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Original article

Volatile profile and sensory descriptors of *Melaleuca quinquenervia* (Cav.) S.T. Blake essential oil from Madagascar as a promising food flavouring

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Summary The present study investigated the volatile and sensory profiles and antioxidant activity of 1,8-cineole type Niaouli essential oil from Madagascar for which limited information is reported in literature. The volatile compounds were identified by Gas chromatography–mass spectrometry and semi-quantitative results were obtained by Gas chromatography with a flame-ionisation detector; the odour descriptors were determined by a trained panel using a Quantitative Descriptive Analysis. The antioxidant activity was determined, too. Several volatiles, most of these here firstly, have been identified; 1,8-cineole (eucalyptus-like) prevailed, followed by viridiflorol and (*E*)-nerolidol. 'Eucalyptus-like, fresh-mint, woody, sweet-floral and spicy' odour descriptors have been defined. Volatiles resulted in agreement with the odour descriptors. The data allowed to characterise the Niaouli essential oil; considering its volatile composition, pleasant sensory notes and antioxidant activity the 1,8-cineole type Niaouli essential oil from Madagascar could be considered a promising flavouring agent and natural preservative in the food industry.

Keywords 1,8-cineole chemotype, antioxidant activity, *Melaleuca quinquenervia* (Cav.) S.T. Blake, Niaouli essential oil, sensory descriptors, volatile composition.

Introduction

The Myrtaceae family, mainly spread in Australia, Southeast Asia, tropical to southern temperate America and Africa, is a rich source of essential oils, with a hundred genera and approximately 3800 species (Wilson *et al.*, 2001). Belonging to the Myrtaceae family, Melaleucas include a diverse range of forms, from woody shrubs to tall timber-producing trees. Leaves vary in size and shape, from short to long, elliptic, cordate and lanceolate; venation is pinnate, longitudinal or longitudinal-pinnate (Brophy *et al.*, 2013). The leaf essential oils of various *Melaleuca* species differ in composition, also due to pedoclimatic conditions; methyl eugenol, (*E*)-methyl isoeugenol, 1,8-cineole, terpinen-4-ol, terpinolene, (*E*)-nerolidol, viridiflorol, ledol, β-caryophyllene have been reported as the main volatile compounds (Wheeler *et al.*, 2007).

Melaleuca quinquenervia (Cav.) S.T. Blake is native and mainly spread in Australia, New Caledonia and Madagascar; it is highly valued for its leaf essential oil named Niaouli oil or 'broad-leaf paperbarks'. Several chemotypes have been described in Madagascar: a chemotype rich in 1,8-cineole; a chemotype rich in (*E*)-nerolidol; a chemotype rich in 1,8-cineole, viridiflorol and terpinolene and a chemotype rich in viridiflorol. Niaouli oil with a high content of 1,8-cineole is the most widespread and appreciated chemotype (Ramanoelina *et al.*, 2008).

The oil has been used in traditional medicine against many pathological conditions and recent studies have demonstrated different pharmacological effects (Nam *et al.*, 2008; Acha *et al.*, 2019). The global market for Niaouli oil has experienced significant growth in recent years mainly due to the increasing consumer interest in natural and holistic health products.

Recently the high antimicrobial and antioxidant activities of niaouli oil as an alternative to chemical preservatives in the food industry have been demonstrated; it has been indicated as an innovative agent for shelf-life prolonging of different types of food products, such as bread, carrot and potato (Valková *et al.*, 2022) due to its antioxidant properties, pleasant flavour and sensory notes.

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To be used as food flavouring, the knowledge of the oil volatile profile is required; this can be obtained using human senses and at the same time, high-resolution advanced analytical techniques. The application of advanced analytical techniques in the volatile studies of essential oils is of great importance since they allow the determination of volatiles even present in trace amounts in a complex mixture.

