

2 Carotenes

β,β -Carotene

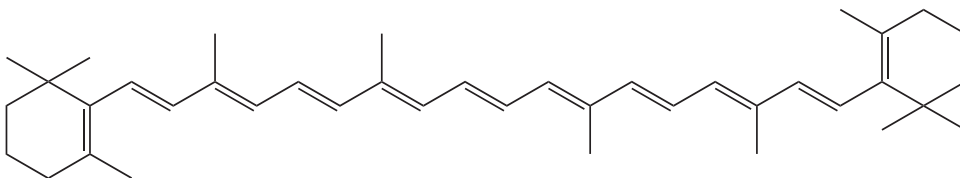
IUPAC: β,β -Carotene

Molecular formula: $C_{40}H_{56}$

Recommended abbreviation: $\beta\beta$ -Car ($\beta\beta$)

(trivial name: β -Carotene)

Molecular weight: 536.87



Biological occurrence

Dominant pigment in chlorophytes, prasinophytes, mesostigmatophytes, rhodophytes and one group of dinoflagellates. Minor in all other algal groups. Also present in plants (notably carrots) [78]

Source culture

Pavlova lutheri (Pavlovophyceae)

Alteration products

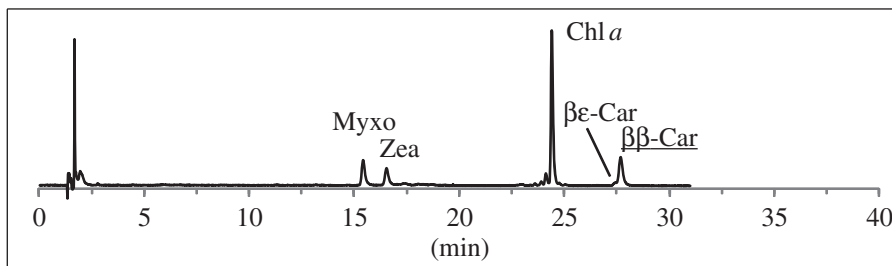
Cis-isomers

Biosynthetically related to

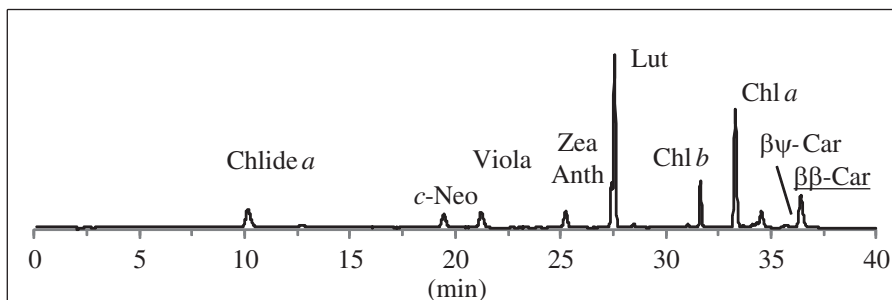
Cantha, Asta, Zea, Viola, Neo, Diato, Diadino, Allo etc.

Occurs together with

HPLC chromatogram of *Synechococcus* sp. (system 1)

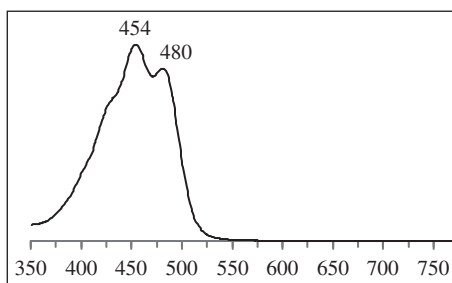
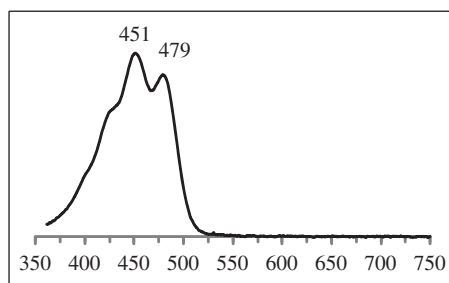
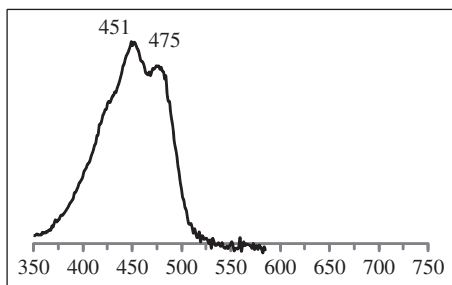
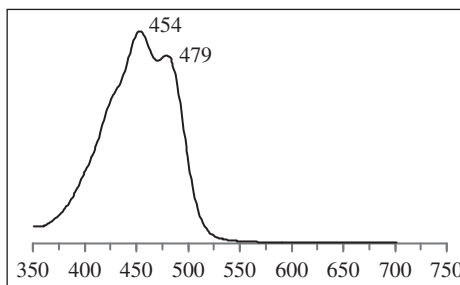


