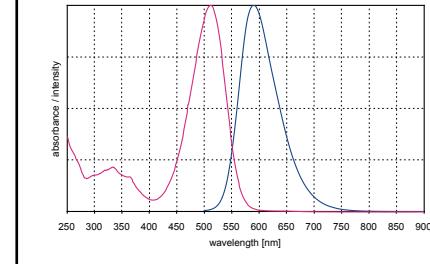
DY-510XL



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	502.59	C ₂₅ H ₃₀ N ₂ O ₇ S	485XL-00
NHS-ester	599.67	C ₂₉ H ₃₃ N ₃ O ₉ S	485XL-01
Amino-derivative	581.13	C ₂₇ H ₃₇ N ₄ O ₆ S * Cl	485XL-02
Maleimide	624.72	C ₃₁ H ₃₆ N ₄ O ₈ S	485XL-03

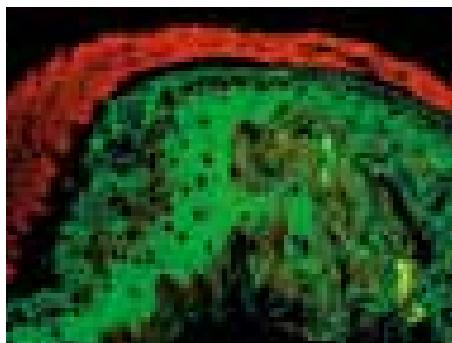


Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	554.67	C ₂₉ H ₃₄ N ₂ O ₇ S	510XL-00
NHS-ester	651.74	C ₃₃ H ₃₇ N ₃ O ₉ S	510XL-01
Amino-derivative	633.21	C ₃₁ H ₄₁ N ₄ O ₆ S * Cl	510XL-02
Maleimide	676.80	C ₃₅ H ₄₀ N ₄ O ₈ S	510XL-03



MegaStokes-Dyes for multi-colour-detection

DY-520XL



Absorption/emission max.:
522 nm / 662 nm (in Ethanol)

Molar absorbance:
 $50,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis

Absorption/emission max.:
526 nm / 666 nm (in Ethanol)

Molar absorbance:
 $50,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

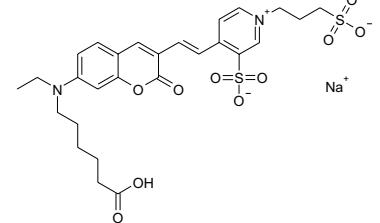
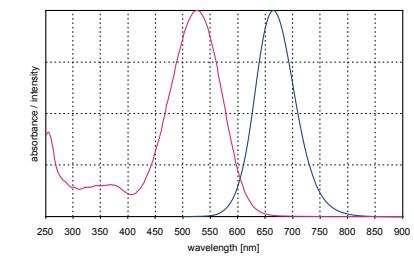
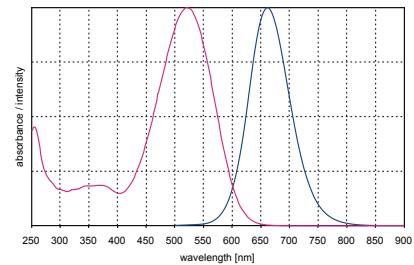
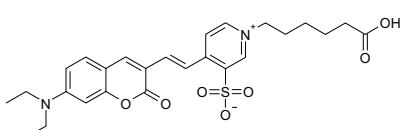
- soluble in water, methanol, DMF, DMSO

DY-521XL



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	514.60	C ₂₆ H ₃₀ N ₂ O ₇ S	520XL-00
NHS-ester	611.68	C ₃₀ H ₃₃ N ₃ O ₉ S	520XL-01
Amino-derivative	593.14	C ₂₈ H ₃₇ N ₄ O ₆ S * Cl	520XL-02
Maleimide	636.73	C ₃₂ H ₃₆ N ₄ O ₈ S	520XL-03

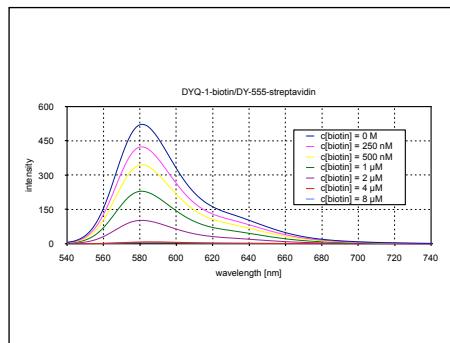
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	630.67	C ₂₇ H ₃₁ N ₂ O ₁₀ S ₂ * Na	521XL-00
NHS-ester	727.75	C ₃₁ H ₃₄ N ₃ O ₁₂ S ₂ * Na	521XL-01
Amino-derivative	650.78	C ₂₉ H ₃₈ N ₄ O ₉ S ₂	521XL-02
Maleimide	752.80	C ₃₃ H ₃₇ N ₄ O ₁₁ S ₂ * Na	521XL-03



Label

Quencher

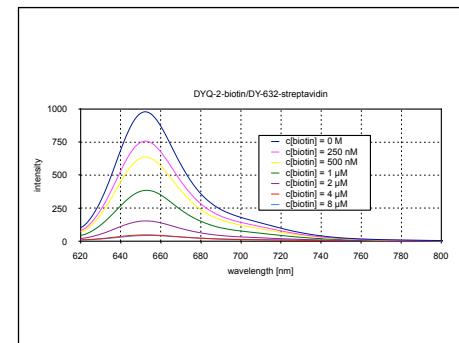
DYQ-1



Absorption max.:
641 nm (in Ethanol)

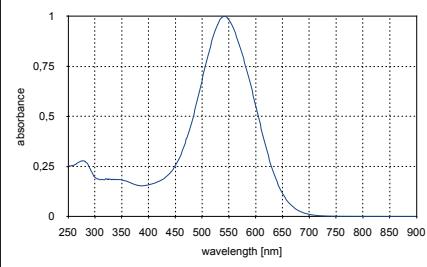
Molar absorbance:
 $90,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
- soluble in water, methanol, ethanol,
DMF, DMSO



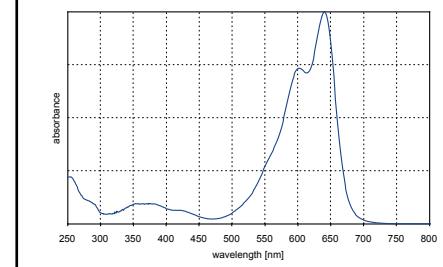
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	788.77	C ₃₂ H ₃₄ N ₆ O ₁₁ S ₂ * 2 Na	DYQ1-00
NHS-ester	885.84	C ₃₆ H ₃₇ N ₇ O ₁₃ S ₂ * 2 Na	DYQ1-01
Amino-derivative	808.87	C ₃₄ H ₄₁ N ₈ O ₁₀ S ₂ * Na	DYQ1-02
Maleimide	910.90	C ₃₈ H ₄₀ N ₈ O ₁₂ S ₂ * 2 Na	DYQ1-03

Structure on request



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	745.81	C ₃₆ H ₃₆ NO ₁₁ S ₂ * Na	DYQ2-00
NHS-ester	842.88	C ₄₀ H ₃₉ N ₂ O ₁₃ S ₂ * Na	DYQ2-01
Amino-derivative	765.91	C ₃₈ H ₄₃ N ₃ O ₁₀ S ₂	DYQ2-02
Maleimide	867.93	C ₄₂ H ₄₂ N ₃ O ₁₂ S ₂ * Na	DYQ2-03

Structure on request

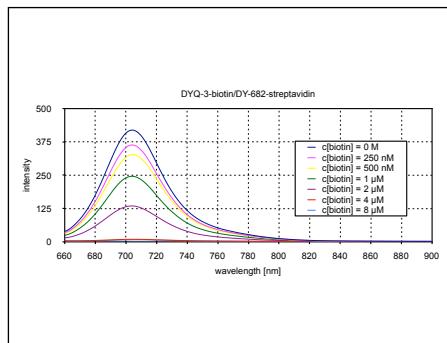


Label

Quencher

DYQ-2

DYQ-3



Absorption max.:
683 nm (in Ethanol)

Molar absorbance:
 $80,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– soluble in water, methanol

Absorption max.:
766 nm (in Ethanol)

