

SUPPLEMENTARY S1: Composition of volatile organic compounds (VOCs) in selected insect and plant oils analyzed by Gas Chromatography coupled to Mass Spectrometry (GC-MS)

				Mean ± SE, (standard error) (ng/μL of oil)					
RT	Compound	RI ^{Cal*}	Idf	<i>Ruspolia differens</i>	<i>Macrotermes spp</i>	<i>Schistocerca gregaria</i>	<i>Ocimum kilimandscharicum</i>	<i>Eucalyptus citriodora</i>	<i>Cymbopogon nardus</i>
Alcohol									
6.17	3-methyl-2-butanol	778	RI,MS		2.52±0.00			2.92±0.19	
6.46	4-methyl-2-pentanol	788	RI,MS			2.80±0.21			
6.46	2-Hexanol	788	Std,RI,MS	2.9±0.21					
10.36	1-Octen-3-ol	945						4.32±0.82	
20.17	1α, 2β, 3α, 5β-Cyclohexanetetrol,	1587				20.38±3.92			
25.42	Z7-C16-enol	2120				98			
Aldehyde									
6.34	Hexanal	783	Std,RI,MS			2.63±0.10			
21.85	(E)-Chrysanthemal	1746	RI,MS					5.03±1.04	
Fatty acid									
8.40	2-Piperidinecarboxylic acid	863	RI,MS	3.6±0.40					
15.63	Nonanoic acid	1247	Std,RI,MS			12.65±3.45			
17.04	n-Decanoic acid	1344	Std,RI,MS			43.72±14.06			
19.32	Dodecanoic acid	1517	Std,RI,MS	9.0±0.74	65.15	1874.21±152.32			
20.73	γ-dodecalactone	1635	RI,MS		12.37±1.12	21.34±0.75			
21.15	Methyl tetradecanoate	1672	Std,RI,MS	5.7±0.09	16.83±2.21				
21.55	Tetradecanoic acid	1709	Std,RI,MS	8.8±0.65	14.09±2.59	380.51±40.45			
21.88	Ethyl tetradecanoate	1749	RI,MS	16.8±1.77					
21.92	Methyl pentadecanoate	1754	Std,RI,MS		6.93±0.63				
22.21	Methyl 2-hydroxydodecanoate	1789	Std,RI,MS			21.92±10.02			
22.59	Pentadecanoic acid	1827	Std,RI,MS		5.95±0.59	29.75±1.39			
22.87	14-methylpentadecanoate	1853	RI,MS		5.98±0.46				
22.91	Ethyl pentadecanoate	1857	RI,MS	5.4±0.12					
23.23	Methyl hexadecanoate	1887	Std,RI,MS	29.2±11.52	84.08±12.83				
23.42	9Z-Hexadecenoic acid	1905	Std,RI,MS	24.1±4.07	14.53±2.66	180.24±32.98			
23.59	Methylhexadecanoate	1924	Std,RI,MS	7.1±0.88	31.23±5.12				
23.70	Ethyl 9-hexadecenoate	1936	RI,MS	82.9±42.04					