As regards the Niaouli essential oil, 1,8-cineole chemotype produced in Madagascar different studies have been carried out on its volatile composition even if this dates back to more than 10 years ago (Ramanoelina *et al*., 1992, 1994, 2005, 2008); moreover, no information is present in the literature on its sensory features, neither on their correlation with the volatile compounds. The composite characteristics of the Niaouli oil are not defined by the International Standardization Organization such happens for other *Melaleuca* varieties, such as *Melaleuca alternifolia* (tea tree oil) terpinen-4-ol type (ISO 4730: 2017).

Considering what was above reported, this research aims to characterise the volatile composition and sensory descriptors of 1,8-cineole chemotype Niaouli essential oil from Madagascar, which represents the most commercially widespread, providing useful elements for its wide application as a flavour ingredient in food products. Advanced analytical techniques such as GC-MS and sensory qualitative descriptive analysis were applied for the essential oil characterisation.

Materials and methods

Materials

Eight samples of *Melaleuca quinquenervia* (Cav.) S.T. Blake essential oil 1,8-cineole type produced in Madagascar, supplied by commercial retailers with attestation of genuineness, have been analysed. According to the producers' information, the essential oils were obtained by hydrodistillation at atmospheric pressure for 3 h, using a Clevenger-type apparatus, and the distilled oils were collected and dried over anhydrous sodium sulphate, filtered snf stored in amber flash. The samples have been analysed in triplicate by Gas chromatography–mass spectrometry (GC-MS) for volatile identification and by Gas chromatography with a flame-ionisation detector (GC-FID) for the quantitative results.

Gas chromatography–mass spectrometry (GC-MS) analysis

The GC-MS analysis was performed using a GC2010 Plus (Shimadzu) gas chromatograph equipped with an SLB-5 ms capillary column (30 m length, 0.25 mm ID, 0.25 μm film thickness), interfaced with a TQMS 8040 triple quadrupole mass spectrometer (Shimadzu). The oven temperature was programmed from 60 °C to 110 °C at a rate of 2 °C/min, from 110 °C to 160 °C at a rate of 3 °C/min and from 160 °C to 260 °C at a rate of 10 °C/min. Helium was used as a carrier gas at a constant flow rate of 0.8 mL/min. The split ratio was 1:200, the injection temperature was set to 260 °C and the transfer line temperature was 250 °C. Mass selective detector operated in electron impact ionisation (EI) mode at 70 eV electron energy, the ion source temperature was set to 230 °C, and the quadrupole temperature was 150 °C. The mass scan range was 40–360 m/z, scan speed of 1250. One microliter of the essential oil, dissolved in hexane (1:100 v/v), was injected. The peak identification has been carried out using mass spectral data, NIST '23' (NIST/EPA/NIH Mass Spectra Library, USA), FFNSC 3.0 database and linear retention indices (LRI) by comparing with those reported in the literature. Linear retention indices were calculated according to the Van den Dool and Kratz equation based on homologue n-alkane hydrocarbons C7–C30 injected under the same operating conditions.

Gas chromatography with flame-ionisation detector (GC-FID) analysis

The GC-FID analysis was performed using a GC2010 Plus (Shimadzu) gas chromatograph equipped with a flame-ionisation detector (FID) set at 290 °C under the same analytical conditions as above reported. The quantification of the compounds was obtained by the internal normalisation method with a response factor equal to unity for all of the sample constituents.

Sensory analysis

A Quantitative Descriptive Analysis (QDA) has been used to identify and quantify the sensory odour descriptors of the Niaouli essential oils. A sensory panel of 12 judges (6 males and 6 females; age range 24–62 years old) has been trained according to the ISO 8586: 2012 at the Aroma and Sensory Laboratory of the Department of Veterinary Sciences of the University of Messina. The members of the panel were selected concerning their experiences in the sensory evaluation of essential oils. The panel participated in eight training preliminary sessions (2-h each, over 2 weeks), the panellists were asked to list the perceived odour descriptors by sniffing a test strip (dipped in the essential oil around 1 cm). During the training sessions, all the samples were evaluated at least twice to generate appropriate descriptors. The final scorecard for odour analysis contained the descriptors selected by at least one-third of the panel. According to this, five attributes were selected and defined by consensus:

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'eucalyptus-like, fresh-minty, woody, sweet-floral and spicy' as reported in Table S1. The samples were evaluated in a sensory booth room at a temperature of 20 °C under fluorescent lighting equivalent to the daylight. Each judge was located in an individual booth during each sensory evaluation session. Each of the eight samples of Niaouli essential oils has been evaluated randomly three times over six sessions over 3 weeks. For each sample oil, three sniffs from the test strip were performed and fresh air was breathed between each sample. The panellists were asked to mark the intensity of the descriptors on a 9 cm line scale length. For each sample, the scorecards were decoded, and the mean values of the descriptors were calculated.

DPPH radical scavenging assay

The antioxidant activity of the Niaouli essential oil, was assessed by measuring its ability to scavenge the DPPH stable radical. Different concentrations (20–100 μL/mL) of each essential oil were prepared in methanol. Three mL of methanolic solution of DPPH (0.004%) to 0.1 mL of each test concentration was added. The resulting mixtures were incubated in the dark for 30 min at room temperature. The decrease in absorbance was measured at 517 nm using a UV-2250 (Shimadzu) spectrophotometer against a black. Scavenging (%) of DPPH free radical by essential oil was calculated as follows:

$$\text{Scavenging } (\%) = \left(A_{\text{blank}} - A_{\text{sample}}/A_{\text{blank}}\right) \times 100$$

where A_{blank} is the absorbance of the control reaction (containing all reagents except the test compound) and A_{sample} is the absorbance of the test oil.

All determinations were performed in triplicate. The DPPH scavenging activity of butylated hydroxytoluene (BHT), a widely used antioxidant recognised as safe for use in foods, was evaluated as a comparison.

Statistical analysis

The means values and standard deviations of volatile compound data were calculated using Excel 2019 software (Microsoft, Milan, Italy).

For the QDA analysis, the confidence interval method was used to eliminate the variance associated with the panellists. The odour descriptors were mean-centred per panellist by scaling with unit variance at a 95% confidence level. Then, the mean values and percentiles of the 8 samples were derived from the values assigned by the 12 panellists, and differences between them were analysed by one-way Analysis of Variance ANOVA and Duncan's multiple range test at a confidence level of 95% using XLstat software 2014 (Addinsoft, Paris, France). Frequency histograms of all odour scores were generated by using Excel 2019 software (Microsoft, Milan, Italy). A spider plot of odours was constructed by using Microsoft Excel 2019 (MS Office 2010, Microsoft Corporation, Seattle, WA, USA) based on the mean scores for the different odours (mean values for 8 samples assessed by 12 panellists).

Results and discussion

Volatile aroma compounds

Table 1 reports the average volatile composition as single compounds and classes of substances of the Niaouli essential oils analysed and a comparison with literature data (Ramanoelina *et al.*, 1992, 1994, 2005, 2008). Figure 1 shows a GC-MS chromatogram of a Niaouli essential oil. Globally, about ninety volatile compounds have been identified by GC-MS analysis, most of these, here for the first time reported. The identified volatiles belonged to different classes of compounds with monoterpenes and sesquiterpenes, both hydrocarbon and oxygenated, being the most represented. Looking at the table the percentage of each volatile agrees with Ramanoelina *et al.* (2008), at least for the common ones. Among monoterpenes, 1,8-cineole (eucalyptus-like) prevailed, as expected; numerous studies have shown that 1,8-cineole has anti-inflammatory, antioxidant, analgesic and potent antimicrobial properties against a wide range of microorganisms, thus has the potential to be an antibacterial agent for meat preservation (Sobhy *et al.*, 2023).

