

Table S1. Identified volatile components in *Jinhua ham* by three extraction methods (SPME, NT and SAFE)

Relative percentage content (%)												
Compounds ¹	RI ²	CAS	Identification method	SPME			NT			SAFE		
				One- year	Two- year	Three- year	One- year	Two- year	Three- year	One- year	Two- year	Three- year
				aging	aging	aging	aging	aging	aging	aging	aging	aging
Aldehyde												
Hexanal	1069.1	66-25-1	RI, MS	5.69±2.26 ^a	3.57±1.83 ^{bc}	0.84±0.11 ^c	1.45±1.37 ^{bc}	2.02±1.35 ^{bc}	0.48±0.44 ^c	1.21±0.91 ^{bc}	0.54±0.15 ^c	0.69±0.21 ^c
Octanal	1353.7	124-13-0	RI, MS	1.15±1.32 ^{ab}	1.08±1.08 ^{ab}	1.53±0.59 ^a	0.08±0.08 ^b	0.00±0.01 ^b	0.28±0.04 ^b	0.14±0.09 ^b	0.05±0.01 ^b	0.10±0.04 ^b
Nonanal	1483.2	124-19-6	RI, MS	2.73±1.37 ^a	1.88±0.25 ^a	1.97±0.47 ^a	0.43±0.30 ^b	0.66±0.41 ^b	0.28±0.04 ^b	0.49±0.24 ^b	0.16±0.03 ^b	0.36±0.11 ^b
Acetaldehyde	531.1	75-07-0	RI, MS	0.35±0.30	0.69±0.63	2.15±0.33	2.50±2.50	-	2.49±1.87	-	-	-
Trans-2-octenaldehyde	1509.4	2548-87-0	RI, MS	0.07±0.04 ^b	0.11±0.03 ^b	0.23±0.04 ^a	0.20±0.08 ^a	0.06±0.04 ^b	0.05±0.02 ^b	0.05±0.03 ^b	0.02±0.01 ^b	0.05±0.02 ^b
2-Methyl-propionaldehyde	595.2	78-84-2	RI, MS	1.21±1.05 ^b	1.60±0.16 ^b	2.97±0.21 ^a	0.89±0.70 ^b	0.68±0.95 ^b	0.29±0.24 ^b	-	-	-
3-Methylbutyraldehyde	740.6	590-86-3	RI, MS	13.29±0.57 ^a	11.50±2.20 ^a	5.64±2.68 ^b	0.36±0.26 ^c	2.48±3.17 ^{bc}	1.58±1.12 ^{bc}	-	-	0.08±0.08 ^c
Benzaldehyde	1553.3	100-52-7	RI, MS	2.17±0.31 ^a	1.82±0.07 ^a	2.04±0.16 ^a	0.49±0.23 ^b	0.69±0.20 ^b	0.47±0.18 ^b	0.58±0.29 ^b	0.28±0.05 ^b	0.34±0.17 ^b
Acetaldehyde, tetramer	2483	108-62-3	RI	0.05±0.05	0.01±0.00	0.01±0.01	0.19±0.27	0.03±0.04	0.01±0.01	-	-	-
Phenylacetaldehyde	1595.6	122-78-1	RI, MS	5.04±0.43 ^b	5.21±0.63 ^b	11.76±2.16 ^a	-	-	0.47±0.33 ^c	4.35±3.87 ^{bc}	1.39±1.22 ^{bc}	3.16±3.36 ^{bc}
Trans-2-decenaldehyde	1630.8	3913-81-3	RI, MS	0.06±0.05 ^b	0.12±0.08 ^{ab}	0.30±0.09 ^a	-	-	0.19±0.16 ^{ab}	0.06±0.08 ^b	0.09±0.12 ^{ab}	-
Propionaldehyde	578.1	123-38-6	RI, MS	0.02±0.03	-	-	-	-	0.00±0.00	-	-	-
Furfural	1514.8	1998-1-1	RI, MS	-	-	0.11±0.16	0.07±0.05	-	0.06±0.04	-	-	0.00±0.00
Methoxy acetaldehyde	2909.8	10312-83-1	RI	0.01±0.01	-	-	-	-	0.22±0.16	-	-	-
Trans-2-heptanaldehyde	1412.9	18829-55-5	RI, MS	-	0.12±0.12	-	-	-	-	-	-	-
2-Undecylenic aldehyde	1759.8	2463-77-6	RI, MS	-	0.02±0.02	-	-	-	-	-	-	-
5-Hexenal	1879	764-59-0	RI	-	-	0.01±0.00	-	-	-	-	-	-