HPLC chromatogram of *Dunaliella salina* (system 2)



UV-Vis spectra (see also reference spectra below)

Solvent	λ_{max} (nm)	Band ratio (% III:II)	Ref.
Acetone	(427), 454, 480	21	[96]
Diethyl ether	(430), 447, 476	5	[133]
Ethanol (see Remarks)	(428), 451, 480	27	[86]
Hexane	(425), 451, 477	29	[96]
Methanol (see Remarks)	(429), 449, 475	25	[160]
Recommended specific absorption coefficient d (L g ⁻¹ cm ⁻¹)		259 (at 453 nm, hexane) [103] 250 (at 454 nm, acetone) [96]	

Reference spectra**In acetone****In hexane****In HPLC solvent system 1****In HPLC solvent system 2****Mass spectra**

Ionization technique	Mass analyser type	Diagnostic ions (m/z, rel. intensity)	Ref.
EI	Magnetic sector	536 [M] ⁺ (44), 444 [M-92] ⁺ (15), 430 [M-106] ⁺ (4), 119 (72), 109 (22), 91 (45), 83 (23), 69 (100)	[144]
Remarks	To aid dissolving in alcohol, first dissolve in a drop or two of hexane		

β,ϵ -Carotene

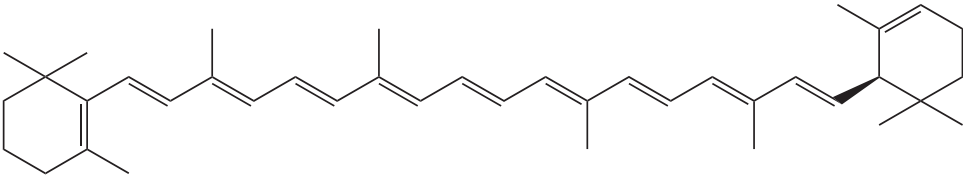
IUPAC: (6'*R*)- β,ϵ -Carotene

Molecular formula: C₄₀H₅₆

Recommended abbreviation: $\beta\epsilon$ -Car ($\beta\epsilon$)

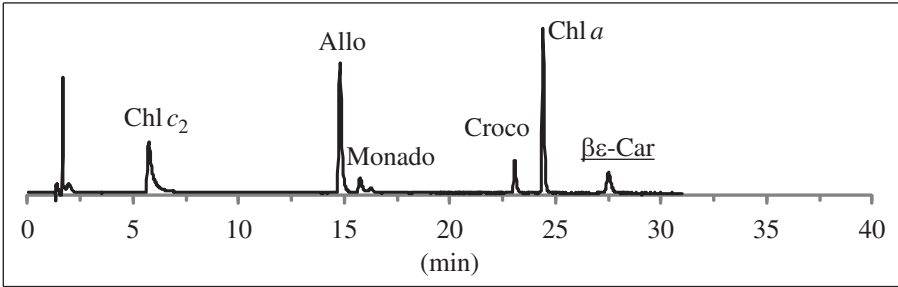
(trivial name: α -carotene)

Molecular weight: 536.87

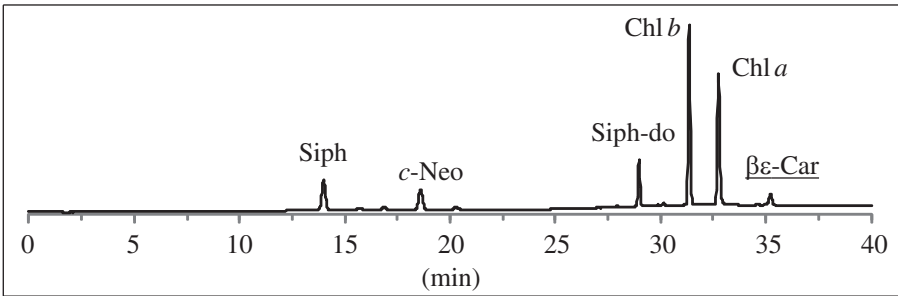


Biological occurrence	Minor or trace pigment in chlorophytes, prasinophytes, cryptophytes, some dinoflagellates, cyanobacteria
Source culture	<i>Chroomonas salina</i> (cryptomonad)
Alteration products	<i>Cis</i> -isomers
Biosynthetically related to	Lut, Loro, Siph
Occurs together with	Lut, Zea, Viola, Anth, Neo, Allo, Croco, Monado