Limonene (citrus, lemon), α-pinene (pine, hearty, woody) and β-pinene (woody, pine) followed in agreement with Ramanoelina *et al.* (2008). Several monoterpenes were, here, for the first time identified, especially monoterpene alcohols and esters such as α-terpenyl acetate at a level of 2.10% which has been found in Niaouli essential oil 1,8-cineole chemotype from New Caledonia (Trilles *et al.*, 2006).

Sesquiterpenes represented 20.72% of the total volatile fraction; viridiflorol and (*E*)-nerolidol were the most abundant, responsible for green-sweet and woody-flower notes, respectively; β-caryophyllene (2.06%) and viridiflorene (1.32%) followed. These are typical sesquiterpenes of *Melaleuca* sp. leaf essential oils with a different amount in the four chemotypes of *Melaleuca quinquenervia*. Viridiflorol and viridiflorene are rare aromadendrane tricyclic sesquiterpenoids with a characteristic skeleton, a dimethyl cyclopropane ring fused to a hydroazulene ring system; they have been detected in the extracts of many Myrtaceae and Lamiaceae plants but until now only two plant viridiflorol synthases have been identified (Ntana *et al.*, 2021). Viridiflorol has recently demonstrated

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Table 1 Average composition as single compound and classes of substances of the 1,8 cineole type Niaouli essential oils from Madagascar