Cis-2-heptanal	1380.7	57266-86-1	RI, MS	-	-	0.97±0.40 ^a	0.17±0.13 ^b	-	-	-	0.06±0.01 ^b	0.29±0.16 ^b
Trans-2-nonenal	1556.1	18829-56-6	RI, MS	-	-	0.13±0.11	-	-	0.04±0.05	-	-	0.03±0.01
Decanal	1530.7	112-31-2	RI, MS	-	-	-	0.03±0.03 ^b	0.06±0.05 ^b	0.33±0.19 ^a	0.06±0.06 ^b	0.01±0.01 ^b	-
Heptanal	1161	111-71-7	RI, MS	-	-	-	1.30±1.71	0.30±0.39	-	0.11±0.06	0.05±0.01	0.09±0.03
2,4-Decendialdehyde	1765.2	2363-88-4	RI, MS	-	-	-	0.13±0.15	-	0.20±0.12	0.14±0.06	-	-
Trans-2-undecylenic aldehyde	1716.7	2463-77-6	RI, MS	-	-	-	0.07±0.07	-	0.09±0.05	-	-	0.01±0.01
3-Hydroxybutyraldehyde	1156.5	107-89-1	RI	-	-	-	0.03±0.05	-	-	-	-	-
2-Methyl-glutaraldehyde	1266.3	123-15-9	RI	-	-	-	0.00±0.00	-	-	-	-	-
4-Ethylbenzaldehyde	1699	4748-78-1	RI, MS	-	-	-	0.03±0.01	-	-	-	-	-
Trans-2,4-decendialdehyde	1716.9	25152-84-5	RI, MS	-	-	-	0.04±0.04	-	-	0.07±0.04	0.06±0.02	0.05±0.05
2,3-Dimethyl-glutaraldehyde	820.8	32749-94-3	RI, MS	-	-	-	-	0.57±0.81	-	-	-	-
2-Butyraldehyde	1058.4	1119-19-3	RI	-	-	-	-	0.00±0.00	-	-	-	0.02±0.02
3,4-Glutaraldehyde	1063.6	4009-55-6	RI	-	-	-	-	0.00±0.00	-	-	-	-
2-Nonenal	1554.6	2463-53-8	RI, MS	-	-	-	-	0.06±0.05	-	0.03±0.01	0.01±0.01	-
2-Methyl-2-butyraldehyde	1317.5	1115-11-3	RI, MS	-	-	-	-	-	0.00±0.00	0.02±0.01	-	-
5-Methyl furfural	1638.3	620-02-0	RI, MS	-	-	-	-	-	0.05±0.06	-	-	-
Undecylaldehyde	1662.7	112-44-7	RI, MS	-	-	-	-	-	0.05±0.06	-	-	-
3-Methyl-benzaldehyde	1679.4	620-23-5	RI, MS	-	-	-	-	-	0.01±0.01	-	-	-
Methylal	2060.6	109-87-5	RI	-	-	-	-	-	0.04±0.05	-	-	-
5-Hydroxymethylfurfural	2544.1	67-47-0	RI	-	-	-	-	-	0.23±0.09	-	-	-
Glutaraldehyde	1004.2	110-62-3	RI, MS	-	-	-	-	-	-	0.48±0.27	0.28±0.09	0.37±0.13
2-Hexenal	1184.6	505-57-7	RI, MS	-	-	-	-	-	-	0.00±0.00	0.00±0.00	0.00±0.00
3-Methyl-2-butyraldehyde	1167.4	107-86-8	RI, MS	-	-	-	-	-	-	0.03±0.02	0.02±0.01	0.04±0.03

2-Hydroxybenzaldehyde	1605.7	1990-2-8	RI, MS	-	-	-	-	-	-	-	0.04±0.02	0.02±0.01	-
2,6-Dimethyl-5-heptanal	1030.7	106-72-9	RI, MS	-	-	-	-	-	-	-	0.01±0.01	-	0.00±0.00
2-Methyl-2-pentenal	1056.7	623-36-9	RI, MS	-	-	-	-	-	-	-	0.00±0.00	-	-
Tetradecane aldehyde	1847.4	124-25-4	RI, MS	-	-	-	-	-	-	-	0.00±0.00	-	-
4-Methoxybenzaldehyde	1940.3	123-11-5	RI, MS	-	-	-	-	-	-	-	-	0.01±0.01	-
Trans-2,6-nondialdehyde	1258	17587-33-6	RI, MS	-	-	-	-	-	-	-	-	-	0.00±0.00
Trans-2,4-heptadienaldehyde	1517	4313-3-5	RI, MS	-	-	-	-	-	-	-	-	-	0.01±0.01
<i>Ketones</i>													
Acetone	600.2	67-64-1	RI, MS	6.37±5.55 ^a	6.11±0.33 ^{ab}	5.64±2.68 ^{ab}	1.68±0.82 ^c	2.19±1.50 ^{bc}	1.07±0.12 ^c	0.91±0.30 ^c	0.78±0.25 ^c	0.49±0.47 ^c	
2-Butanone	710.8	78-93-3	RI, MS	4.09±0.56 ^a	2.39±0.54 ^b	2.62±0.44 ^b	0.76±0.39 ^c	-	1.38±0.97 ^{bc}	-	-	-	
3-Undecanone	1578.1	2216-87-7	RI	0.01±0.01	0.02±0.00	-	-	-	-	-	-	-	
3-Nonone	1478.2	925-78-0	RI, MS	-	1.22±0.20	-	-	-	-	-	-	-	
2,3-Octadione	1416.8	585-25-1	RI, MS	-	1.22±0.20 ^a	-	-	-	0.06±0.06 ^b	-	0.06±0.02 ^b	-	
2,3-Pentanedione	1552.4	600-14-6	RI, MS	-	-	0.00±0.00 ^b	-	-	-	0.02±0.02 ^{ab}	0.02±0.01 ^{ab}	0.04±0.02 ^a	
6-Methyl-5-heptene-2-one	1425.4	110-93-0	RI, MS	-	1.57±0.43	-	-	-	-	-	-	-	
2-Pentadecanone	1990.5	2345-28-0	RI, MS	-	-	-	0.23±0.03	0.17±0.03	-	-	-	-	
2-Heptanenone	2194.7	2922-51-2	RI, MS	-	-	-	0.04±0.03	0.07±0.04	-	-	-	-	
2-Octanone	1518.9	111-13-7	RI, MS	-	-	-	0.04±0.01	-	0.02±0.03	-	-	-	
3-Pentanone	873.2	96-22-0	RI, MS	-	-	-	0.30±0.40	-	-	-	-	-	
2,5-Hexanedione	1359.3	110-13-4	RI, MS	-	-	-	0.04±0.04	-	-	-	-	-	
Amyl ketone	1968.9	942-92-7	RI	-	-	-	0.06±0.00	-	-	-	-	-	
4-Methyl-2-heptanone	620.6	6137-06-0	RI, MS	-	-	-	-	0.01±0.02	-	-	-	-	
2,3-Butanedione	894.4	431-03-8	RI, MS	-	-	-	-	0.16±0.22	-	-	-	-	

[illegible]