23.85	Methyl heptadecanoate	1952	Std,RI.MS		8.96±0.71				
23.90	Ethyl hexadecanoate	1957	RI.MS	270.0±105.51					
23.94	14-methyl hexadecanoate	1962	Std,RI.MS		7.48±0.57				
24.00	Methyl 10Z-heptadecenoate	1968	RI.MS		11.27±1.12				
24.21	Methyl heptadecanoate	1991	Std,RI.MS		12.13±1.22				
24.41	<i>n</i> -Hexadecanoic acid	2013	Std,RI.MS	192.5±101.38	56.20±13.79	1046.22±50.20			
24.41	methyl heptadecanoate	2013	Std,RI.MS	119.2±88.77					
24.56	Heptadecanoic acid	2028	Std,RI.MS	25.1±3.99		41.53±10.01			
24.62	Methyl 8-heptadecenoate	2035	RI.MS	18.9±3.80					
24.68	9Z,12Z-Octadecadienoic acid	2041	Std,RI.MS		10.18±1.05				
24.91	Methyl 9E-Octadecenoate	2066	Std,RI.MS	103.5±46.54	34.43±26.48				
24.92	Methyl 9Z-Octadecenoate	2067	Std,RI.MS		159.23±25.03				
25.14	Methyl octadecenoate	2090	Std,RI.MS	64.2±22.02	30.55±4.02				
25.32	Oleic Acid	2109	Std,RI.MS	378.6±187.01					
25.33	Methyl linoleate	2111	Std,RI.MS		48.56±12.12				
25.46	Octadecanoic acid	2125	Std,RI.MS		41.72±13.85				
25.55	Ethyl Oleate	2135	RI.MS	698.4±194.30					
25.74	Ethyl octadecenoate	2155	RI.MS	429.6±145.41					
26.34	(9Z,12Z,15Z)-octadecatrienoic acid	2222	Std,RI.MS	5.3±0.08					
26.36	Ethyl 5,8,11,14-eicosatetraenoate	2225	RI.MS	32.1±11.02	10.06±1.36				
26.61	Ethyl nonadecanoate	2254	RI.MS	41.4±9.44					
26.90	Ethyl eicosanoate	2288	RI.MS	48.7±6.59	15.55±2.57				
27.24	5Z,8Z,11Z,14Z-eicosatetraenoic acid	2329	RI.MS	42.1±7.45					
27.68	(5Z,8Z,11Z,14Z,17Z)-eicosapentaenoic acid	2382	RI.MS	31.9±5.32					
Ketone									
6.78	Octane	800	Std,RI.MS	2.7±0.08					
6.91	6,6-Dimethylhepta-2,4-diene	805	RI.MS		2.52±0.00	3.09±0.25			
8.16	o-Xylene	854	Std,RI.MS	4.3±0.45	3.95±0.64	4.87±1.08	4.43±0.96	3.73±0.42	
8.17	p-Xylene	854	Std,RI.MS	4.2±0.72	4.46±0.68	3.91±0.48	3.49±0.44	4.53±0.37	3.22±0.12
8.47	2,4,6-Trimethyl-3-heptene	866	RI.MS	4.9±0.09					
8.47	1,3-Bis(cyclopentyl)-1-cyclopentanone	866	RI.MS			3.40±0.01			
8.60	3-methyl 2E-Pentene	871	RI.MS					3.39±0.57	
8.75	Ethylbenzene	877	Std,RI.MS					3.09±0.26	