Molar absorbance:
 $180,000 \text{ M}^{-1}\text{cm}^{-1}$

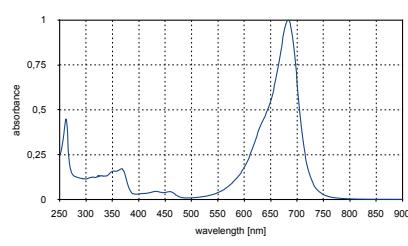
Comments:
– soluble in water, methanol

DYQ-4



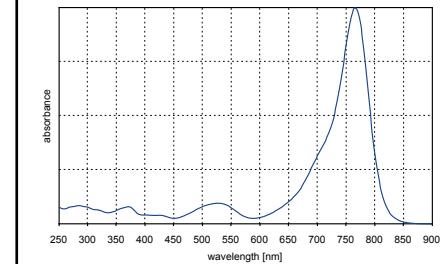
Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	802.86	C ₃₅ H ₃₆ N ₂ O ₁₁ S ₃ * 2 Na	DYQ3-00
NHS-ester	899.93	C ₃₉ H ₃₉ N ₃ O ₁₃ S ₃ * 2 Na	DYQ3-01
Amino-derivative	822.96	C ₃₇ H ₄₃ N ₄ O ₁₀ S ₃ * Na	DYQ3-02
Maleimide	924.98	C ₄₁ H ₄₂ N ₄ O ₁₂ S ₃ * 2 Na	DYQ3-03

Structure on request



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	796.94	C ₄₃ H ₄₅ N ₂ O ₉ S ₂ * Na	DYQ4-00
NHS-ester	894.02	C ₄₅ H ₄₈ N ₃ O ₁₁ S ₂ * Na	DYQ4-01
Amino-derivative	817.04	C ₄₃ H ₅₂ N ₄ O ₈ S ₂	DYQ4-02
Maleimide	919.15	C ₄₇ H ₅₁ N ₄ O ₁₀ S ₂ * Na	DYQ4-03

Structure on request



Label

Quencher

DYQ-425

Absorption spectrum (in PBS buffer pH 7.2)		Description	
Absorption max.	425 nm (PBS)		
Molar absorbance	24,500 M ⁻¹ cm ⁻¹ (PBS)		
Comments	polar water soluble quencher		
Available Modifications	Molecular weight	Molecular formula	
Carboxylic acid	784.37 g/mol	C ₂₃ H ₂₂ Br ₂ N ₄ O ₁₀ S ₂ * 2 Na	
NHS-ester	881.44 g/mol	C ₂₇ H ₂₅ Br ₂ N ₅ O ₁₂ S ₂ * 2 Na	
Biotin-derivative	1140.86 g/mol	C ₃₉ H ₅₀ Br ₂ N ₈ O ₁₃ S ₃ * 2 Na	

Label

Quencher

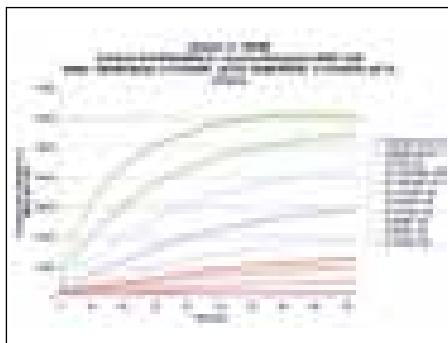
DYQ-700

Absorption spectrum (in PBS buffer pH 7.2)		Description	
Absorption max.	696 nm (PBS)		
Molar absorbance	48,500 M ⁻¹ cm ⁻¹ (PBS)		
Comments	polar water soluble quencher		
Available Modifications	Molecular weight	Molecular formula	
Carboxylic acid	808.92 g/mol	C ₃₉ H ₄₁ N ₆ O ₈ S ₂ * Na	
NHS-ester	905.99 g/mol	C ₄₃ H ₄₄ N ₇ O ₁₀ S ₂ * Na	
Biotin-derivative	1165.41 g/mol	C ₅₅ H ₆₉ N ₁₀ O ₁₁ S ₃ * Na	

DYQ-505

Absorption spectrum (in PBS buffer pH 7.2)		Description	
Absorption max.	504 nm (PBS)		
Molar absorbance	39,000 M ⁻¹ cm ⁻¹ (PBS)		
Comments	polar water soluble dark quencher		
Available Modifications	Molecular weight	Molecular formula	
Carboxylic acid	713.75 g/mol	C ₃₁ H ₃₃ N ₅ O ₈ S ₂ * 2 Na	
NHS-ester	810.82 g/mol	C ₃₅ H ₃₆ N ₆ O ₁₀ S ₂ * 2 Na	
Biotin-derivative	1070.24 g/mol	C ₄₇ H ₆₁ N ₉ O ₁₁ S ₃ * 2 Na	

DYQ-660



Absorption max.:
660 nm (in Ethanol)

Molar absorbance:
 $140,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, ethanol, DMF, DMSO
- bright solid state emission
- suitable for microarray experiments, FISH microscopy, gel electrophoresis

Absorption max.:
662 nm (in Ethanol)

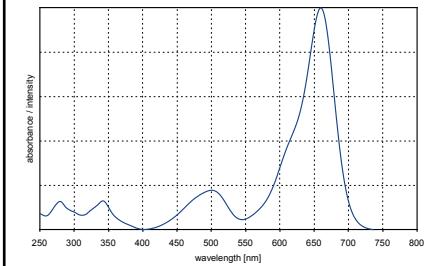
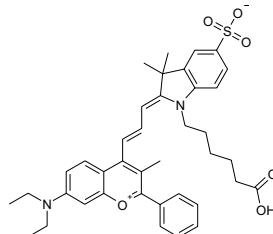
Molar absorbance:
 $140,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

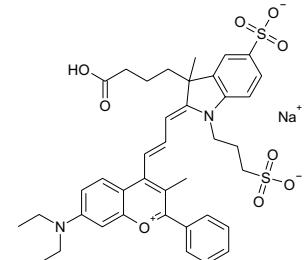
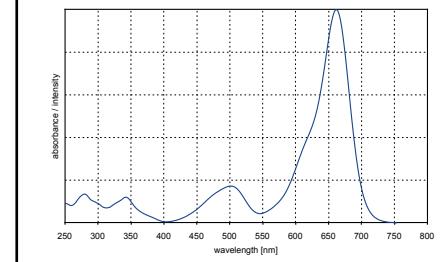
- soluble in water, methanol, DMF, DMSO
- bright solid state emission
- suitable for protein labeling, microarray experiments, FISH microscopy, gel electrophoresis
- improved water solubility

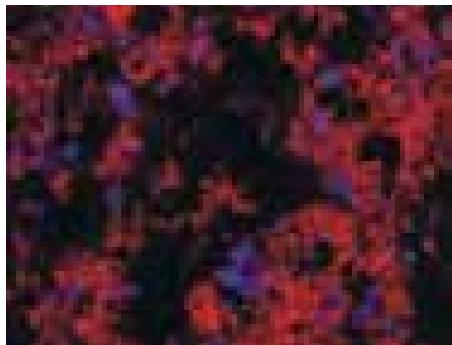


Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	668.86	C ₃₉ H ₄₄ N ₂ O ₆ S	660Q-00
NHS-ester	765.93	C ₄₃ H ₄₇ N ₃ O ₈ S	660Q-01
Amino-derivative	747.40	C ₄₁ H ₅₁ N ₄ O ₅ S * Cl	660Q-02
Maleimide	790.99	C ₄₅ H ₅₀ N ₄ O ₇ S	660Q-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	770.90	C ₃₉ H ₄₃ N ₂ O ₉ S ₂ * Na	661Q-00
NHS-ester	867.97	C ₄₃ H ₄₆ N ₃ O ₁₁ S ₂ * Na	661Q-01
Amino-derivative	791.01	C ₄₁ H ₅₀ N ₄ O ₈ S ₂	661Q-02
Maleimide	893.03	C ₄₅ H ₄₉ N ₄ O ₁₀ S ₂ * Na	661Q-03



FatRed

Absorption/emission max.:
656 nm / 680 nm (in Ethanol)

Molar absorbance:
 $95,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:

- soluble in methanol, ethanol, DMF, DMSO



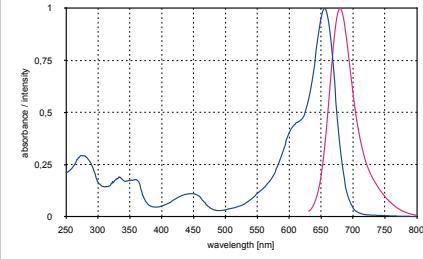
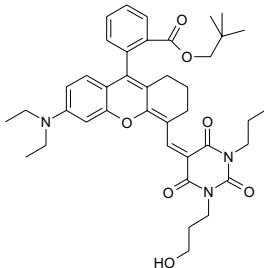
Absorption/emission max.:
513 nm / 580 nm (in Ethanol)

Molar absorbance:
 $120,000 \text{ M}^{-1}\text{cm}^{-1}$

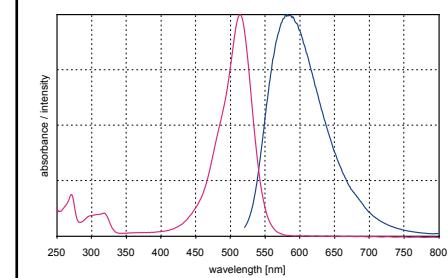
Comments:

- soluble in water, methanol, ethanol, DMF, DMSO

Molecular weight (g · mol ⁻¹)	Molecular formula	Product- number
697.88	C ₄₁ H ₅₁ N ₃ O ₇	FatRed-00



Molecular weight (g · mol ⁻¹)	Molecular formula	Product- number
751.94	C ₃₇ H ₅₀ N ₃ O ₈ S ₂ * Na	Stain01-00



Structure on request

Special Dyes

Protein-Stains

V07-05125

Absorption and emission spectra (in PBS buffer pH 7.2)		Description
	Absorption max.	520 nm (PBS)
	Emission max.	565 nm (PBS)
	Molar absorbance	45,000 M ⁻¹ cm ⁻¹ (in methanol)
	Comments	polar water soluble stain, bright fluorescence in presence of proteins (weak fluorescence in water/buffer without protein)
	Molecular weight	Molecular formula
	701.77 g/mol	C ₃₃ H ₃₇ NO ₉ S ₂ * 2 Na

Special Dyes

Protein-Stains

V07-06050

Absorption and emission spectra (in methanol)		Description
	Absorption max.	684 nm (methanol)
	Emission max.	717 nm (methanol)
	Molar absorbance	81,000 M ⁻¹ cm ⁻¹ (methanol)
	Comments	polar water soluble stain, bright fluorescence in presence of proteins (weak fluorescence in water/buffer without protein)
	Molecular weight	Molecular formula
	931.12 g/mol	C ₄₇ H ₅₉ N ₂ O ₁₂ S ₂ * Na

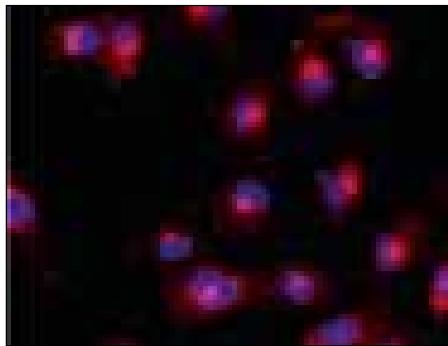
V07-05167

Absorption and emission spectra (in methanol)		Description
	Absorption max.:	520 nm (methanol)
	Emission max.:	563 nm (methanol)
	Molar absorbance:	125,000 M ⁻¹ cm ⁻¹ (methanol)
	Comments:	polar water soluble stain, bright fluorescence in presence of proteins (weak fluorescence in water/buffer without protein)
	Molecular weight	Molecular formula
	806.93 g/mol	C ₃₉ H ₄₇ N ₂ O ₁₁ S ₂ * Na

V07-07150

Absorption and emission spectra (in methanol)		Description
	Absorption max.	615 nm (methanol)
	Emission max.	648 nm (methanol)
	Molar absorbance	80,000 M ⁻¹ cm ⁻¹ (methanol)
	Comments	polar water soluble stain, bright fluorescence in presence of proteins (weak fluorescence in water/buffer without protein)
	Molecular weight	Molecular formula
	802.99 g/mol	C ₄₄ H ₅₄ N ₂ O ₁₀ S

V02-07027



Absorption/emission max.:
477 nm / 522 nm (in Ethanol)

Molar absorbance:
60,000 M⁻¹cm⁻¹

Comments:

- soluble in acetone, ethanol, DMF, DMSO
- suitable as DNA-stain

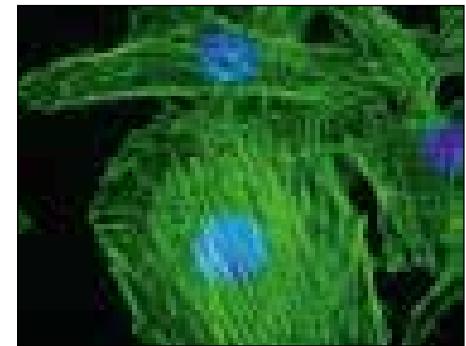
Absorption/emission max.:
490 nm / 537 nm (in Ethanol)

Molar absorbance:
60,000 M⁻¹cm⁻¹

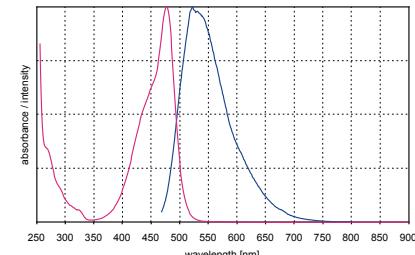
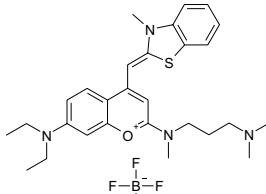
Comments:

- soluble in acetone, ethanol, DMF, DMSO
- suitable as DNA-stain

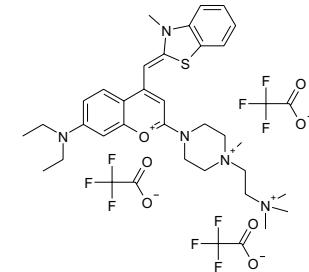
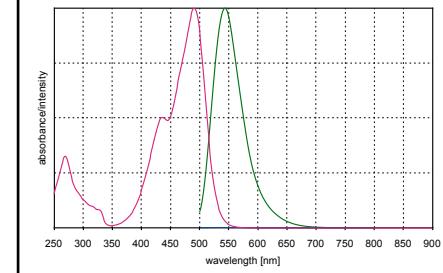
V02-08078



Molecular weight (g · mol ⁻¹)	Molecular formula	Product- number
564.50	C ₂₈ H ₃₇ N ₄ OS * BF ₄	V02-07027



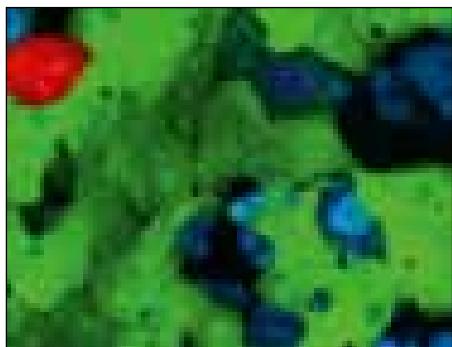
Molecular weight (g · mol ⁻¹)	Molecular formula	Product- number
887.88	C ₃₂ H ₄₆ N ₅ OS * 3 C ₂ F ₃ O ₂	V02-08078



Special Dyes

DNA-Stains

V13-01184



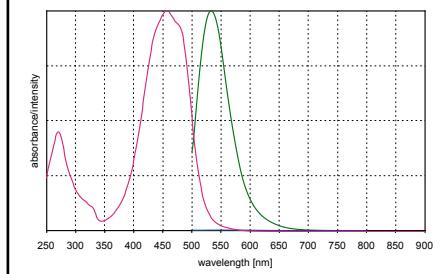
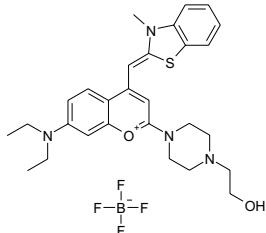
Absorption/emission max.:
481 nm / 526 nm (in Ethanol)

Molar absorbance:
80,000 M⁻¹cm⁻¹

Comments:

- soluble in acetone, ethanol, DMF, DMSO
- suitable as DNA-stain

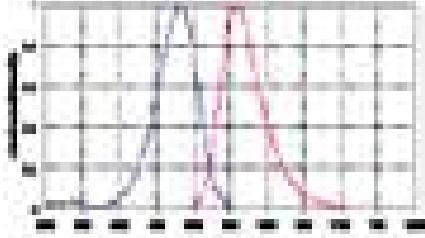
Molecular weight (g · mol ⁻¹)	Molecular formula	Product- number
578.48	C ₂₈ H ₃₅ N ₄ O ₂ S * BF ₄	V13-01184



Special Dyes

MegaStokes-Dyes

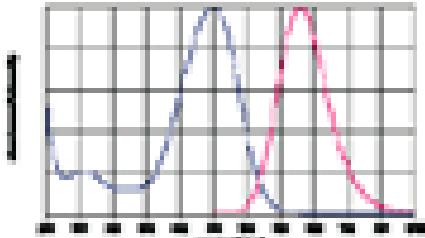
V07-07059

Absorption and emission spectra (in ethanol)		Description	
		Absorption max.	485 nm (ethanol)
		Emission max.	558 nm (ethanol)
		Molar absorbance	62,000 M ⁻¹ cm ⁻¹ (ethanol)
		Comments	Nonpolar MegaStokes Dye for multi-colour-detection. Hydroxy-substituted as precursor for phosphoramidite
		Molecular weight	Molecular formula
		440.24 g/mol	C ₂₁ H ₂₅ N ₂ O ₃ * BF ₄