4,6-Dimethyl-2-heptanone	1106.4	19549-80-5	RI	-	-	-	-	-	-	-	-	0.00±0.01
<i>Alcohols</i>												
N-octanol	1574.3	111-87-5	RI, MS	0.45±0.24 ^a	0.55±0.24 ^a	0.64±0.09 ^a	0.04±0.74 ^b	-	-	0.33±0.31 ^{ab}	-	-
Benzyl alcohol	1849.9	100-51-6	RI, MS	0.06±0.05 ^d	0.09±0.01 ^d	0.11±0.03 ^{cd}	-	0.04±0.03 ^d	0.04±0.01 ^d	0.22±0.08 ^{ab}	0.19±0.04 ^{bc}	0.31±0.09 ^a
Phenylethanol	1884.4	1960-12-8	RI, MS	0.31±0.10 ^c	0.59±0.42 ^{bc}	0.95±0.43 ^{bc}	-	0.08±0.02 ^c	0.03±0.02 ^c	1.87±1.56 ^{ab}	0.95±0.29 ^{bc}	2.99±1.04 ^a
2-Ethyl-1-hexanol	1541.1	104-76-7	RI, MS	0.08±0.01 ^a	0.04±0.04 ^b	0.03±0.01 ^b	-	-	-	0.03±0.02 ^b	0.02±0.02 ^b	0.02±0.01 ^b
2-Ethyl-1-decanol	1845.4	21078-65-9	RI, MS	0.02±0.02	0.02±0.01	0.02±0.01	-	-	-	-	-	-
Furfuryl alcohol	1668.4	98-00-0	RI, MS	-	0.05±0.04	0.04±0.01	0.13±0.15	0.24±0.07	-	-	-	-
Ethanol	781.2	64-17-5	RI, MS	3.34±1.01	4.26±2.86	-	1.88±0.64	-	3.9±3.22	-	-	-
2-Butanol	2821.5	78-92-2	RI	0.03±0.04	0.54±0.86	-	-	0.13±0.19	-	0.00±0.00	0.03±0.01	0.09±0.05
Trans-2-octene-1-ol	1600.2	18409-17-1	RI, MS	0.85±0.69	2.27±0.03	-	-	-	-	-	-	-
Hexanol	1443.7	111-27-3	RI, MS	0.58±0.62	-	0.24±0.06	0.05±0.06	0.12±0.02	0.04±0.05	0.10±0.15	0.06±0.02	0.06±0.02
1-Octene-3-ol	1520.9	3391-86-4	RI, MS	-	4.47±0.60 ^a	3.67±0.11 ^b	-	0.23±0.05 ^c	0.39±0.48 ^c	0.61±0.49 ^c	0.37±0.06 ^c	0.59±0.18 ^c
3-Methylthiopropanol	1693.5	505-10-2	RI, MS	-	0.06±0.04	0.08±0.02	-	-	-	-	-	-
5-Nonanol	1742	623-93-8	RI	-	-	0.02±0.00	-	-	-	-	-	-
2-Methoxy-ethanol	786.7	109-86-4	RI, MS	-	-	0.02±0.02	-	-	0.01±0.01	-	-	-
2-Methyl-3-pentanol	1084.4	565-67-3	RI	-	-	0.29±0.26	-	-	-	0.07±0.04	-	0.02±0.02
1-Pentene-3-ol	1158.6	616-25-1	RI, MS	-	-	0.36±0.08 ^a	-	-	-	0.22±0.07 ^b	0.13±0.03 ^b	0.19±0.05 ^b
1-Dodecanol	1936.5	112-53-8	RI, MS	-	-	-	0.03±0.02	0.08±0.03	0.12±0.04	0.02±0.02	0.08±0.12	-
2,3-Butanediol	1577.3	513-85-9	RI, MS	-	-	-	1.69±0.58	1.21±0.87	0.04±0.04	-	2.42±2.68	-
N-pentanol	1328.4	71-41-0	RI, MS	-	-	-	0.09±0.13 ^c	0.41±0.05 ^a	0.26±0.24 ^{ab}	0.16±0.13 ^{ab}	0.06±0.02 ^c	0.09±0.03 ^c
3-Methyl-1-butanol	1190.3	123-51-3	RI, MS	-	-	-	-	0.32±0.29	0.41±0.58	0.25±0.20	0.10±0.06	0.19±0.06
Isopropanol	740.4	67-63-0	RI, MS	-	-	-	0.10±0.12	-	0.06±0.06	-	-	-