9.97	2,6-dimethyl-4-hepten-3-one	928	RI.MS	6.9±0.15	4.89±0.00	4.89±0.00		
24.29	Eicosane	2000	Std,RI.MS		9.43±1.76			
Monoterpene								
9.02	Tricyclene	887	Std,RI.MS			4.56±0.60		
9.18	α -thujene	893	Std,RI.MS			3.84±0.10		
9.62	α -phellandrene	912	Std,RI.MS			50.46±13.97		
9.64	α -Pinene	913	Std,RI.MS	4.5±0.60	3.70±0.38	12.96±2.51	4.53±0.62	3.77±0.35
10.17	4-Carene	937	Std,RI.MS				8.35±1.62	
10.21	β -Pinene	939	Std,RI.MS			10.35±1.58		
10.56	β -Myrcene	954	Std,RI.MS			6.41±1.29	6.91±0.55	6.20±0.12
11.23	δ -3-Carene	985	Std,RI.MS	9.0±0.48	7.55±1.16		8.73±0.69	6.64±0.42
11.24	γ -Terpinene	985	Std,RI.MS	8.1±0.34				
11.25	Limonene	985	Std,RI.MS			143.90±91.50	31.13±7.84	16.68±0.75
11.52	o-Cymene	998	Std,RI.MS			5.82±1.36		
11.63	Eucalyptol	992	Std,RI.MS			37.37±21.31		
12.35	L-Fenchone	1040	Std,RI.MS			6.73±1.82		
12.70	Terpinolene	1064	Std,RI.MS				5.97±0.36	
12.72	α -Terpinene	1066	Std,RI.MS					5.51±0.04
12.75	(E)-Rose oxide	1068	RI.MS				6.23±0.46	
12.89	Linalool	1077	Std,RI.MS			38.41±21.10	10.88±1.73	7.65±0.43
12.93	<i>trans</i> -p-Mentha-2,8-dienol	1080	RI.MS			12.21±3.42		
13.31	Isopulegol	1105	RI.MS				11.50±4.52	12.01±1.47
13.34	Camphor	1107	Std,RI.MS			382.74±136.49		
13.42	Citronellal	1111	Std,RI.MS				183.61±38.04	100.03±36.59
13.49	Isoborneol	1115	Std,RI.MS			8.94±2.59		
13.63	Borneol	1124	Std,RI.MS			11.92±2.84		
13.81	Terpinen-4-ol	1134	Std,RI.MS			31.37±15.23	7.86±2.09	
14.07	α -Terpineol	1149	Std,RI.MS			15.13±6.63	9.14±2.57	
14.12	Myrtenol	1152	Std,RI.MS			15.16±3.74		
14.57	Citronellol	1179	Std,RI.MS				94.96±27.90	53.96±7.46
14.96	Geraniol	1202	Std,RI.MS				117.42±32.81	71.21±7.87
16.27	3-Cyclopentylcyclopentene	1289	RI.MS		5.11±0.08			
16.37	Citronellyl acetate	1296	RI.MS				20.29±3.74	
16.74	2,6-dimethyl-2,6-octadiene	1322	RI.MS					13.23±0.61

18.12	1,6-Octadien-3-ol, 3,7-dimethyl-, formate	1423	RI.MS				5.42±0.43
	sesquiterpene						
16.35	α -Cubebene	1295	Std,RI.MS		5.79±1.27		
16.72	α -Copaene	1321	Std,RI.MS		8.82±2.23		
16.85	β -Bourbonene	1330	RI.MS		8.72±2.45		
16.92	β -Elemene	1335	RI.MS			22.94±4.87	
17.15	Muurolo-3,5-diene	1352	RI.MS	5.3±0.22	5.43±1.00		
17.17	α -Gurjenene	1353	RI.MS		5.64±0.99		
17.27	β -Elemene	1360	RI.MS		6.55±1.17		14.58±1.17
17.30	β -ylangene	1363	RI.MS			10.11±1.82	
17.64	Isolatedene	1387	RI.MS			7.31±0.58	
17.66	(E)-Caryophyllene	1389	Std,RI.MS		13.28±1.95	7.49±0.38	
17.79	Copaene	1398	Std,RI.MS		7.77±1.25		6.75±0.30
18.02	γ -Muuroloene	1415	Std,RI.MS		5.34±0.95	7.29±0.55	5.3±0.35
18.10	Germacrene D	1422	Std,RI.MS		7.18±2.59	9.24±0.95	9.46±0.41
18.13	Humulene	1424	Std,RI.MS		7.44±2.43		5.85±0.17
18.30	α -Muuroloene	1437	RI.MS		4.71±1.00	8.16±1.28	7.55±0.28
18.46	Germacrene B	1449	RI.MS			11.07±2.28	
18.50	α -Cadinene	1452	RI.MS			10.85±3.38	
18.50	γ -Cadinene	1452	RI.MS		7.29±1.40		
18.59	Calamene	1459	RI.MS		6.87±1.15		
18.67	β -Muuroloene	1466	Std,RI.MS			11.89±1.89	
18.85	α -Calacorene	1479	RI.MS		6.47±1.29		
18.86	γ -Cadinene	1480	RI.MS			8.11±1.07	8.79±0.65
18.95	δ -Cadinene	1487	RI.MS			20.01±2.22	13.78±1.11
19.27	Elemol	1513	RI.MS			20.51±7.50	15.94±1.46
19.35	Caryophyllene oxide	1519	Std,RI.MS		15.15±6.01		
19.62	Dauca-5,8-diene	1541	RI.MS			8.29±1.71	
19.64	(E)-Muurolo-3,5-diene	1543	RI.MS		4.90±0.98		7.2±0.26
19.91	α -Eudesmol	1565	RI.MS			7.16±0.45	12.95±1.07
20.01	β -Cadinene	1574	RI.MS		6.81±1.87		
20.29	γ -Eudesmol	1597	RI.MS			11.31±1.67	
20.40	tau-Cadinol	1606	RI.MS			7.80±2.58	
20.42	Amorpha-4,7(11)-diene	1608	RI.MS			12.15±6.90	