	Compounds	LRI	LRI lit.	Area %	SD	Area % Ref. 4	Ref.
	Monoterpene hydrocarbons						
3	α-Thujene	924	924	0.20	0.03	nd	1
4	α-Pinene	932	932	9.38	0.52	10.92	1,2,3,4
5	α-Fenchene	947	944	0.03	0.01	nd	–
6	Camphene	948	948	0.11	0.01	nd	2
8	Sabinene	971	969	0.01	0.01	nd	2
9	β-Pinene	977	977	2.73	0.14	2.84	2,3,4
11	Myrcene	988	987	0.89	0.05	0.89	1,2,3,4
12	*p*-Menta-1(7),8-diene	1006	1008	0.07	0.01	nd	–
13	α-Phellandrene	1007	1003	0.07	0.01	nd	1,2
14	δ-3-Carene	1009	1010	0.01	0.01	nd	2
15	α-Terpinene	1016	1016	0.27	0.02	nd	2
16	*p*-Cymene	1023	1025	1.82	0.08	1.06	2,3,4
17	Limonene	1030	1030	9.18	0.84	8.12	1,2,3,4
19	(*E*)-β-Ocimene	1043	1046	0.04	0.01	nd	–
20	γ-Terpinene	1056	1056	1.17	0.07	1.71	2,3,4
21	Terpinolene	1084	1086	0.59	0.03	0.68	1,2,3,4
22	*p*-Cymenene	1090	1090	0.05	0.01	nd	–
	Total			26.62	1.86	26.22	
	Sesquiterpene hydrocarbons						
44	Isoledene	1371	1371	0.01	0.01	nd	–
45	α-Copaene	1376	1374	0.13	0.01	nd	2
46	α-Gurjunene	1406	1406	0.17	0.02	0.04	1,2,3,4
47	β-Caryophyllene	1418	1417	2.06	0.14	1.37	1,2,3,4
48	Aromadendrene	1438	1438	0.06	0.01	nd	2
49	α-Guajene	1439	1432	0.17	0.01	nd	–
50	Amorpha-4,11-diene	1441	–	0.03	0.01	nd	–
51	Selina-5,11-diene	1445	1441	0.02	0.01	nd	–
52	α-Humulene	1454	1452	0.37	0.04	nd	1,2
53	Alloaromadendrene	1459	1460	0.59	0.05	nd	2
54	Cadina-1(6),4-diene	1472	1475	0.03	0.01	nd	–
55	Selina-4,11-diene	1473	–	0.04	0.01	nd	–
56	γ-Muurolene	1475	1477	0.05	0.02	nd	–
57	β-Selinene	1488	1490	0.46	0.06	nd	1
58	Viridiflorene	1491	1493	1.32	0.17	0.73	1,2,3,4
59	α-Selinene	1495	1494	0.30	0.04	nd	–
60	α-Muurolene	1498	1499	0.07	0.01	nd	–
61	β-Bisabolene	1508	1509	0.01	0.01	nd	–
62	γ-Cadinene	1513	1513	0.25	0.03	0.23	1,2,3,4
63	δ-Cadinene	1519	1522	0.30	0.01	0.23	1,2,3,4
64	*trans*-Calamenene	1521	1520	0.03	0.01	nd	2
65	*trans*-Cadina-1,4-diene	1532	1530	0.02	0.01	nd	2
66	α-Cadinene	1537	1541	0.03	0.01	nd	2
67	(*E*)-α-Bisabolene	1541	1547	0.01	0.01	nd	–
	Total			6.53	0.72	2.60	
	Oxygenated monoterpenes						
	Carbonyl compounds						
33	*cis*-Pinocamphone	1179	–	0.06	0.01	nd	–
	Alcohols						
24	Linalool	1102	1101	0.24	0.03	0.24	1,2,3,4
26	Fenchyl alcohol	1120	1120	0.06	0.01	nd	2
27	*trans*-Pinocarveol	1141	1141	0.04	0.01	nd	–
28	Isopulegol	1149	1145	0.11	0.02	nd	–
29	Camphene hydrate	1156	1150	0.02	0.01	nd	–
30	*iso*-Isopulegol	1159	1160	0.05	0.01	nd	–
31	δ-Terpineol	1171	1165	0.20	0.01	nd	–
32	Borneol	1172	1173	0.06	0.01	nd	–
34	Terpinen-4-ol	1181	1179	1.05	0.26	0.71	1,2,3,4
35	*p*-Cymen-8-ol	1188	1183	0.05	0.01	nd	–
36	α-Terpineol	1196	1199	7.10	0.84	6.35	1,2,3,4
37	Citronellol	1229	1228	0.07	0.01	nd	–
38	Geraniol	1253	1253	0.02	0.01	nd	–
43	Eugenol	1353	1356	0.05	0.01	nd	–
	Total			9.12	1.25	7.30	
	Epoxides						
25	*cis*-Rose oxide	1112	1112	0.02	0.01	nd	–
	Esters						
40	δ-Terpineol acetate	1314	1316	0.04	0.01	nd	–
41	Myrtenyl acetate	1324	1328	0.02	0.01	nd	–
42	α-Terpinyl acetate	1348	1350	2.10	0.34	nd	–
	Total			2.16	0.36		
	Eters						
18	1,8-Cineole	1033	1033	44.06	1.19	55.38	1,2,3,4
	Oxygenated sesquiterpenes						
	Alcohols						
68	*epi*-Longipinalol	1557	1557	0.04	0.01	nd	–
69	(*E*)-Nerolidol	1562	1563	3.18	0.29	2.62	1,2,3,4
70	Palustrol	1570	1570	0.25	0.03	nd	–
71	Thujopsan-2-α-ol	1582	1579	0.38	0.05	nd	–
73	Viridiflorol	1594	1591	2.59	0.57	4.98	1,2,3,4
74	Cubeban-11-ol	1597	1592	0.26	0.03	nd	–
75	Copaborneol	1606	1592	0.11	0.07	nd	–
77	Eremoligenol	1630	1631	0.04	0.01	nd	–
78	*epi*-Cubenol	1635	1631	0.02	0.01	nd	–
79	γ-Eudesmol	1639	1637	0.06	0.01	nd	–
80	α-Caryophylladienol	1645	1645	0.03	0.01	nd	–
81	*epi*-α-Cadinol	1650	1656	0.08	0.01	nd	3
82	Cadin-4-en-10-ol	1652	1658	0.02	0.01	nd	-
83	α-Muurolol	1655	1643	0.01	0.01	0.06	4
84	β-Eudesmol	1664	1664	0.16	0.02	nd	–
85	Bulnesol	1673	1672	0.02	0.01	nd	–
	Total			9.97	1.15	7.66	
	Epoxides						
72	Caryophyllene oxide	1586	1584	0.09	0.01	0.29	1,2,4
76	Humulene epoxide II	1612	1603	0.05	0.01	nd	–
	Total			0.14	0.02	0.29	
	Aliphatic compounds						
	Alcohols						
1	(*Z*)-4-Hexen-1-ol	873	879	0.02	0.01	nd	–
	Ketones						
10	6-Methyl-5-hepten-2-one	986	986	0.03	0.01	nd	–
	Hydrocarbons						