1-Butanol	556	71-36-3	RI, MS	-	-	-	0.24±0.33	-	0.16±0.23	0.11±0.08	0.03±0.04	0.07±0.04
1-Heptanol	1517.5	111-70-6	RI, MS	-	-	-	0.11±0.11	-	-	-	0.02±0.02	-
2,2'-Oxydiethanol	1931.3	111-46-6	RI, MS	-	-	-	0.05±0.03	-	-	-	-	-
1,4-Butanediol	784.9	110-63-4	RI	-	-	-	-	0.08±0.11	-	-	-	-
2-Ethyl-1-butanol	1029.3	97-95-0	RI, MS	-	-	-	-	0.02±0.02	-	-	-	-
1-Methoxy-2-propanol	1057.7	107-98-2	RI, MS	-	-	-	-	0.07±0.05	-	0.06±0.04	0.10±0.03	0.12±0.07
2-Methyl-1-propanol	1086.3	78-83-1	RI, MS	-	-	-	-	0.15±0.21	-	0.07±0.06	0.09±0.06	0.15±0.02
N-propanol	1094.6	71-23-8	RI, MS	-	-	-	-	-	0.02±0.02	-	-	-
3-Butene-2-ol	1292	598-32-3	RI	-	-	-	-	-	0.01±0.01	-	-	-
2-Ethoxy-ethanol	1423.2	110-80-5	RI, MS	-	-	-	-	-	0.01±0.01	-	-	-
2-Methyl-3-hexanol	1427.6	617-29-8	RI	-	-	-	-	-	0.03±0.04	-	-	-
4-Octanol	1524.3	589-62-8	RI	-	-	-	-	-	0.00±0.01	-	-	-
1-Nonanol	1733	143-08-8	RI, MS	-	-	-	-	-	0.02±0.02	-	-	-
2-Furan methanol	1733.4	98-00-0	RI, MS	-	-	-	-	-	0.06±0.06	-	-	-
3-(Methylthio)-1-propanol	1793.1	505-10-2	RI, MS	-	-	-	-	-	0.01±0.01	-	-	0.25±0.05
2-(2-Hydroxypropoxy)-1-propanol	1974.7	106-62-7	RI, MS	-	-	-	-	-	0.03±0.03	-	-	-
4-Heptanol	1978.3	589-55-9	RI, MS	-	-	-	-	-	0.03±0.02	-	-	-
2-Methylbenzyl alcohol	2075.6	89-95-2	RI	-	-	-	-	-	0.01±0.01	-	-	-
N-decanol	2259.4	112-30-1	RI, MS	-	-	-	-	-	0.01±0.01	-	-	-
Hexadecanol	2446.8	36653-82-4	RI, MS	-	-	-	-	-	0.08±0.03	-	0.10±0.14	-
2-Pentanol	2702.1	6032-29-7	RI, MS	-	-	-	-	-	0.02±0.02	-	-	0.06±0.06
3-Pentene-2-ol	1146.7	1569-50-2	RI, MS	-	-	-	-	-	-	0.12±0.06 ^a	0.04±0.03 ^b	0.02±0.02 ^b
2,2'-Oxybis ethanol	1895.2	111-46-6	RI, MS	-	-	-	-	-	-	0.01±0.01	0.01±0.01	0.05±0.05

2-Methyl-3-butene-2-ol	1043.1	115-18-4	RI, MS	-	-	-	-	-	-	0.20±0.08	0.10±0.03	0.44±0.33
3-Methyl-3-butene-1-ol	1226	763-32-6	RI, MS	-	-	-	-	-	-	0.01±0.00	0.02±0.01	-
2-Methyl-2-propanol	1210.7	75-65-0	RI, MS	-	-	-	-	-	-	0.01±0.01	-	-
3-Pentanol	1535.9	584-02-1	RI, MS	-	-	-	-	-	-	0.00±0.01	-	-
Octadecyl alcohol	2614.2	112-92-5	RI, MS	-	-	-	-	-	-	0.01±0.02	-	-
Cis-3-hexenol	1430.4	928-96-1	RI, MS	-	-	-	-	-	-	-	0.02±0.03	-
3-Octanol	2469.9	589-98-0	RI, MS	-	-	-	-	-	-	-	-	0.15±0.06
2-Methyl-1-pentene-3-ol	2041.8	2088-7-5	RI, MS	-	-	-	-	-	-	-	-	2.45±0.85
<i>Acids</i>												
Acetic acid	1518.2	64-19-7	RI, MS	4.85±1.29 ^c	4.02±0.99 ^c	9.33±2.90 ^{bc}	19.18±13.77 ^{abc}	29.51±4.20 ^a	25.08±11.85 ^{ab}	4.52±3.14 ^c	9.27±2.91 ^{bc}	12.15±10.56 ^{bc}
Butyrate	1608	107-92-6	RI, MS	2.76±0.80 ^{cd}	1.31±0.20 ^{cd}	1.61±1.53 ^{cd}	5.96±1.42 ^{bc}	0.77±0.26 ^d	3.52±1.93 ^{cd}	11.69±4.08 ^a	5.04±0.58 ^{bcd}	8.45±4.93 ^{ab}
Caproic acid	1823.1	142-62-1	RI, MS	6.11±0.93 ^b	2.65±0.73 ^c	3.19±0.61 ^c	3.04±0.69 ^c	0.95±0.28 ^c	0.50±0.15 ^c	-	12.25±2.07 ^a	11.4±3.04 ^a
Octanoic acid	2036.1	124-07-2	RI, MS	0.76±0.07 ^b	0.10±0.02 ^b	0.09±0.01 ^b	0.87±0.05 ^b	0.83±0.27 ^b	0.28±0.00 ^b	8.31±2.49 ^a	7.60±1.68 ^a	7.04±2.58 ^a
Decanoic acid	2240.1	334-48-5	RI, MS	0.12±0.07 ^c	0.07±0.02 ^c	0.26±0.10 ^c	1.21±0.44 ^{bc}	2.10±0.15 ^{abc}	0.38±0.27 ^c	3.32±2.97 ^{ab}	4.03±1.22 ^a	3.77±1.42 ^a
2-Methylpropionic acid	1575.4	79-31-2	RI, MS	3.98±0.97 ^{abc}	5.03±1.46 ^{abc}	5.41±1.19 ^{abc}	12.60±1.13 ^a	1.25±0.35 ^c	3.59±1.98 ^{bc}	10.97±10.66 ^{ab}	-	7.20±6.56 ^{abc}
3-Methylbutanoic acid	1649.5	503-74-2	RI, MS	10.14±2.16 ^a	11.90±5.29 ^a	10.02±0.96 ^a	-	1.73±0.56 ^b	3.59±1.97 ^b	0.02±0.01 ^b	-	-
Propionic acid	1561.6	1979-9-4	RI, MS	0.07±0.12	0.99±1.58	-	1.20±0.43	0.16±0.03	0.76±0.23	0.40±0.17	-	-
Phenylacetic acid	2472.8	103-82-2	RI, MS	-	0.01±0.00 ^b	0.05±0.02 ^b	0.80±0.35 ^a	0.47±0.33 ^{ab}	-	-	-	-
4-Methylvaleric acid	1781.2	646-07-1	RI, MS	-	0.03±0.01 ^{ab}	0.03±0.01 ^{ab}	-	0.05±0.01 ^{ab}	-	0.05±0.03 ^{ab}	0.03±0.01 ^b	0.07±0.03 ^a
Heptanic acid	1929.9	111-14-8	RI, MS	0.12±0.01	-	-	0.27±0.01	0.39±0.28	-	0.52±0.43	-	-
Valeric acid	2140.4	109-52-4	RI, MS	-	0.03±0.01 ^c	-	0.73±0.14 ^c	0.28±0.11 ^a	0.08±0.06 ^b	-	0.00±0.00 ^c	-
Formic acid	1117.6	64-18-6	RI, MS	-	-	0.03±0.03	0.65±0.66	-	-	-	-	-
Nonanoic acid	2045.1	112-05-0	RI, MS	-	-	0.03±0.01 ^c	0.25±0.09 ^b	0.41±0.08 ^a	-	0.01±0.01 ^c	-	0.01±0.00 ^c