Sterols						
26.74	10-oxo-methyl octadecanoate	2269	RI,MS		11.51±1.44	
33.64	γ-Tocopherol	3009	Std,RI,MS	17.2±7.26		
34.11	Cholestan-3-ol, (3, beta,5, beta.)-	3027	RI,MS		22.96±6.59	
34.72	Cholesterol	3049	Std,RI,MS	60.0±52.60	41.84±9.24	
36.58	Campesterol	3297	Std,RI,MS		29.26±13.24	
37.26	Stigmasterol	2595	Std,RI,MS		28.24±17.70	
37.61	Stigmastanol	3366	Std,RI,MS		13.35±4.35	
38.47	β-Sitosterol	3385	Std,RI,MS		24.97±10.22	
Triterpene						
22.05	2,6,10,14-tetramethyl hexadecane	1770	RI,MS		7.47±1.12	
30.96	Squalene	2787	Std,RI,MS	48.0±18.94	25.41±4.66	105.40±9.95
31.34	Peucelinendiol	2824	RI,MS			6.49±1.03

Relative mass values presented as Mean M± SE, (SE =standard error) of triplicate determinations (*ng/μL* of oil) from three independent analyses. Id^f Definitive compound identification via authentic standard; if not authenticated then compound identity is tentative by RI match to only one column. Ie Std, comparison with available analytical standards; RI, linear with those calculated with ADAMS (2007) [20] and NIST (2020) [21] libraries. comparison to published mass spectral data (MS). *Retention index for HP-5 MS column using a homologous series of n-alkanes and a linear GC ramp program.

Supplementary Table 2a: Repellency (%) (±SE) at hours after treatment and CPT of 1:1 <i>C.nardus</i> /Macrotermes spp against <i>Ae. aegypti</i> in laboratory test																			
Treatment	Volunteer	N	0min		30min		60min		120min		150min		180min		240min		270min		CPT
			N	R(%)	N	R(%)	N	R(%)	N	R(%)	N	R(%)	N	R(%)	N	R(%)	N	R(%)	
50%CTN/50%TMT	V-1	12	0	100	0	100	0	100	0	100	8	96	10	95	20	90	22	89	2.5
	V-2	17	0	100	0	100	0	100	0	100	0	100	0	100	15	92.5	23	88.5	3
	V-3	23	0	100	0	100	0	100	0	100	11	94.5	16	92	16	92	27	86.5	2
	V-4	25	0	100	0	100	0	100	0	100	0	100	0	100	10	95	20	90	3
	V-5	27	0	100	0	100	0	100	0	100	0	100	0	100	22	89	34	83	3
	V-6	34	0	100	0	100	0	100	0	100	0	100	0	100	32	84	36	82	3.5
	V-7	11	0	100	0	100	0	100	0	100	0	100	0	100	40	80	40	80	3
	V-8	19	0	100	0	100	0	100	0	100	0	100	5	97.5	25	87.5	36	82	2.5
	AVG	21.0	0.0	100.0	0.0	100.0	0.0	100.0	0.0	100.0	2.4	98.8	3.9	98.1	22.5	88.8	29.8	85.1	2.8
	StdErr	2.76	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	1.58	0.79	2.16	1.08	3.44	1.72	2.70	1.35	0.46