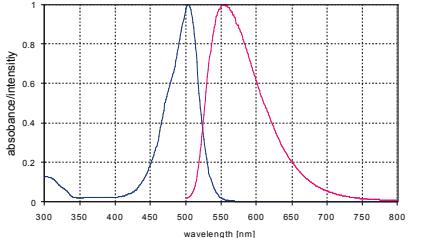
Special Dyes

MegaStokes-Dyes

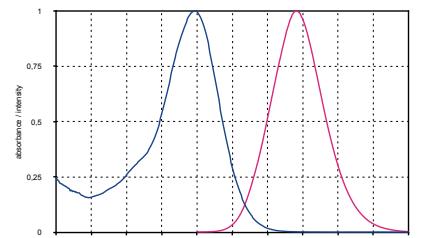
V07-08145

Absorption and emission spectra (in methanol)		Description	
		Absorption max.	503 nm (methanol)
		Emission max.	637 nm (methanol)
		Molar absorbance	53,000 M ⁻¹ cm ⁻¹ (methanol)
		Comments	lipophilic MegaStokes Dye soluble in nonpolar solvents like acetone, acetonitrile, DMF, dichloromethane, THF or DMSO
		Molecular weight	Molecular formula
		624.89 g/mol	C ₃₆ H ₅₂ N ₂ O ₅ S

V10-02026

Absorption and emission spectra (in methanol)		Description	
		Absorption max.	502 nm (methanol)
		Emission max.	556 nm (methanol)
		Molar absorbance	120,000 M ⁻¹ cm ⁻¹ (methanol)
		Comments	Nonpolar MegaStokes Dye for multi-colour-detection. Hydroxy-substituted as precursor for phosphoramidite
		Molecular weight	Molecular formula
		460.58 g/mol	C ₂₈ H ₃₂ N ₂ O ₄

V19-02035

Absorption and emission spectra (in methanol)		Description	
		Absorption max.	497 nm (methanol)
		Emission max.	641 nm (methanol)
		Molar absorbance	36,000 M ⁻¹ cm ⁻¹ (methanol)
		Comments	lipophilic MegaStokes Dye soluble in nonpolar solvents like acetone, acetonitrile, DMF, dichloromethane, THF or DMSO
		Molecular weight	Molecular formula
		520.66 g/mol	C ₂₉ H ₃₂ N ₂ O ₅ S

Special Dyes

MegaStokes-Dyes

V07-10147

Absorption and emission spectra (in PBS)		Description	
		Absorption max.	501 nm (PBS)
		Emission max.	645 nm (PBS)
		Molar absorbance	39,000 M ⁻¹ cm ⁻¹ (PBS)
		Comments	Cationic MegaStokes Dye with intercalating properties
		Molecular weight	Molecular formula
		442.00 g/mol	C ₂₅ H ₃₂ N ₃ O ₂ * Cl

Special Dyes

MegaStokes-Dyes

V11-04139

Absorption and emission spectra (in methanol)		Description	
		Absorption max.	635 nm (methanol)
		Emission max.	674 nm (methanol)
		Molar absorbance	113,000 M ⁻¹ cm ⁻¹ (methanol)
		Comments	Cationic MegaStokes Dye with intercalating properties. Excitation also at 530nm possible.
		Molecular weight	Molecular formula
		634.57 g/mol	C ₃₇ H ₄₃ N ₂ O ₂ * BF ₄

V07-10159

Absorption and emission spectra (in PBS)		Description	
		Absorption max.	534 nm (PBS)
		Emission max.	697 nm (PBS)
		Molar absorbance	31,000 M ⁻¹ cm ⁻¹ (PBS)
		Comments	Cationic MegaStokes Dye with intercalating properties
		Molecular weight	Molecular formula
		492.06 g/mol	C ₂₉ H ₃₄ N ₃ O ₂ * Cl

V11-04141

Absorption and emission spectra (in methanol)		Description	
		Absorption max.	653 nm (methanol)
		Emission max.	701 nm (methanol)
		Molar absorbance	113,000 M ⁻¹ cm ⁻¹ (methanol)
		Comments	Cationic MegaStokes Dye with intercalating properties. Excitation also at 530nm possible.
		Molecular weight	Molecular formula
		654.56 g/mol	C ₃₉ H ₃₉ N ₂ O ₂ * BF ₄

Special Dyes

Direct from Research

V07-08026

Absorption and emission spectra (in PBS buffer pH 7.2)		Description	
		Absorption max.	356 nm (PBS)
		Emission max.	455 nm (PBS)
		Molar absorbance	$18,000 \text{ M}^{-1}\text{cm}^{-1}$ (methanol)
		Quantum yield	0.55 (PBS)
		Comments	polar water soluble dye, very bright blue fluorescence, hydrophilic alternative to coumarine dyes e.g. (7-Amino-4-methyl-3-coumarinyl)-acetic acid, (7-Hydroxy-4-methyl-3-coumarinyl)-acetic acid, Marina Blue™ or (7-Dimethylamino-4-coumarinyl)-acetic acid
Available Modifications	Molecular weight	Molecular formula	
Carboxylic acid	632.62 g/mol	$\text{C}_{27}\text{H}_{26}\text{N}_2\text{O}_9\text{S}_2 * 2 \text{ Na}$	
NHS-ester	729.70 g/mol	$\text{C}_{31}\text{H}_{29}\text{N}_3\text{O}_{11}\text{S}_2 * 2 \text{ Na}$	

Special Dyes

Direct from Research

V07-10170

Absorption and emission spectra (in PBS buffer pH 7.2)		Description	
		Absorption max.	368 nm (PBS)
		Emission max.	474 nm (PBS)
		Molar absorbance	$13,200 \text{ M}^{-1}\text{cm}^{-1}$ (PBS)
		Quantum yield	0.85 (PBS)
		Comments	polar water soluble dye, very bright blue fluorescence, hydrophilic alternative to coumarine dyes e.g. Marina Blue™ or (7-Dimethylamino-4-coumarinyl)-acetic acid
Available Modifications	Molecular weight	Molecular formula	
Carboxylic acid	542.52 g/mol	$\text{C}_{22}\text{H}_{19}\text{N}_2\text{O}_9\text{S}_2 * \text{Na}$	
NHS-ester	639.60 g/mol	$\text{C}_{26}\text{H}_{22}\text{N}_3\text{O}_{11}\text{S}_2 * \text{Na}$	

V10-02153

Absorption and emission spectra (in PBS buffer pH 7.2)		Description	
		Absorption max.	362 nm (PBS)
		Emission max.	459 nm (PBS)
		Molar absorbance	$25,000 \text{ M}^{-1}\text{cm}^{-1}$ (methanol)
		Quantum yield	0.52 (PBS)
		Comments	polar water soluble dye, very bright blue fluorescence, hydrophilic alternative to coumarine dyes e.g. (7-Amino-4-methyl-3-coumarinyl)-acetic acid, (7-Hydroxy-4-methyl-3-coumarinyl)-acetic acid, Marina Blue™ or (7-Dimethylamino-4-coumarinyl)-acetic acid
Available Modifications	Molecular weight	Molecular formula	
Carboxylic acid	632.62 g/mol	$\text{C}_{27}\text{H}_{26}\text{N}_2\text{O}_9\text{S}_2 * 2 \text{ Na}$	
NHS-ester	729.70 g/mol	$\text{C}_{31}\text{H}_{29}\text{N}_3\text{O}_{11}\text{S}_2 * 2 \text{ Na}$	

V07-11049

Absorption and emission spectra (in PBS buffer pH 7.2)		Description	
		Absorption max.	380 nm (PBS)
		Emission max.	484 nm (PBS)
		Molar absorbance	$13,000 \text{ M}^{-1}\text{cm}^{-1}$ (PBS)
		Quantum yield	0.85 (PBS)
		Comments	polar water soluble dye, very bright blue fluorescence, hydrophilic alternative to coumarine dyes e.g. Marina Blue™ or (7-Dimethylamino-4-coumarinyl)-acetic acid
Available Modifications	Molecular weight	Molecular formula	
Carboxylic acid	542.52 g/mol	$\text{C}_{22}\text{H}_{19}\text{N}_2\text{O}_9\text{S}_2 * \text{Na}$	
NHS-ester	639.60 g/mol	$\text{C}_{26}\text{H}_{22}\text{N}_3\text{O}_{11}\text{S}_2 * \text{Na}$	

Special Dyes

Direct from Research

V10-04059

Absorption and emission spectra (in PBS buffer pH 7.2)		Description	
Absorption max.	383 nm (PBS)	Emission max.	516 nm (PBS)
Molar absorbance	19,000 M ⁻¹ cm ⁻¹ (PBS)	Quantum yield	0.70 (PBS)
Comments	polar water soluble dye, very bright green fluorescence, large Stokes shift, suitable also for multi-colour-detection, alternative to Pacific Orange™	Available Modifications	Molecular weight
Carboxylic acid	448.50 g/mol	C ₂₁ H ₂₄ N ₂ O ₇ S	Molecular formula
NHS-ester	545.57 g/mol	C ₂₅ H ₂₇ N ₃ O ₉ S	