9-Decenoic acid	2196.1	14436-32-9	RI, MS	-	-	0.02±0.00	-	-	-			
Dodecanoic acid	2277.7	143-07-7	RI, MS	-	-	0.02±0.02 ^c	1.07±0.42 ^b	2.00±0.47 ^a	0.34±0.09 ^c	-	-	0.18±0.18 ^c
3-Methylvaleric acid	1718.8	105-43-1	RI	-	-	0.01±0.00	0.02±0.02	0.03±0.02	-	-	-	-
3-Methyl-2-butenic acid	1595.8	541-47-9	RI, MS	-	-	0.01±0.01	-	-	-	0.02±0.01	0.01±0.01	-
Oxalate	1130.5	144-62-7	RI	-	-	-	0.01±0.01	0.01±0.02	0.03±0.02	-	-	-
Tridecanoic acid	2468.2	638-53-9	RI, MS	-	-	-	0.09±0.01	0.08±0.05	0.03±0.04	-	-	-
Tetradecanoic acid	2580.4	544-63-8	RI, MS	-	-	-	7.99±2.45 ^b	15.58±3.35 ^a	0.22±0.07 ^c	-	-	0.05±0.09 ^c
Pentadecanoic acid	2623.8	1002-84-2	RI, MS	-	-	-	0.06±0.02	0.62±0.21	0.58±0.51			
Hexadecylic acid	2704	1957-10-3	RI, MS	-	-	-	0.55±0.19	0.48±0.12	17.27±24.18	0.29±0.28	-	1.63±2.81
2-Methylacrylic acid	1116.3	79-41-4	RI	-	-	-	0.03±0.04	0.14±0.06	-	-	-	-
2-Acrylic acid	1614.3	1979-10-7	RI	-	-	-	0.06±0.05	0.29±0.03	-	-	-	-
Heptadecanoic acid	2818.8	506-12-7	RI, MS	-	-	-	0.59±0.37	1.29±0.29	-	-	-	-
3-Butyric acid	601.2	2345-51-9	RI	-	-	-	-	0.07±0.04	0.01±0.01	-	-	-
9-Tetraenoic acid	2608.4	544-64-9	RI, MS	-	-	-	-	0.54±0.27	0.36±0.27	-	-	-
2-Methyl butyric acid	1637.8	116-53-0	RI, MS	-	-	-	8.44±2.98	-	-	-	4.99±1.32	5.29±4.65
Benzoic acid	2345.7	65-85-0	RI, MS	-	-	-	0.90±0.99	-	-	-	-	-
Octadecanoic acid	2896.1	1957-11-4	RI, MS	-	-	-	12.75±4.57 ^a	-	-	0.02±0.02 ^b	-	0.34±0.31 ^b
3-Methylvaleric acid	1755.7	105-43-1	RI	-	-	-	-	0.03±0.02 ^b	-	0.08±0.04 ^a	0.04±0.01 ^{ab}	0.07±0.02 ^{ab}
3-Decenoic acid	2274.6	15469-77-9	RI	-	-	-	-	0.27±0.29	-	-	-	-
Undecanonic acid	2365.1	112-37-8	RI, MS	-	-	-	-	8.77±12.15	-	-	-	-
Linoleic acid	2859.6	60-33-3	RI, MS	-	-	-	-	3.15±0.63	-	-	-	-
2-Ethyl-hexanoic acid	2042	149-57-5	RI, MS	-	-	-	-	-	0.07±0.01	-	-	0.00±0.00
2-Methyl-4-pentenoic acid	2459.8	1575-74-2	RI	-	-	-	-	-	0.26±0.12	-	-	-