HPLC chromatogram of *Chroomonas salina* (system 1)

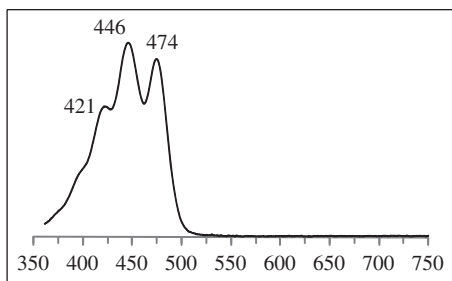


HPLC chromatogram of *Codium fragile* (system 2)

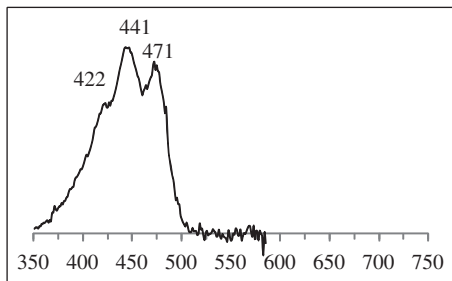
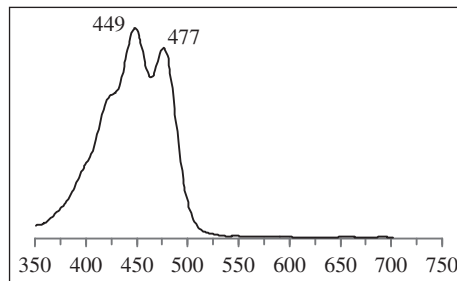


UV-Vis spectra (see also reference spectra below)

Solvent	$\lambda_{\max}$ (nm)	Band ratio (% III:II)	Ref.
Acetone	(424), 448, 476	56	[96]
Ethanol (see remarks)	423, 444, 473	61	[86]
Hexane	423, 446, 474	71	[57]
Methanol (see remarks)	(422), 442, 471	61	[160]
Recommended specific absorption coefficient		270 (at 446 nm, hexane) [57]	
d (L g⁻¹ cm⁻¹)		270 (at 448 nm, acetone) [96]	

Reference spectra**In hexane**

For spectrum in acetone, see [109]

In HPLC solvent system 1**In HPLC solvent system 2****Mass spectra**

Ionization technique	Mass analyser type	Diagnostic ions (m/z, rel. intensity)	Ref.
EI	Magnetic sector	536 [M] ⁺ (100), 480 [M-56] ⁺ (2), 444 [M-92] ⁺ (25), 430 [M-106] ⁺ (18), 388, 378, 374	[144]
Remarks	To aid dissolving, add a drop or two of hexane before adding the alcohol		

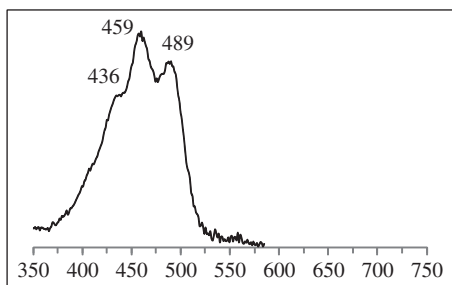
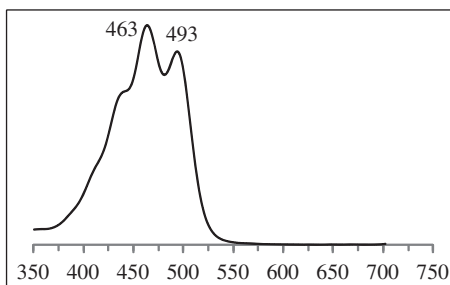
UV-Vis spectra (see also reference spectra below)

Solvent	λ_{max} (nm)	Band ratio (% III:II)	Ref.
Acetone	(439), 461, 491	n.d.	[66]
Cyclohexane	439, 464, 496	59	[69]
Ethanol (see remarks)	(440), 460, 489	23	[83]
Recommended specific absorption coefficient d (L g ⁻¹ cm ⁻¹)		276 (at 462 nm, hexane) [181] 319 (at 459 nm, petroleum ether) [126]	

Reference spectra

For spectrum in acetone, see [109]

For spectrum in hexane, see [109]