Special Dyes

Direct from Research

V07-11082

Absorption and emission spectra (in PBS buffer pH 7.2)		Description	
Absorption max.	375 nm (PBS)	Emission max.	543 nm (PBS)
Molar absorbance	16,000 M ⁻¹ cm ⁻¹ (PBS)	Quantum yield	0.62 (PBS)
Comments	polar water soluble dye, very bright green fluorescence, large Stokes shift, suitable also for multi-colour-detection, alternative to Pacific Orange™	Available Modifications	Molecular weight
Carboxylic acid	447.47 g/mol	C ₂₀ H ₂₁ N ₃ O ₇ S	Molecular formula
NHS-ester	544.54 g/mol	C ₂₄ H ₂₄ N ₄ O ₉ S	

V07-11176

Absorption and emission spectra (in PBS buffer pH 7.2)		Description	
Absorption max.	404 nm (PBS)	Emission max.	535 nm (PBS)
Molar absorbance	18,000 M ⁻¹ cm ⁻¹ (PBS)	Quantum yield	0.73 (PBS)
Comments	polar water soluble dye, very bright green fluorescence, large Stokes shift, suitable also for multi-colour-detection, alternative to Pacific Orange™	Available Modifications	Molecular weight
Carboxylic acid	592.58 g/mol	C ₂₆ H ₂₁ N ₂ O ₉ S ₂ * Na	Molecular formula
NHS-ester	689.66 g/mol	C ₃₀ H ₂₄ N ₃ O ₁₁ S ₂ * Na	

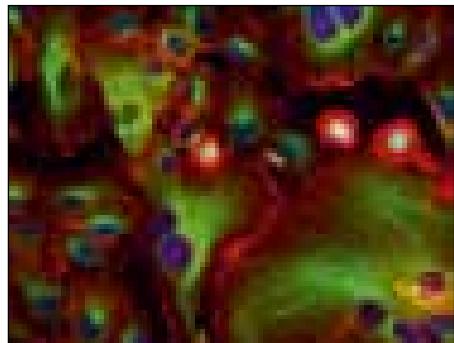
V19-01139

Absorption and emission spectra (in PBS buffer pH 7.2)		Description	
Absorption max.	397 nm (PBS)	Emission max.	572 nm (PBS)
Molar absorbance	18,000 M ⁻¹ cm ⁻¹ (PBS)	Quantum yield	0.28 (PBS)
Comments	polar water soluble dye, bright orange fluorescence, large Stokes shift, suitable also for multi-colour-detection, alternative to Cascade Yellow™ or Pacific Orange™	Available Modifications	Molecular weight
Carboxylic acid	490.54 g/mol	C ₂₂ H ₂₆ N ₄ O ₇ S	Molecular formula
NHS-ester	587.61 g/mol	C ₂₆ H ₂₉ N ₅ O ₉ S	

Special Dyes

Direct from Research

V08-13007



Absorption/emission max.:
699 nm / 723 nm (in Ethanol)

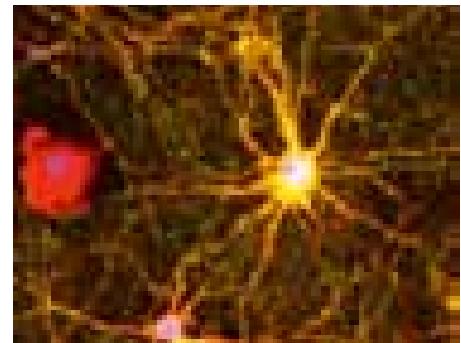
Molar absorbance:
 $140,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– useful for multicolour detection

Absorption/emission max.:
656 nm / 684 nm (in Ethanol)

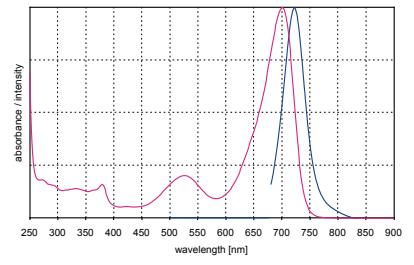
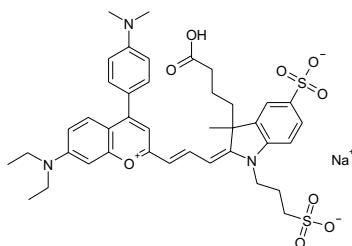
Molar absorbance:
 $140,000 \text{ M}^{-1}\text{cm}^{-1}$

Comments:
– useful for multicolour detection

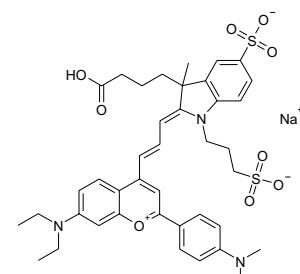
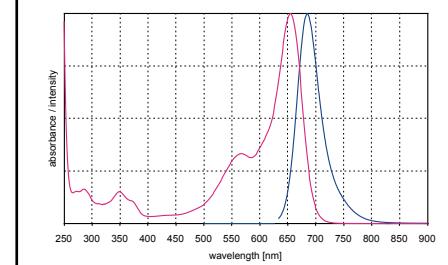


V13-04159

Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	799.95	C ₄₀ H ₄₆ N ₃ O ₉ S ₂ * Na	V08-13007-00
NHS-ester	897.02	C ₄₄ H ₄₉ N ₄ O ₁₁ S ₂ * Na	V08-13007-01
Amino-derivative	820.05	C ₄₂ H ₅₃ N ₅ O ₈ S ₂	V08-13007-02
Maleimide	922.06	C ₄₆ H ₅₂ N ₅ O ₁₀ S ₂ * Na	V08-13007-03



Available Modification	Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
Carboxylic acid	799.95	C ₄₀ H ₄₆ N ₃ O ₉ S ₂ * Na	V13-04159-00
NHS-ester	897.02	C ₄₄ H ₄₉ N ₄ O ₁₁ S ₂ * Na	V13-04159-01
Amino-derivative	820.05	C ₄₂ H ₅₃ N ₅ O ₈ S ₂	V13-04159-02
Maleimide	922.06	C ₄₆ H ₅₂ N ₅ O ₁₀ S ₂ * Na	V13-04159-03



Special Dyes

Direct from Research

V07-07103

Absorption and emission spectra (in methanol)		Description	
	Absorption max.	863 nm (methanol)	
	Emission max.	890 nm (methanol)	
	Molar absorbance	178,000 M ⁻¹ cm ⁻¹ (methanol)	
	Comments	IR-dye	
Available Modifications		Molecular weight	Molecular formula
Carboxylic acid	907.10 g/mol	C ₄₈ H ₅₅ N ₂ O ₁₀ S ₂ * Na	
NHS-ester	1004.17 g/mol	C ₅₂ H ₅₈ N ₃ O ₁₂ S ₂ * Na	

Special Dyes

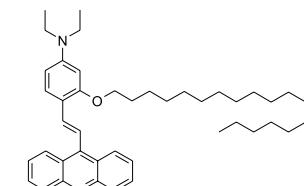
Sensor Dyes

DY-660-S

Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
620.96	C ₄₃ H ₆₀ N ₂ O	660-S



Absorption max.: 665 nm (in Ethanol)
Molar absorbance: 20,000 M⁻¹cm⁻¹
Comments: soluble in ethanol, acetonitril, DMF, DMSO



V07-07113

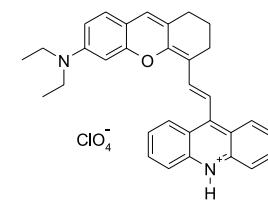
Absorption and emission spectra (in methanol)		Description	
	Absorption max.	888 nm (methanol)	
	Emission max.	935 nm (methanol)	
	Molar absorbance	151,000 M ⁻¹ cm ⁻¹ (methanol)	
	Comments	IR-dye	
Available Modifications		Molecular weight	Molecular formula
Carboxylic acid	923.16 g/mol	C ₄₈ H ₅₅ N ₂ O ₉ S ₃ * Na	
NHS-ester	1020.24 g/mol	C ₅₂ H ₅₈ N ₃ O ₁₁ S ₃ * Na	

DY-840-S

Molecular weight (g · mol ⁻¹)	Molecular formula	Product-number
559.01	C ₃₂ H ₃₁ N ₂ O · ClO ₄	840-S



Absorption max.: 840 nm (in Ethanol)
Molar absorbance: 65,000 M⁻¹cm⁻¹
Comments: soluble in ethanol, acetonitril, DMF, DMSO