Cis-9-hexadecenoic acid	2899.3	373-49-9	RI, MS	-	-	-	-	-	12.06±5.00	-	-	-
2-Octenoic acid	2153.6	1470-50-4	RI	-	-	-	-	-	-	0.07±0.08	-	-
2-Methyl-2-butenic acid	2376.2	13201-46-2	RI	-	-	-	-	-	-	0.02±0.01	-	-
3-Butenoic acid	1731.2	625-38-7	RI	-	-	-	-	-	-	-	0.00±0.00	-
4-Hexenoic acid	1938.7	35194-36-6	RI, MS	-	-	-	-	-	-	-	-	0.01±0.01
4-Pentenoic acid	1757.3	591-80-0	RI	-	-	-	-	-	-	-	-	0.06±0.02
Ethyl-9,12-octadienoic acid	2454.9	7619-8-1	RI, MS	-	-	-	-	0.77±0.21	-	-	-	-
<i>Esters</i>												
Methyl hexanoate	1174.7	106-70-7	RI, MS	2.75±2.10 ^a	1.11±1.44 ^{ab}	1.12±0.83 ^{ab}	-	-	-	-	-	-
Ethyl hexanoate	1246.2	123-66-0	RI, MS	0.50±0.56 ^a	0.21±0.14 ^{ab}	0.16±0.03 ^{ab}	-	-	0.00±0.00 ^b	0.00±0.00 ^b	0.02±0.01 ^b	0.01±0.00 ^b
Methyl octanoate	1479.9	111-11-5	RI, MS	0.73±0.45	0.42±0.46	0.55±0.29	-	-	-	-	-	-
Ethyl caproate	1513.7	106-32-1	RI, MS	0.15±0.09	0.11±0.06	0.11±0.03	-	-	-	0.02±0.01	0.05±0.02	0.24±0.33
Methyl decanoate	1592.7	110-42-9	RI, MS	0.09±0.04 ^{ab}	0.11±0.10 ^a	0.13±0.01 ^a	-	-	-	-	-	0.01±0.00 ^b
Ethyl decanoate	1629.8	110-38-3	RI, MS	0.03±0.02 ^b	0.06±0.03 ^b	0.11±0.06 ^a	-	0.01±0.01 ^b	-	0.01±0.00 ^b	0.01±0.01 ^b	0.01±0.00 ^b
Dehydropropionolactone	1990.4	2381-87-5	RI, MS	0.02±0.01 ^c	0.02±0.01 ^c	0.09±0.02 ^{bc}	0.05±0.03 ^c	-	0.01±0.01 ^c	0.57±0.50 ^b	0.58±0.20 ^b	1.51±0.55 ^a
Methyl isovalerate	1010.9	556-24-1	RI, MS	2.59±1.59	1.54±0.79	1.53±1.52	-	-	-	-	-	-
Methyl phenylacetate	1737	101-41-7	RI	0.07±0.04	0.05±0.03	0.18±0.18	-	-	-	-	-	-
Panthenolide	2000	599-04-2	RI, MS	-	0.35±0.06 ^b	2.16±0.56 ^a	-	0.17±0.03 ^b	-	-	-	-
Methyl butyrate	924	623-42-7	RI, MS	0.64±0.26	-	0.59±0.75	-	-	-	-	-	-
Ethyl phenylpropionate	1862.3	2021-28-5	RI, MS	0.03±0.01	0.01±0.00	-	-	-	-	-	-	-
Ethyl acetate	686.1	141-78-6	RI, MS	-	0.67±0.22	0.59±0.18	-	-	-	-	-	-
Ethyl phenylacetate	1764.2	101-97-3	RI, MS	-	0.02±0.00	0.03±0.01	-	-	-	-	0.03±0.01	0.04±0.01
Ethyl dodecanoate	1835	106-33-2	RI, MS	-	0.00±0.00	0.01±0.00	-	-	-	-	-	0.07±0.06

Propyl acetate	890	109-60-4	RI, MS		5.93±2.28 ^a	-	-	0.01±0.02 ^b	0.26±0.21 ^b	—	0.56±0.23 ^b	0.44±0.25 ^b
Methyl-2-methylbutyrate	1002.2	868-57-5	RI	-	0.45±0.25	-	-	-	-	-	-	-
Butyl acetate	1060.6	123-86-4	RI, MS	-	0.59±0.13 ^a	-	-	-	-	0.02±0.01 ^b	0.04±0.01 ^b	0.04±0.02 ^b
Methyl nonanoate	1542.2	1731-84-6	RI, MS	-	0.01±0.01	-	-	-	-	-	-	-
Ethyl-2-hydroxypropionate	2629.5	97-64-3	RI	-	0.00±0.00	-	-	0.00±0.00	0.02±0.02	-	-	-
Butyrolactone	1584.3	96-48-0	RI, MS	-	-	3.63±0.87 ^b	-	0.67±0.49 ^b	2.06±1.14 ^b	-	7.83±3.63 ^a	-
Methyl valerate	1071.5	624-24-8	RI, MS	-	-	0.05±0.02	-	-	-	-	-	-
Isobutyl acetate	1058.7	110-19-0	RI, MS	-	-	0.31±0.11 ^a	0.02±0.02 ^b	0.02±0.03 ^b	-	-	-	-
Methyl-2-methylbutyrate	1002.1	868-57-5	RI	-	-	0.92±1.09	-	-	-	-	-	-
2-Hydroxy-γ-butyrolactone	2130.3	19444-84-9	RI, MS	-	-	0.00±0.00	-	0.08±0.01	-	-	-	-
Ethyl tetradecanoate	2021.3	124-06-1	RI, MS	-	-	-	0.02±0.02	0.05±0.01	0.03±0.03	-	-	-
Ethyl octadecanoate	2402.1	111-61-5	RI, MS	-	-	-	0.14±0.11 ^a	0.16±0.01 ^a	0.02±0.02 ^{ab}	-	-	0.00±0.00 ^b
Ethyl linoleate	2454.6	544-35-4	RI, MS	-	-	-	0.15±0.08	-	0.03±0.02			
Methyl tetradecanoate	1980.2	124-10-7	RI, MS	-	-	-	0.03±0.01	0.02±0.00	-	-	-	0.01±0.00
Methyl hexadecate	2181.7	112-39-0	RI, MS	-	-	-	0.23±0.06 ^{ab}	0.26±0.09 ^a	-	0.07±0.03 ^c	0.11±0.07 ^{bc}	0.08±0.04 ^c
Ethyl-9-hexadecanoate	2239.4	54546-22-4	RI, MS	-	-	-	0.05±0.04	0.14±0.10	-	-	-	-
Methyl formate	2924	107-31-3	RI	-	-	-	-	0.65±0.47	0.01±0.01	-	-	-
Ethyl palmitate	2218.6	628-97-7	RI, MS	-	-	-	-	0.30±0.22 ^a	0.06±0.04 ^b	0.01±0.00 ^b	0.02±0.03 ^b	0.04±0.01 ^b
Methyl acetate	1350	79-20-9	RI, MS	-	-	-	-	0.15±0.11	0.00±0.00	-	-	-
Ethyl butyrate	1062.7	4341-76-8	RI	-	-	-	-	0.00±0.00	-	-	-	-
Octyl hexadecate	2616.9	16958-85-3	RI	-	-	-	-	0.17±0.14	-	-	-	-
Methyl oleate	2384.4	112-62-9	RI, MS	-	-	-	-	0.34±0.12	-	-	-	-
Ethyl oleate	2416.3	111-62-6	RI, MS	-	-	-	-	0.51±0.14	-	-	-	-

