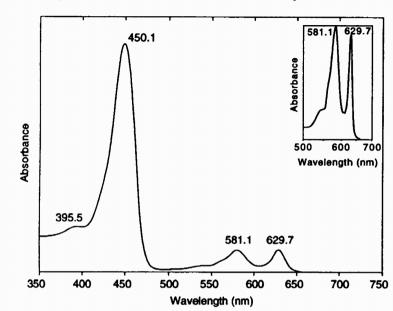


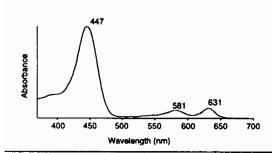
HPLC peak 6a

Standard spectrum in reference solvent: diethyl ether

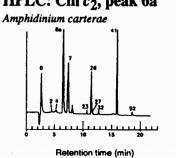


Molecular structure

Diode array spectrum in SCOR eluant



HPLC: Chl c2, peak 6a



Chlorophyll c_2

Property		Data		
Name:	(Trivial) (IUPAC)	Chlorophyll c ₂ 3 ¹ ,3 ² ,8 ¹ ,8 ² ,17 ¹ ,17 ² -hexadehydro-13 ² - methoxycarbonyl-phytoporphyrinato-Mg(II); see Hynninen (1991)		
SCOR abbreviation:		Chl c ₂		
Occurrence:		Chromophyte algae, brown seaweeds (Jeffrey, 1976, 1989)		
Colour:		Light green on TLC; emerald green (concentrated solution)		
Molecular formula:		$C_{35}H_{28}N_4O_5Mg$		
Molecular weight:		608.94		
Specific extinction coefficient: † α (1 g ⁻¹ cm ⁻¹)		374 (at 443.8 nm in 90% acetone + 1% pyridine) 40.4 (at 630.9 nm in 90% acetone + 1% pyridine) 37.2 (at 629.6 nm in 100% acetone + 1% pyridine) Jeffrey (1972)		

[†] Where chlorophylls $c_1 \& c_2$ are not resolved, the mean of the extinction coefficients was used. In 90% acetone, $\alpha = (44.8 + 40.4)/2 = 42.6 \text{ 1 gm}^{-1} \text{ cm}^{-1}$.

Molar extinction coefficient: ε (1 mol⁻¹ cm⁻¹)

227.7x10³ (at 443.8 nm in 90% acetone + 1% pyridine) 24.6x10³ (at 630.9 nm in 90% acetone + 1% pyridine) 22.7x10³ (at 629.6 nm in 100% acetone + 1% pyridine)

*Soret (blue maximum): red ratio

Calculated from a above

UV-vis spectra:

Fluorescence spectra:

Solvent	Absorbance m	axima (nm)	Band ratio*	Reference
100% Acetone	444.6 581.3	629.6	8.62	Jeffrey (1972)
Diethyl ether	450.1 581.1	629.7	10.35	SCOR WG 78 data
HPLC Eluant	447 581	631	8.48	SCOR WG 78: Wright et al. (1991) method

Solvent	Excitation (nm)	Emission (nm)	Reference
100% Acetone	453	635, 696	Jeffrey (1972)
100% Ethanol	450	638, 694	Jeffrey (1972)

Corresponding phytoporphyrin **Alteration products:** Culture from which SCOR Amphidinium carterae (dinoflagellate) data were obtained:

Jeffrey (1989); Scheer (1991) Additional reference(s):