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Table 1 (*Continued*)

	Compounds	LRI	LRI lit.	Area %	SD	Area % Ref. 4	Ref.
86	Eicosane	1997	2000	0.53	0.06	nd	–
	Aromatic compounds						
	Aldehydes						
7	Benzaldehyde	962	962	0.29	0.05	0.19	1,2,3,4
	Esters						
23	Methyl benzoate	1096	1094	0.06	0.01	nd	–
	Nitrogen compounds						
39	1-Nitro-2-Phenyl ethane	1293	1300	0.07	0.01	nd	–
	Hydrocarbons						
2	Styrene	897	897	0.02	0.01	nd	–

LRI = Linear Retention Index calculated on SLB-5 capillary column; LRI lit. = Linear Retention Index from Literature; Area % = average of all the essential oil samples analysed in triplicate; Area % Ref. 4 = average composition of 1,8-cineole type Niaouli essential oil from madagascal reported by Ramanoelina *et al.*, 2008; Ref. = (1) Ramanoelina *et al.*, 1992; (2) Ramanoelina *et al.*, 1994; (3) Ramanoelina *et al.*, 2005; (4) Ramanoelina *et al.*, 2008.
nd, not determined; SD, Standard Deviation.

anti-neoplastic properties inducing apoptosis of breast, lung and brain cancer cells (Akiel *et al.*, 2022). (*E*)-Nerolidol is an acyclic sesquiterpene widely spread in the essential oil of various medicinal plants which are well known for their pharmacological and biological activities; for this reason, the U.S. Food and Drug Administration (FDA) allows its use as a food flavouring agent (Chan *et al.*, 2016). β-caryophyllene is a bicyclic sesquiterpene widely spread in nature, too; it is responsible for woody, sweet, spicy notes and it has wide applications as food flavouring. Moreover, sesquiterpenes are described as possessing anti-inflammatory and antibacterial properties including enhancement of bacterial susceptibility to antibiotics, and to exhibit anticarcinogenic and antioxidant effects (Cincotta *et al.*, 2015). Among the sesquiterpenes identified for the first time, α-selinene and α-bisabolene isomers were present. α-Selinene is one of the major sesquiterpenes in the celery volatile profile (Naghneh *et al.*, 2023) while α-bisabolene is a primary component in black pepper (Eddin *et al.*, 2022).

Sensory analysis

The QDA revealed different perceived odours with different intensities (Fig. 2a). According to the judges' evaluation, the oils have been characterised by the following descriptors: 'woody, fresh-mint, eucalyptus-like, sweet-floral and spicy'. The descriptors 'eucalyptus-like and woody' odour showed the highest mean scores followed by 'fresh-minty' odour.

Further, 'eucalyptus-like' showed the highest median score, with the narrowest lowest-to-highest range and interquartile range; the descriptor of 'fresh-minty' odour was second to this (Fig. 2b); their intensity was the most stable and the highest.