Cyclopentyl-4-ethylbenzoate	1742.9	14779-78-3	RI, MS	-	-	-	-	4.93±3.18	-	-	-	-
2-Methyl acrylate	880.4	922-67-8	RI	-	-	-	-	-	0.01±0.01	-	-	-
2-Vinyl acrylate	1149.4	2177-18-6	RI	-	-	-	-	-	0.01±0.01	-	-	-
Vinyl formate	1261.1	692-45-5	RI	-	-	-	-	-	0.27±0.38	-	-	-
Butyl valerate	1553.7	591-68-4	RI, MS	-	-	-	-	-	0.00±0.00	-	-	-
Heptyl formate	1563	112-23-2	RI	-	-	-	-	-	0.65±0.56	-	-	-
Butyl-2-butenolate	1573.2	7299-91-4	RI, MS	-	-	-	-	-	0.02±0.02	-	-	-
Ethyl-dodecanoate	1700.3	106-33-2	RI, MS	-	-	-	-	-	0.00±0.00	-	-	-
Triethyl phosphate	1741.9	78-40-0	RI, MS	-	-	-	-	-	0.02±0.01	-	-	-
Methyl-2-hydroxy-2-methyl-propionate	1967.4	2110-78-3	RI	-	-	-	-	-	0.03±0.04	-	-	-
Isopropyl myristate	2124.9	110-27-0	RI, MS	-	-	-	-	-	0.07±0.07	-	-	-
Diethyl-1,4-phthalate	2418	636-09-9	RI	-	-	-	-	-	0.16±0.14	-	-	-
Diisobutyl phthalate	2581.6	84-69-5	RI, MS	-	-	-	-	-	0.02±0.02	-	-	-
Dibutyl phthalate	2706.9	84-74-2	RI, MS	-	-	-	-	-	0.08±0.06	-	-	0.96±0.56
Ethyl-p-hydroxybenzoate	2915.7	120-47-8	RI, MS	-	-	-	-	-	0.03±0.02	-	-	-
Decyl decanoate	2429.8	1654-86-0	RI, MS	-	-	-	-	-	-	0.10±0.03	0.10±0.05	0.05±0.03
Ethyl-2-methyl butyrate	1054.3	7452-79-1	RI, MS	-	-	-	-	-	-	0.02±0.02	0.03±0.03	0.07±0.03
Ethyl-3-methyl butyrate	1066.6	108-64-5	RI, MS	-	-	-	-	-	-	0.01±0.01	0.01±0.00	0.02±0.01
Dimethyl phthalate	2200.6	131-11-3	RI, MS	-	-	-	-	-	-	0.05±0.01	0.04±0.02	0.04±0.02
Diethyl phthalate	2265.8	84-66-2	RI, MS	-	-	-	-	-	-	0.03±0.01	0.04±0.03	0.04±0.01
2-Hydroxy-γ- Butyrolactone	2077.7	19444-84-9	RI, MS	-	-	-	-	-	-	0.03±0.02 ^b	0.04±0.02 ^b	0.15±0.06 ^a
Butyl butyrate	1188	109-21-7	RI, MS	-	-	-	-	-	-	0.00±0.00	0.00±0.00	-
Butyl 2-butenolate	1194.6	7299-91-4	RI, MS	-	-	-	-	-	-	0.00±0.00	0.00±0.00	-

Ethyl heptanate	1378.1	106-30-9	RI, MS	-	-	-	-	-	-	-	0.00±0.00	0.00±0.00
Dimethyl adipate	1745.7	627-93-0	RI, MS	-	-	-	-	-	-	-	-	-
Ethyl-2-hydroxypropionate	1389.6	97-64-3	RI, MS	-	-	-	-	-	-	-	0.00±0.00	0.00±0.00
Methyl acetate	1198.2	79-20-9	RI, MS	-	-	-	-	-	-	0.02±0.02	-	-
Ethyl-4-ethoxybenzoate	2082.2	23676-09-7	RI, MS	-	-	-	-	-	-	0.06±0.04	-	-
Ethyl tetradecanoate	1972.1	124-06-1	RI, MS	-	-	-	-	-	-	0.02±0.01	-	-
Ethyl propionate	1105.5	105-37-3	RI, MS	-	-	-	-	-	-	-	0.00±0.00	-
Methyl heptanate	1008.7	106-73-0	RI, MS	-	-	-	-	-	-	-	0.00±0.00	-
Butyl-3-methyl butyrate	1009.8	106-27-4	RI, MS	-	-	-	-	-	-	-	0.02±0.00	-
Ethyl valerate	1119.3	539-82-2	RI, MS	-	-	-	-	-	-	-	-	0.00±0.00
Methyl-2-butyrate	1018.2	623-43-8	RI, MS	-	-	-	-	-	-	-	-	0.01±0.01
Butyl-2-methyl butyrate	1153.5	15706-73-7	RI, MS	-	-	-	-	-	-	-	-	0.00±0.00
Ethyl-2-methylvalerate	1764.1	39255-32-8	RI, MS	-	-	-	-	-	-	-	-	0.03±0.01
<i>Alkane</i>												
Pentane	494	109-66-0	RI	0.04±0.04	0.05±0.04	0.09±0.01	0.06±0.05	0.18±0.26	0.03±0.04	0.05±0.01	0.03±0.01	0.03±0.00
N-hexane	505.2	110-54-3	RI	0.01±0.01 ^b	0.02±0.00 ^b	0.01±0.01 ^b	-	-	-	34.46±17.91 ^a	32.53±10.52 ^a	16.18±4.26 ^b
Decane	999.6	124-18-5	RI	4.25±1.05 ^a	2.52±0.73 ^b	1.57±0.25 ^b	-	0.03±0.04 ^c	0.39±0.56 ^c	-	-	0.24±0.11 ^c
Undecanone	1023.9	1120-21-4	RI	0.10±0.04	0.08±0.03	0.02±0.02	-	-	0.02±0.03	-	-	-
Dodecane	1194.2	112-40-3	RI	2.36±1.50 ^a	2.69±1.98 ^a	0.24±0.04 ^b	-	-	0.01±0.01 ^b	0.11±0.05 ^b	0.18±0.05 ^b	0.04±0.02 ^b
Tridecane	1054.1	629-50-5	RI	0.26±0.03 ^a	0.13±0.06 ^b	0.13±0.06 ^b	-	-	-	0.02±0.01 ^c	0.02±0.01 ^c	0.02±0.01 ^c
Tetradecane	1499.9	629-59-4	RI	0.62±0.22 ^a	0.40±0.04 ^a	0.44±0.17 ^a	-	-	-	0.02±0.01 ^b	0.13±0.03 ^b	0.17±0.12 ^b
Cetane	1599.7	544-76-3	RI	0.04±0.04 ^{ab}	0.01±0.01 ^b	0.04±0.01 ^{ab}	0.05±0.05 ^{ab}	0.06±0.06 ^{ab}	0.06±0.08 ^{ab}	0.14±0.08 ^{ab}	0.17±0.11 ^a	-
3-Methylheptane	565.6	589-81-1	RI	0.12±0.11	0.12±0.05	0.09±0.01	-	-	-	-	-	-