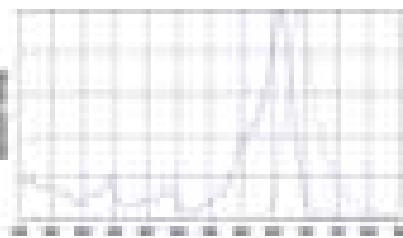


Special Dyes

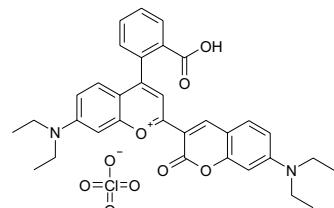
Laser dyes

DY-660-X

Molecular weight (g · mol ⁻¹)	Molecular formula	Product- number
637.09	C ₃₃ H ₃₃ N ₂ O ₅ · ClO ₄	660-X

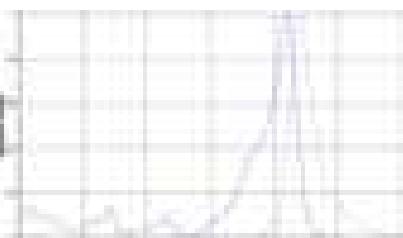


Absorption/
emission max.: 662 nm / 695 nm
(in Ethanol)
Molar absorbance: 80,000 M⁻¹cm⁻¹
Comments: soluble in dichloro-
methane, acetonitril,
DMF, DMSO

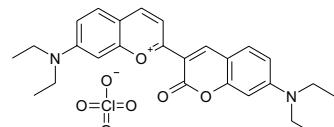


DY-665-X

Molecular weight (g · mol ⁻¹)	Molecular formula	Product- number
516.98	C ₂₆ H ₂₉ N ₂ O ₃ · ClO ₄	665-X



Absorption/
emission max.: 670 nm / 692 nm
(in dichloromethane)
Molar absorbance: 80,000 M⁻¹cm⁻¹
Comments: soluble in dichloro-
methane, acetonitril,
DMF, DMSO



Dyomics offers a wide range of labeled haptens like biotin, phalloidin, mycotoxins, amino acids or nucleotides. These haptens can easily be coupled to all of our fluorescent labels.

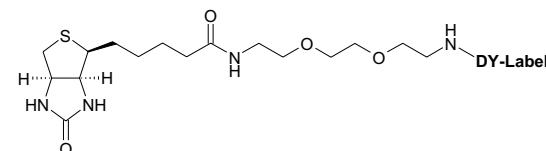
If you have need for a special hapten please ask us. We will do the coupling with our superior dyes.

Haptens

Fluorescently labeled biotins

The binding pairs streptavidin – avidin / biotin or antibody / antigen are very common in cellbiology and molecular biology in applications like *in situ* hybridisation, histochemistry or flow cytometry. With such systems it is possible to probe poor detectable molecules.

The building block for preparing biotin – DY-dye – conjugates is the biotin-PEO3-amine. PEO3 is a spacer consisting of ethyleneglycol units. This spacer between the biotin and the dye separates the biotin moiety from its point of attachment, enhances the ability to bind to the binding sites of avidin or streptavidin and improves the hydrophilic character of the biotin conjugate.



Unit size
500 µg (higher amounts on request)

Fluorescently labeled phalloidins

Phalloidin binds very selectively to actin and allows the generation of impressive images in combination with a fluorophor. 300 units (the standard sales unit) are sufficient for staining 300 microscopy slides. Methanol is an appropriate solvent for the labeled phalloidins.

For all our labels phalloidin-conjugates are available. Please add -33 to the order number of the dye of your choice.



Unit size
10 nmol/300 units (higher amounts on request)

Fluorescently labeled streptavidin

We offer streptavidin conjugates of all our DY-labels and MegaStokes-Dyes. Each streptavidin conjugate will be delivered with a batch specific absorption spectrum. The conjugates will be shipped in PBS, pH 7.5, 100 mM at a concentration of 1 mg/ml.

Unit size	1 mg at 1 mg/ml (higher amounts on request)
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Fluorescently labeled antibodies for secondary detection

Secondary antibodies are very common in applications like fluorescence microscopy and flow cytometry. Dyomics offers several fluorescently labeled polyclonal antibodies for secondary detection, among them

- goat anti-rabbit IgG
- goat anti-mouse IgG &
- goat anti-human IgG

(all affinity purified, H&L). The IgGs will be delivered dissolved in PBS, pH 7.5, 100 mM at a concentration of 1 mg/ml. A batch specific absorption spectrum will be provided.

Unit size	1 mg at 1 mg/ml (higher amounts on request)
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Others

If you need an antibody or another protein not listed in this catalogue feel free to contact Dyomics. We can assist you in finding a supplier and we can arrange the import, too. Thus, Dyomics can act as a one stop source for your secondary detection reagents.

DY-Starter-Kit

To allow customers to evaluate a set of DY-labels covering the spectrum from 550 to 800 nm, we offer the **DY-StarterKit**, which contains seven 1 mg portions of DY-NHS-esters of your choice.

Product-number	SK-1
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RedPack

Dyomics has special expertise in the field of red fluorescent labels. To allow customers to test and select the optimal red fluorescent label for their individual application, we offer the **RedPack**, a collection of seven red fluorescent NHS-esters (1 mg of DY-630, DY-632, DY-633, DY-634, DY-636, DY-649 & DY-652).

Product-number	RP-1
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ChipPack

The **ChipPack™** contains 20 µg aliquots of DY-547-NHS/DY-647-NHS-esters. This packing size is convenient for labeling usual amounts of cDNA obtained by reverse transcription with enzymatic incorporation of aminoallyl-dUTP.

size	# of aliquots	Product-number
S	10 x 2	CP-S
M	25 x 2	CP-M
L	50 x 2	CP-L

Protein Labeling Kits

Dyomics offers labeling kits in cooperation with emp Biotech. The core components of these kits are the CentriSep purification columns (Princeton Separation, Inc.) which allow a fast and easy separation of the conjugate from unconjugated dye.

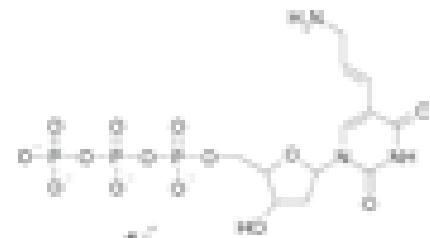
Unit size	Product-number
5 reactions	LK-05
10 reactions	LK-10

For an individual combination of labels for your special need do not hesitate to ask for your personal offer.

Aminoallyl-dUTP for enzymatic incorporation

Molecular formula	C ₁₂ H ₁₆ N ₃ O ₁₄ P ₃ * 4 Li
MW	546.95 gmol ⁻¹
Concentration	10 mM in water

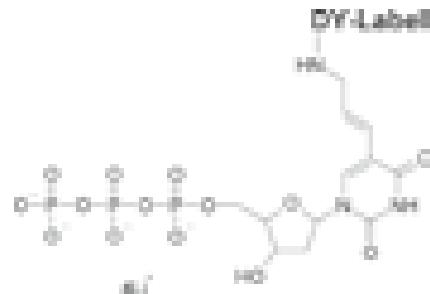
unit size	Product-number
1 mg	010-01
5 mg	010-05
10 mg	010-10



Fluorescently labeled nucleotides

For all our labels aminoallyl-conjugates are available. Please add -34 to the order-number of the dye of your choice.

Unit size	100 µl
Concentration	1 mM aqueous solution
Product-number	XXX-34



Protocols

Recommendations / instructions for the labeling of proteins with NHS-esters

Preface:

Fluorescently labelled antibodies are versatile tools in bioanalytics and medical diagnostics. They find their use, among others, in technologies like fluorescence microscopy, flow cytometry, protein microarrays and microtiter plate based detection schemes. There are many suppliers for labelled antibodies, and even more for unconjugated proteins. Kits for the labelling of proteins are also available from various sources. One crucial issue, especially in academic institutions, is the price of the kits or conjugated proteins which is prohibitive for some experiments. The following instructions want to describe an economic method of labelling proteins with simple recipes and instrumentation which allow a full control of all ongoing processes.

The recipes:

Sodium bicarbonate buffer for dialysis and labelling (10x /500 mM)

Dissolve 21 g (250 mmol) of sodium hydrogen carbonate in 400 ml distilled water. Add 1 g of sodium azide (0.2 per cent, works as bacteriozide). Adjust the pH with a concentrated aqueous solution of sodium hydroxide to pH 8.3. Prior to use, dilute the buffer by adding 9 parts water to one part of the concentrated stock solution (v/v).