'Eucalyptus-like' odour is mainly related to the high content of 1,8-cineole or eucalyptol while 'fresh-minty' note could be related to monoterpene alcohols such as α-terpineol. 'Woody' odour could be attributable to the presence of monoterpenes and sesquiterpene hydrocarbons including α and β-pinene, *p*-cymene, γ-terpinene, β-caryophyllene and (*E*)-nerolidol; the 'sweet-floral' odour could be related to oxygenated terpenes such as (*E*)-nerolidol whereas 'spicy' to sesquiterpenes such as β-caryophyllene and germacrene D which are responsible for 'woody and spicy' notes. Hence, the typical aromatic characteristics of Niaouli oil were mostly attributed to the dominant 'eucalyptus-like' and 'fresh-minty' odour.

Antioxidant activity

The antioxidant activity of the Niaouli essential oils was assessed using the DPPH assay. The % inhibition was found to increase from 4.66% to 20.02% with the increase in the concentration of the essential oil from 10 to 100 μL/mL. The oils were found to exhibit similar antioxidant activity if compared to BHT (Table 2). This could be due to other constituents such as terpenoids with reactive sites such as hydroxyl

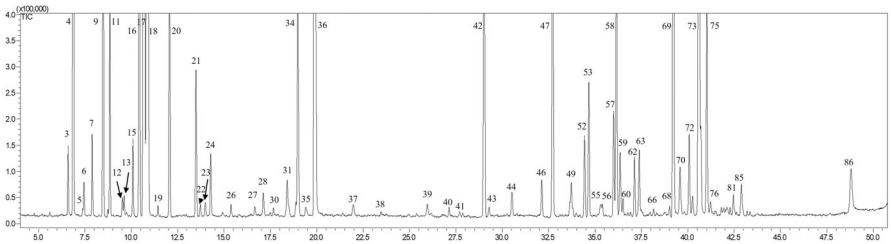


Figure 1 HS-SPME-GC-MS chromatogram of 1,8-cineole type Niaouli essential oils from Madagascar. The identification of peaks is reported in Table 1.

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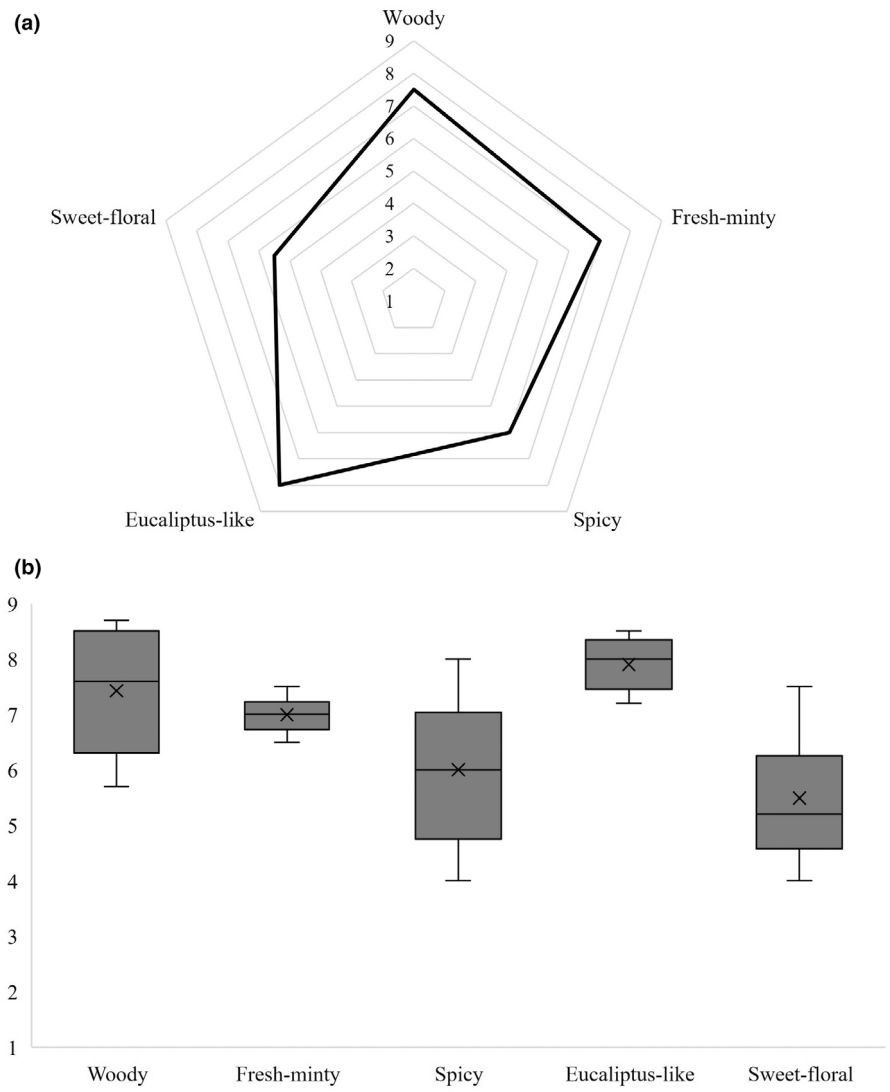


