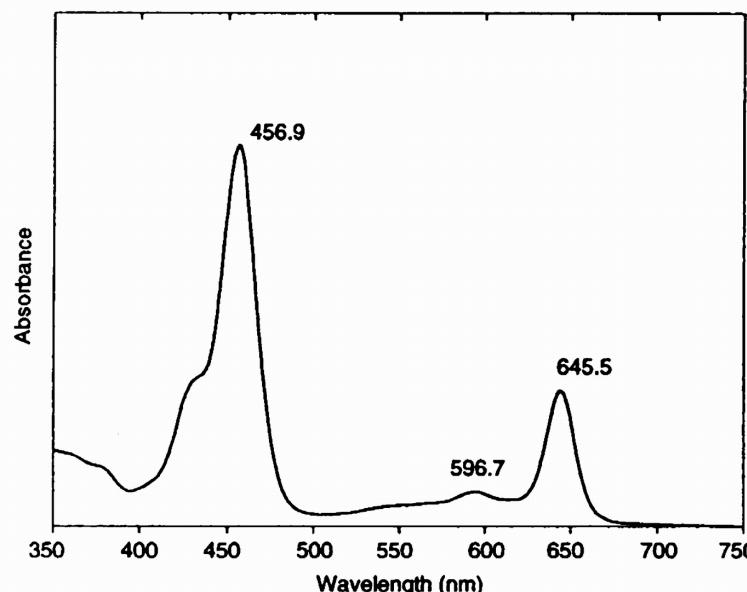


# Chlorophyll b

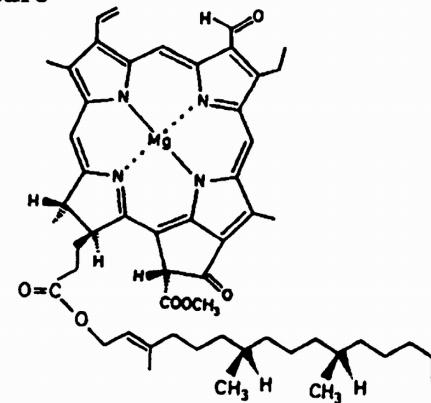
## HPLC peak 37

# Chlorophyll b

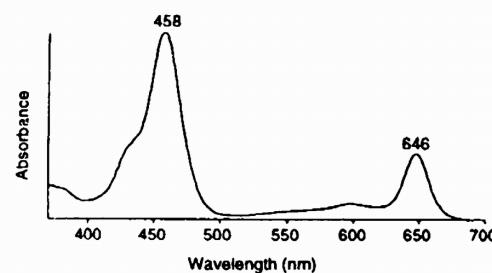
Standard spectrum in reference solvent: acetone (100%)



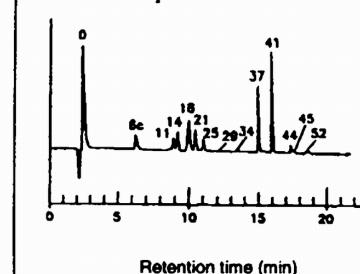
### Molecular structure



### Diode array spectrum in SCOR eluant



### HPLC: Chl b, peak 37 *Micromonas pusilla*



### Property

### Data

Name:	(Trivial) (IUPAC)	Chlorophyll b
		Trivial name sufficient, see Hynninen (1991)
SCOR abbreviation:	Chl b	
Occurrence:	Higher plants, green algae, symbiotic prochlorophytes	
Colour:	Olive green (TLC); emerald green (concentrated solution)	
Molecular formula:	C <sub>55</sub> H <sub>70</sub> N <sub>4</sub> O <sub>6</sub> Mg	
Molecular weight:	907.49	
Specific extinction coefficient: $\alpha$ (1 g <sup>-1</sup> cm <sup>-1</sup> )	51.36 (at 646.8 nm in 90% acetone) 62.00 (at 643.3 nm in diethyl ether) Jeffrey & Humphrey (1975)	
Molar extinction coefficient: $\epsilon$ (1 mol <sup>-1</sup> cm <sup>-1</sup> )	46.61 x 10 <sup>3</sup> (at 646.8 nm in 90% acetone) 56.26 x 10 <sup>3</sup> (at 643.3 nm in diethyl ether) Calculated from $\alpha$ above	
UV-vis spectra:		

Solvent	Absorbance maxima ( nm)			Band ratio*	Reference
100% Acetone	456.9	596.7	645.5	2.82	SCOR WG 78 data
Diethyl ether	455	595	644	2.81	Smith & Benitez (1955)
Diethyl ether	453	593	642	2.77	Hynninen & Lötjönen (1983)
Ethanol	464		649		Griffiths (1991)
HPLC Eluant	458	596	646	2.83	SCOR WG 78: Wright <i>et al.</i> (1991) method

### Fluorescence spectra:

\*Soret (blue maximum): red ratio

Solvent	Excitation ( nm)	Emission ( nm)	Reference
Diethyl ether	453	646	Boardman & Thorne (1971)
100% Acetone	459	652, 710	Jeffrey (1972)
100% Acetone	453	652	SCOR WG 78 data

### Alteration products:

Chlorophyllide b, pheophytin b, pheophorbide b, epimers, allomers, pyro-pheophytin b

### Culture from which SCOR data were obtained:

*Dunaliella tertiolecta* (chlorophyte),  
*Micromonas pusilla* (prasinophyte)

### Additional reference(s):

Scheer (1991)