100 mM PBS-buffer pH 7.4 for elution and storage

Dissolve 2.19 g sodium dihydrogen phosphate ($\text{NaH}_2\text{PO}_4 \cdot \text{H}_2\text{O}$, 15.9 mmol), 14.97 g disodium hydrogen phosphate ($\text{Na}_2\text{HPO}_4 \cdot 2\text{H}_2\text{O}$, 84.1 mmol), 5.8 g sodium chloride (100 mmol) and 0.32 g sodium azide in 1 l distilled water.

Additional materials and instruments required:

- Protective gloves
- 1 pipet 0.5–10 µl and corresponding tips (white)
- 1 pipet 10–100 µl and corresponding tips (yellow)
- Centrifuge for Eppendorf vials
- Ultrasonification bath (recommended, but not required)
- 1 Pasteur pipet
- Dry DMF (e.g. Fluka P0-no. 40228)
- Sephadex G-25 medium (GE Healthcare)
- Column for size exclusion chromatography (e.g. 5 ml or 10 ml graduated pipettes)
- Silanized glass fiber wadding (e.g. Macherey & Nagel P0-no. 718 002)
- 1 glass bar (at least as long as the column)
- 4 beakers (size can vary)
- 1 tripod with clamp
- Shaker for Eppendorf tubes
- 2 Eppendorf vials (1.5 ml)
- UV-spectrometer

Only when dialysis is required:

- QuixSep Micro dialyzer (Orange Scientific n.v./s.a./ Belgium)
- Dialyzing membrane (also available from Orange Scientific)
- Magnetic stirrer
- Stirring bar

Safety remark:

Always wear protective gloves, lab coat and splash goggles for your personal safety when following the instructions below. In addition, make yourself familiar with the MSDSs which come with the products and chemicals.

Antibody preparation:

Commercially available proteins are usually supplied in aqueous buffer or lyophilised from it. It is essential for a successful labelling reaction that there are no primary or secondary amines present in the reaction solution (otherwise the amine containing buffer components would be labelled instead of the target protein). Thus it is highly recommended to remove buffer components like Tris or ammonia by dialysis against the labelling buffer (50 mM sodium hydrogen carbonate, see recipe above). The QuixSep microdialyzer chamber (see <http://www.orangesci.com/dialqs.htm> for detailed instructions) offers therefore a convenient procedure which is also suitable for low volumes ranging from 100 µl up to 5 ml. Simply fill your protein solution into the dialysis chamber, place a dialysis membrane which was allowed to equilibrate in the labelling buffer for a few minutes on it and fix it by pushing the collar over it. Dialysis can now be done in a 250 ml beaker against the labelling buffer under constant stirring with a magnetic stirrer. Exchange the labelling buffer twice within three hours. Collect the protein solution from the vial by piercing a pipet through the membrane, aspirating the sample and placing it into an Eppendorf vial. Now you have your protein ready for the labelling reaction.

If you are sure that there are no amines present in your protein solution you can also simply add an appropriate volume of the labelling buffer stock solution to your protein to adjust the pH around 8.3.

To enable a batch-to-batch consistency of the degree of labelling, it is necessary to keep the dye to protein ratio constant for each labelling reaction. The following calculation is an example for the use of DY-647-NHS-ester (MW 761.85 g·mol⁻¹, 1 mg) for labelling 1 mg of an IgG (MW 150,000 g·mol⁻¹). Notwithstanding the fact that an optimal degree of labelling depends on the individual application for the conjugate and can require optimisation, a sixfold molar excess of dye for the labelling reaction is a good value to start with.

The molar ratio between 1 mg dye and 1 mg antibody is 197 (150,000 / 761.85). In order to work with volumes lower than 100 µl, 98.5 µl DMF (197 / 2) should be given into the plastic vial containing the 1 mg of DY-647-NHS-ester. To assist proper dissolving of the NHS-ester you can vortex the vial. Alternatively, ultrasonification of the solution for a couple of seconds is also a convenient way to dissolve the reactive dye. It can happen that some solvent and dye is contained in the screw cap of the vial. This material can be recovered by shortly centrifuging the vial.

DMF is superior to the more common DMSO as solvent for reactive dyes for several reasons. It is less hydroscopic, has a lower boiling point (which is important if you want to remove the solvent again from unused dye in a speedvac), and often dyes are better soluble in it. These advantages should override fears about the toxicity of this versatile solvent.

To start the labelling reaction, transfer 3 µl of the NHS-ester solution to the IgG in the reaction buffer. If you can control the concentration of the protein, a concentration of 5 mg/ml is recommended. In our case, the 1 mg IgG should have been dissolved in 200 µl of 50 mM bicarbonate buffer. Lower concentrations lead to a reduced labelling efficiency due to the fact that the hydrolysis of the NHS-ester competes the labelling reaction, and the less protein is present the more the hydrolysis dominates.

Allow the labelling to be done in the Eppendorf vial placed in a shaker over the course of two hours. After this time, the NHS-ester should be either bound to the protein or have been hydrolized which makes the use of a stopping reagent obsolete.

The Sephadex G-25 medium is a classic gel filtration material. Upon addition of elution buffer, it swells by a factor of five in volume. Thus, it is recommended to put at least 6 ml of elution buffer onto 1.2 g Sephadex in a beaker for filling a 5 ml-column. Double the amounts for a 10 ml column. A 5 ml-column is recommended for reaction solution volumes up to 250 µl. Please use a 10 ml-column for volumes up to 1 ml.

To prepare the size-exclusion chromatography column, fix the column vertically in the clamp of the tripod. Put a small amount of the silanized glass fiber wadding into the column and push it with the help of the glass bar to the small opening at the ground. The wadding allows the buffer to flow through it while it holds back the gel.

Now pipet once elution buffer into the column, immediately followed by the swollen gel in elution buffer which had been allowed before to equilibrate for at least one hour. Fill the column entirely with the gel and allow thereby the material to settle. If the Sephadex in the beaker runs dry due to the aspiration of the buffer, simply add some more of the elution buffer to liquidize the gel again. Keep a 1 cm region at the top of the column free of the gel.

Now transfer your coloured labelling solution slowly onto the column and allow it to sink into the gel. Wash your pipet carefully (or use a new one) to remove all traces of dye or protein which would contaminate your elution buffer. After the labelling solution has sunk fully into the column, start to elute the conjugate by slowly adding the elution buffer drop by drop onto the column. You can add bigger portions once there is a dye free zone on top of the column. When the reaction solution is eluted, you will see a separation between the labelled protein which runs ahead as a quite sharp band while the free dye is slowly smearing behind.

Once the conjugate arrives at the bottom of the column collect it in an Eppendorf vial. While selecting the size of the vial keep in mind that the volume of the protein conjugate solution increases by a factor of ca. 2.5 on the column. Dispose the used Sephadex gel after collecting the conjugate and wash the column. It would be very time-consuming to wash the free dye from the gel, and traces of it could impurify the next conjugate. It is really not worth the time, and the price for 1 g Sephadex is less than €3.

To characterize your protein conjugate solution it is recommended to record an absorption spectrum of it over the full wavelength range covering both the protein absorption at 280 nm as well as the maximum wavelength of the dye together with the shape of the dye's absorption band. Avoid thereby absorbance values exceeding "2" because such values are hard to interpret. Use cuvettes with smaller pathlengths instead or dilute a small fraction of your conjugate solution appropriately with the elution buffer, e.g. by a factor of 20.

A dye to protein (D/P) ratio can be calculated according to a formula derived from the combination of the Lambert-Beer law for 280 nm and the absorption maximum of the dye. A high uncertainty in this formula is the molar absorbance of the dye at its absorption maximum. Usually, the molar absorbance is determined in organic solvents because this parameter is concentration dependent in aqueous solution.

Thus, one can not be sure if the calculated D/P ratio reflects the experimental truth. The absorption spectrum over the full UV-Vis range is a more reliable analytical information and should be used to compare conjugates among each other.

The 95.5 µl of the DY-647-NHS-ester which remain from the reaction should be stored at -20°C. An exact shelf life of this solution can not be given. The most crucial factor for the stability of the NHS-ester is to protect it from humidity/moisture. Thus, allow the frozen vial to come completely to room temperature prior to opening it again. A simple test for the remaining activity of the dissolved reactive dye could be to couple it to a small amount (0.5 mg) of a cheap protein like bovine serum albumin (or, in times of BSE, streptavidin) according to the instructions given above when using the dye the first time. Later on, the coupling can be repeated, and the absorption spectrum of the conjugate gives an information about how the NHS-ester degraded over the storage time. In correspondence, the amount of NHS-ester can be adjusted for the labelling reaction with the desired protein.

When labelling proteins with colourless haptens like biotin-NHS-ester, a first coupling with a coloured dye and a cheap protein should be performed on the column whereby the volume of the eluent which leaves the column prior to the conjugate should be measured. Assuming that the elution times for proteins are nearly identical (an allowable assumption since it is about size exclusion chromatography), the same dead volume can be disposed prior to collecting the desired colourless conjugate in a second run on the same column with fresh Sephadex gel.