In HPLC solvent system 1**In HPLC solvent system 2****Mass spectra**

Ionization technique	Mass analyser type	Diagnostic ions (m/z, rel. intensity)	Ref.
EI	Magnetic sector	536 [M] ⁺ (100), 467 [M-69] ⁺ (8), 444 [M-92] ⁺ (20), 430 [M-106] ⁺ (25), 407 (m*, M → M-69)	[30]
Remarks	To aid dissolving, add a drop or two of hexane before adding the alcohol		

ϵ,ϵ -Carotene

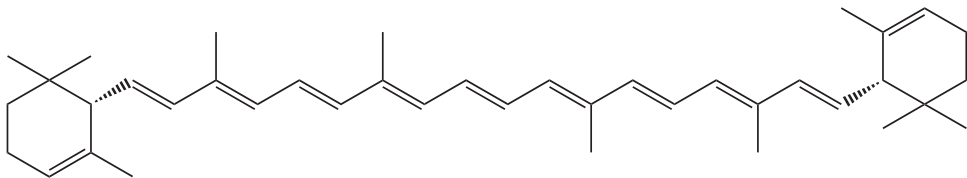
IUPAC: (6*S*,6'*S*)- ϵ,ϵ -Carotene

Molecular formula: C₄₀H₅₆

Recommended abbreviation: $\epsilon\epsilon$ -Car (ϵ)

(trivial name: ϵ -carotene)

Molecular weight: 536.87

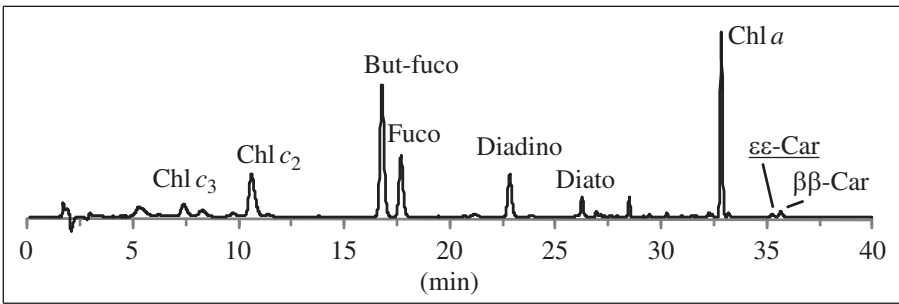


Biological occurrence	Significant pigment in pelagophytes but not always present (see Chapter 1). Occasional traces in cultures of chlorophytes, diatoms, cryptomonads and prasinophytes Pigment Type 3
Source culture	<i>Pelagococcus subviridis</i> (chrysophyte)
Alteration products	<i>Cis</i> -isomers
Biosynthetically related to	Lyc
Occurs together with	

HPLC chromatogram (system 1)

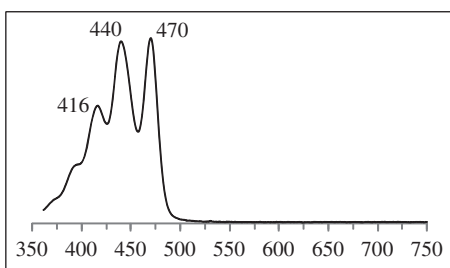
NO DATA AVAILABLE

HPLC chromatogram of *Pelagomonas calceolata* (system 2)

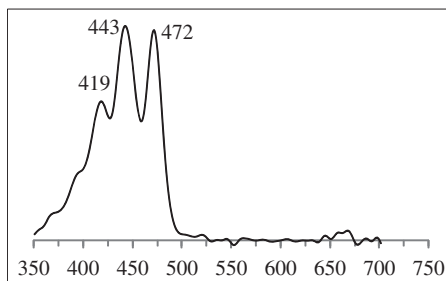


UV-Vis spectra (see also reference spectra below)

Solvent	λ_{max} (nm)	Band ratio (% III:II)	Ref.
Cyclohexane	416, 440, 470	n.d.	[157]
95% Ethanol	417, 440, 470	96	[36]
Hexane	415, 439, 469	101	[23]
Recommended specific absorption coefficient d (L g ⁻¹ cm ⁻¹)		312 (at 440 nm, petroleum ether) [143]	

Reference spectra**In hexane****In HPLC solvent system 1**

NO DATA AVAILABLE

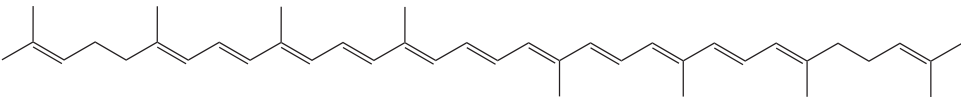
In HPLC solvent system 2**Mass spectra**