Figure 2 Graphical representation of the average data of QDA analysis (a) and score percentiles box for the odours (b) assigned by 12 panellists of all the 1,8-cineole type Niaouli essential oils from Madagascar analysed.

Table 2 Antioxidant activity of Niaouli essential oils measured in terms of DPPH radical scavenging

	Concentration (μL/mL)	Niaouli oils	SD	BHT
DPPH (%)	10	4.66	0.23	6.18
	20	6.76	0.59	9.22
	50	11.06	0.97	17.42
	100	20.02	1.34	31.98

BHT, antioxidant activity of butylated hydroxytoluene; SD, Standard Deviation.

groups useful for electron transfer which can contribute to the antioxidant potential of Niaouli essential oil (Bakkali *et al.*, 2008).

Conclusions

The consumer request for natural ingredients in food and beverages is increasing and the market of natural flavourings is expected to witness significant growth in the next years. In this context, the characterisation of the 1,8-cineole type Niaouli essential oil from Madagascar using the volatile and odour profile results in considerable interest. Using advanced analytical techniques, many volatiles, especially sesquiterpenes, have been identified. The Quantitative Descriptive Sensory Analysis has been able to identify the descriptors that characterised the oil; volatile compounds and odour descriptors resulted in agreement. Considering its volatile composition, the pleasant sensory notes of

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'eucalyptus-like' and 'fresh-minty' and, the antioxidant activity, the 1,8-cineole type Niaouli essential oil from Madagascar could be considered as a flavouring agent and natural preservative in the food industry mainly in the field of candy, baked products, meat and beverages.

Author contributions

Maria Merlino: Methodology; investigation; writing – original draft. **Fabrizio Cincotta:** Investigation; data curation; visualization; supervision. **Antonella Verzera:** Conceptualization; writing – review and editing; methodology; data curation; resources. **Anthea Miller:** Formal analysis. **Marco Torre:** Formal analysis. **Concetta Condurso:** Conceptualization; writing – review and editing; supervision.

Conflict of interest

There is no conflict of interest.

Ethical approval

Ethics approval was not required for this research. All participants involved in the sensory analysis signed an informed consent according to the principles of the Declaration of Helsinki before the beginning of the study.

Data availability statement

The data that support the findings of this study are available from the corresponding author upon reasonable request.

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Niaouli essential oil with a high content of 1,8-cineole is the most widespread and appreciated chemotype. Different studies have been carried out on Niaouli essential oil with a high content of 1,8-cineole volatile composition even if this dates back to more than ten years ago (Ramanoelina *et al*., 1994).

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Supporting Information

Additional Supporting Information may be found in the online version of this article:

Table S1. Lexicon used to describe the sensory characteristics of the 1,8-cineole type Niaouli essential oils from Madagascar.

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