Services

Aside from our standard product portfolio we offer the customizing of fluorescent labels as well as the synthesis of conjugates on demand. Together with reliable partners we can supply fluorescently labeled DNA-oligonucleotides, fluorogenic, peptide-based enzyme substrates and fluorescently labeled small molecules such as β -estradiol or nucleotides. Feel free to contact Dyomics for a quotation. We are sure that our cost/performance ratio will convince you.

Recommended literature

DY-681 and DY-781 labeled primers

Ritz C.M., Wissman V., Microsatellite Analyses of Artificial and Spontaneous Dogrose Hybrids Reveal the Hybridogenic Origin of Rosa micrantha by the Contribution of Unreduced Gametes, *Journal of Heredity*, doi:10.1093/jhered/esq124.

Review

Napp J., Mathejczyk J.E., Alves F., Optical imaging *in vivo* with a focus on paediatric disease: technical progress, current preclinical and clinical applications and future perspectives, *Pediatr Radiol*, (2011) DOI 10.1007/s00247-010-1907-0.

DY-682-labeled secondary antibodies

Koplin A., Preissler S., Ilina Y., Koch M., Scior A., Erhardt M., Deuerling E., A dual function for chaperones SSB-RAC and the NAC nascent polypeptide-associated complex on ribosomes, *The Journal of Cell Biology*, Vol. 189(2010)57–68.

DY-682 and DY-800 labeled secondary antibodies

Ferreira D.N., Katayama I.A., Oliveira I.B., Rosa K.T., Furukawa L.N.S., Coelho M.S., Casarini D.E., Heimann J.C., Salt-Induced Cardiac Hypertrophy and Interstitial Fibrosis Are Due to a Blood Pressure-Independent Mechanism in Wistar Rats, *The Journal of Nutrition*, 140(2010)1742-51.

DY-601XL, DY-613, V07-04118 and V07-04146 as stains in acid media

Brockmann S., Arnold T., Schweder B., Bernhard G., Visualizing Acidophilic Microorganisms in Biofilm Communities Using Acid Stable Fluorescence Dyes, *Journal of Fluorescence*, 20(2010)943-51.

DY-520XL in TIRF microscopy

Spendier K., Carroll-Portillo A., Lidke K.A., Wilson B.S., Timlin J.A., Thomas J.L., Distribution and Dynamics of Rat Basophilic Leukemia Immunoglobulin E Receptors (Fc3RI) on Planar Ligand-Presenting Surfaces, *Biophysical Journal*, 99(2010)388-397

DY-549 and DY-649 on protein microarrays

Schröder C., Jacob A., Tonack S., Radon T.P., Sill M., Zucknick M., Rüffer S., Costello E., Neoptolemos J.P., Crnogorac-Jurcevic T., Bauer A., Fellenberg K., Hoheisel J.D., Dual-color Proteomic Profiling of Complex Samples with a Microarray of 810 Cancer-related Antibodies, *Mol Cell Proteomics*, 9(2010)1271-1280.

Review

Paganin-Gioanni A., Bellard E., Paquereau L., Ecochard V., Golzio M., Teissié J., Fluorescence imaging agents in cancerology, *Radiol Oncol* 2010; 44(3): 142-148.

DY-520XL in multicolour microscopy

Carroll-Portillo A., Spendier K., Pfeiffer J., Griffiths G., Li H., Lidke K.A., Oliver J.M., Lidke D.S., Thomas J.L., Wilson B.S., Timlin J.A., Formation of a Mast Cell Synapse: FcRI Membrane Dynamics upon Binding Mobile or Immobilized Ligands on Surfaces, *The Journal of Immunology*, 184(2010)1328-1338.

DY-752 as imaging agent

Pöschinger T., Janunts E., Brünner H., Langenbucher A., CCD-based projectional imaging of fluorescent probes in tissue-like media: experimental setup and characterization, *Zeitschrift für Medizinische Physik*, 2010, in press.

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DY-636 on labeled particles

Jiang X., Weise1 S., Hafner M., Röcker C., Zhang F., Parak W.J., Nienhaus G.U., Quantitative analysis of the protein corona on FePt nanoparticles formed by transferrin binding, *J. R. Soc. Interface* 6 February 2010 vol. 7 no. Suppl 1 S5-S13.

DY-682 in tumor imaging

Gong H., Kovar J., Little G., Chen H., Olive D.M., In Vivo Imaging of Xenograft Tumors Using an Epidermal Growth Factor Receptor-Specific Affibody Molecule Labeled with a Near-infrared Fluorophore, *Neoplasia*. 2010 February; 12(2): 139-149.

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Owczarek-Lipska M., Dolf G., Guziewicz K.E., Leeb T., Schelling C., Posthaus H. Braunschweig M.H., Bovine cardiac troponin I gene (TNNI3) as a candidate gene for bovine dilated cardiomyopathy, *Archiv Tierzucht*, 52(2009)113-123.

DY-781-dUTP in PCR application

Braunschweig M.H., Quantification of global DNA methylation with infrared fluorescence in liver and muscle tissues of differentially fed boars, *Luminescence*, 24(2009)213-216.

DY-633-HaloTag

Kosaka N., Ogawa M., Choyke P.L., Karassina N., Corona C., McDougall M., Lynch D., Hoyt C., Levenson R., Los G.V., Kobayashi H., In Vivo Stable Tumor-Specific Painting in Various Colors Using Dehalogenase-Based Protein-Tag Fluorescent Ligands, *Bioconjug Chem.*, 20(2009)1367-1374.

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Vollmer T., Kleesiek K., Dreier J., Lipopolysaccharide-binding protein (LBP) gene polymorphisms: Rapid genotyping by real-time PCR and association with infective endocarditis, *Clinical Biochemistry* 42 (2009) 1413-1419.

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Pola R., Studenovsky M., Pechar M., Ulbrich K., Hovorka O., Vetvicka D., Rihova B., HPMA-copolymer conjugates targeted to tumor endothelium using synthetic oligopeptides, *Journal of Drug Targeting*, 2009.

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Mindt T.L., Jungi V., Wyss S., Friedli A., Pla G., Novak-Hofer I., Grünberg J., Schibli R., Modification of Different IgG1 Antibodies via Glutamine and Lysine using Bacterial and Human Tissue Transglutaminase, *Bioconjugate Chem.*, 19(2008)271-278.

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DY-676 for in vivo imaging

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Lisy M.R., Goermer A., Thomas C., Pauli J., Resch-Genger U., Kaiser W.A., Hilger I., In Vivo Near-infrared Fluorescence Imaging of Carcinoembryonic Antigen-expressing Tumor Cells in Mice, *Radiology*, 247(2008)779-787.

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Riederer B.M., Non-covalent and covalent protein labeling in two-dimensional gel electrophoresis, *Journal of Proteomics*, 71(2008)231-244.

DY-547-Streptavidin

Mora J.R., Zielinsky T.L., Nelson B.P., Getts R.C., Protein detection enhanced by 3DNA dendrimer signal amplification, *BioTechniques*, 44(2008)815-818.

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Kusnezow W., Banzon V., Schröder C., Schaal R., Hoheisel J.D., Rüffer S., Luft P., Duschl A., Syagailo Y.V., Antibody microarray-based profiling of complex specimens: systematic evaluation of labeling strategies, *Proteomics*, 7(2007)1786-1799.

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Acknowledgement

Dyomics wants to thank the following scientists and partners for contributing impressive images and pictures to this catalogue:

Monika Bäumle, Fluka AG, Buchs, Switzerland

Anette Eitner & Reimar Krieg, University of Jena, Jena Germany

Ulrich Hassiepen, Novartis Institutes for BioMedical Research, Basel, Switzerland

Ingrid Hilger, Institute for Diagnostic and Interventional Radiology, Jena, Germany

Sandra Orthaus, Institute of Molecular Biotechnology, Jena, Germany

Katharina Pachmann, KIM II, Jena, Germany

Lana Popova, Cell Signaling Inc., Danver, MA, USA

Beat Riederer, Institut de Biologie Cellulaire et de Morphologie, Lausanne, Switzerland

Jens Sobek, Functional Genomics Center Zurich, Switzerland

Kai Sohn, Fraunhofer IGB, Stuttgart, Germany

Applied Spectral Imaging, Ltd., Migdal Ha'Emek, Israel

Covalys Biosciences AG, Witterswil, Switzerland

SIRS-Lab GmbH, Jena, Germany

unique-m.o.d.e. AG, Jena, Germany

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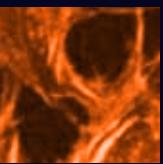
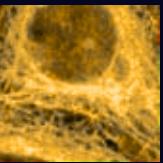
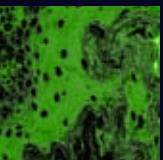
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