Ionization technique	Mass analyser type	Diagnostic ions (m/z, rel. intensity)	Ref.
EI	Magnetic sector	536 [M] ⁺ (100), 480 [M-56] ⁺ (5), 444 [M-92] ⁺ (28), 430 [M-106] ⁺ (8), 388 [M-92-56] ⁺ (30)	[143]
Remarks		Opposite stereochemistry than βε-Car and Lut [23, 46]	

ψ,ψ-Carotene (Lycopene)
IUPAC: ψ,ψ-Carotene
Molecular formula: C₄₀H₅₆

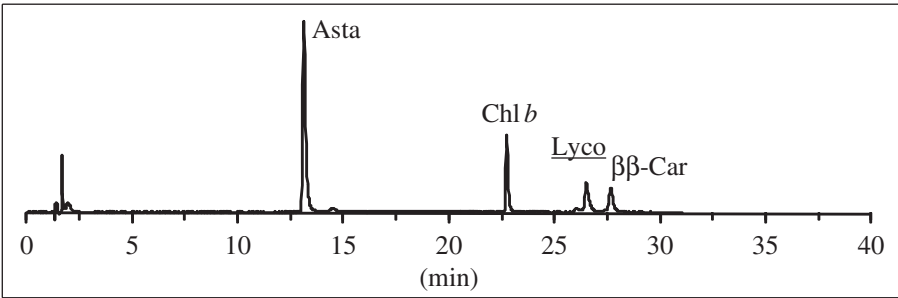
Recommended abbreviation: Lyco (Ly)

Molecular weight: 536.87



Biological occurrence	Minor or trace pigment in mesostigmatophytes (also characteristic of tomatoes)
Source culture	<i>Mesostigma viride</i> (mesostigmatophyte)
Alteration products	<i>Cis</i> -isomers
Biosynthetically related to	The biosynthetic precursor to all carotenoids
Occurs together with	

HPLC chromatogram of mixed standards (system 1)

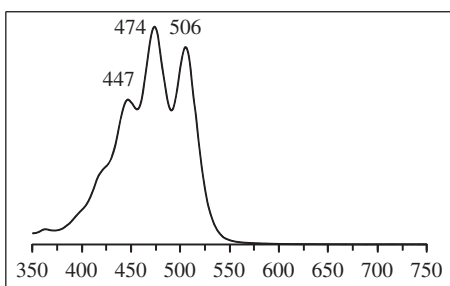
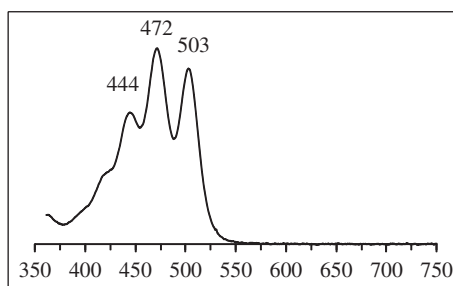
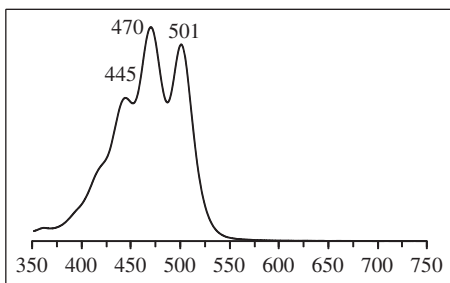


HPLC chromatogram (system 2)

NO DATA AVAILABLE

UV-Vis spectra (see also reference spectra below)

Solvent	λ_{max} (nm)	Band ratio (% III:II)	Ref.
Acetone	448, 474, 506	84	[2]
Ethanol	443, 472, 502	n.d.	[113]
Hexane	448, 473, 504	n.d.	[14]
Recommended specific absorption coefficient d (L g ⁻¹ cm ⁻¹)		347 (at 473 nm, hexane) [181] 345 (at 474 nm, acetone) [2]	

Reference spectra**In acetone****In hexane****In HPLC solvent system 1****In HPLC solvent system 2**

NO DATA AVAILABLE

Mass spectra

Ionization technique	Mass analyser type	Diagnostic ions (m/z, rel. intensity)	Ref.
EI	Magnetic sector	536 [M] ⁺ (22), 467 [M-69] ⁺ (4), 444 [M-92] ⁺ (2), 430 [M-106] ⁺ (4), 109 (19), 91 (47), 69 (100)	[55]
Remarks	Mesostigmatophytes are found in freshwater environments		