Supplementary Table 2b: Repellency (%) (±SE) at hours after treatment and CPT of 1:1 <i>C.nardus</i> /Macrotermes spp against <i>Cx quinquefasciatus</i> in laboratory test																			
Treatment	Volunteer	N	0min		30min		60min		120min		150min		180min		240min		270min		CPT
			N	R(%)	N	R(%)	N	R(%)	N	R(%)	N		N	R(%)	N	R(%)	N	R(%)	
50%CTN/50%TMT	V-1	20	0	100	0	100	0	100	0	100	0	100	4	98	9	95.5	26	87	2.5
	V-2	23	0	100	0	100	0	100	0	100	2	99	2	99	9	95.5	10	95	2
	V-3	36	0	100	0	100	0	100	0	100	0	100	0	100	8	96	14	93	3
	V-4	40	0	100	0	100	0	100	0	100	0	100	0	100	7	96.5	56	72	3
	V-5	40	0	100	0	100	0	100	0	100	0	100	4	98	14	93	16	92	2.5
	V-6	50	0	100	0	100	0	100	0	100	0	100	4	98	16	92	24	88	2.5
	V-7	25	0	100	0	100	0	100	0	100	0	100	6	97	13	93.5	27	86.5	2.5
	V-8	41	0	100	0	100	0	100	0	100	0	100	6	97	14	93	20	90	2.5
	AVG	34.4	0.0	100.0	0.0	100.0	0.0	100.0	0.0	100.0	0.3	99.9	3.3	98.4	11.3	94.4	24.1	87.9	2.6
	StdErr	3.73	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.25	0.13	0.84	0.42	1.19	0.60	5.02	2.51	0.32

Supplementary Table 2c: Repellency (%) (±SE) at hours after treatment and CPT of 1:1 *C.nardus*/Macrotermes spp against *An. gambiae* in laboratory test

Treatment	Volunteer	N	0min		30min		60min		120min		150min		180min		240min		270min		CPT
			N	R(%)	N	R(%)	N	R(%)	N	R(%)	N	R(%)	N	R(%)	N	R(%)	N	R(%)	
50%CTN/50%TMT	V-1	20	0	100	0	100	0	100	0	100	0	100	0	100	0	100	6	97	4
	V-2	11	0	100	0	100	0	100	0	100	0	100	0	100		100	10	95	3
	V-3	47	0	100	0	100	0	100	0	100	0	100	0	100	8	96	12	94	4
	V-4	50	0	100	0	100	0	100	0	100	0	100	0	100	6	97	8	96	3
	V-5	43	0	100	0	100	0	100	0	100	0	100	0	100	0	100	14	93	3.5
	V-6	38	0	100	0	100	0	100	0	100	0	100	0	100	10	95	16	92	3
	V-7	19	0	100	0	100	0	100	0	100	0	100	0	100	0	100	18	91	3.5
	V-8	37	0	100	0	100	0	100	0	100	0	100	0	100	0	100	22	89	4
	AVG	33.1	0.0	100.0	0.0	100.0	0.0	100.0	0.0	100.0	0.0	100.0	0.0	100.0	3.4	98.5	13.3	93.4	3.5
	StdErr	5.13	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	1.57	0.76	1.89	0.94	0.46

The number (*N*) of mosquitoes landing on the arm of each volunteer was counted per hour for 5 h. Repellency (*R*)was calculated each hour and complete protection time (CPT) was determined by calculating the number of minutes from the time of repellent application to the first mosquito landing.