3-Methylundecanone	1159.8	1002-43-3	RI	1.35±0.79	1.35±1.10	0.08±0.02	-	-	-	-	-	-
5-Methylnonane	841.8	15869-85-9	RI	0.09±0.01 ^a	0.05±0.00 ^b	0.04±0.01 ^b	-	-	-	-	-	-
4-Methyltetradecane	1462	25117-24-2	RI	0.06±0.01	0.05±0.02	0.06±0.01	-	-	-	-	-	-
3-Methyltridecane	1467.8	6418-41-3	RI, MS	0.48±0.07 ^a	0.21±0.07 ^b	0.28±0.06 ^b	-	-	-	-	0.02±0.01 ^c	-
4-Ethyloctane	829.5	15869-86-0	RI	0.14±0.01	0.11±0.04	-	-	-	-	-	-	-
3-Methylpentadecane	1584	2882-96-4	RI	0.06±0.01	0.04±0.02	-	-	-	-	-	-	-
Octane	585.8	111-65-9	RI	-	0.31±0.08	0.21±0.04	0.47±0.36	-	0.07±0.08	-	-	-
4-Ethyldecane	1134.1	1636-44-8	RI	0.13±0.02	-	-	-	-	-	-	-	-
2,5-Dimethyldecane	1050.2	17312-50-4	RI	0.45±0.40	-	-	-	-	-	-	-	-
5-Methylundecanone	1141.9	1632-70-8	RI	0.37±0.23	-	-	-	-	-	-	-	-
2-Methylcetane	1565.3	1560-92-5	RI	0.01±0.01	-	-	-	-	-	-	-	-
4-Ethylheptane	1031.1	2216-32-2	RI	-	0.19±0.18	-	-	-	-	-	-	-
5-Propyldecane	1132.5	17312-62-8	RI	-	0.14±0.04	-	-	-	-	-	-	-
4-Methyldodecane	1156.6	6117-97-1	RI	-	0.09±0.10	-	-	-	-	-	-	-
3-Methyloctane	1098.5	2216-33-3	RI	-	-	0.22±0.21	-	-	-	-	-	-
5-Methyldecane	1044.4	13151-35-4	RI	-	-	0.30±0.07	-	-	-	-	-	-
3,3-Dimethyloctane	1043.8	4110-44-5	RI	-	-	0.08±0.02	-	-	-	-	-	-
2,2-Dimethylbutane	1057	75-83-2	RI, MS	-	-	0.09±0.13	-	-	-	-	-	-
4-Ethyltetradecane	1576.5	55045-14-2	RI	-	-	0.03±0.00	-	-	-	-	-	-
2,5-Dimethyltridecane	1400.3	56292-66-1	RI	-	-	0.11±0.03	-	-	-	-	-	-
2,3,5-Trimethylheptane	1102.4	20278-85-7	RI, MS	-	-	0.33±0.07	-	-	-	-	-	-
Butane	670.9	106-97-8	RI	-	-	-	-	0.01±0.01	0.01±0.01	-	-	-

Heptane	517.4	142-82-5	RI	-	-	-	0.12±0.14 ^b	-	0.02±0.03 ^b	1.09±0.85 ^a	0.21±0.28 ^{ab}	0.80±0.70 ^{ab}
Propane	719.1	74-98-6	RI	-	-	-	0.28±0.33	0.00±0.00	-	-	-	-
3-Methyl-heptadecane	1754	6418-44-6	RI, MS	-	-	-	0.01±0.01	0.03±0.03	-	-	-	-
2,3-Dimethylbutane	716.2	79-29-8	RI	-	-	-	-	0.18±0.25	-	-	-	-
3-Methyl-heptane	612.5	589-81-1	RI	-	-	-	-	-	0.10±0.10	-	-	-
2-Nitro-propane	783	79-46-9	RI	-	-	-	-	-	0.00±0.00	-	-	-
Neopentane	1486.7	463-82-1	RI	-	-	-	-	-	0.02±0.02	-	-	-
4-Propyl-heptane	2821.1	55044-10-5	RI	-	-	-	-	-	0.02±0.03	-	-	-
Heptadecane	1634.5	629-78-7	RI	-	-	-	-	-	-	0.06±0.03	0.06±0.03	0.03±0.01
Octadecane	1731.4	593-45-3	RI	-	-	-	-	-	-	0.04±0.01		
2-Methyltridecane	1410.7	1560-96-9	RI	-	-	-	-	-	-	-	0.00±0.00	0.01±0.00
2-Undecanone	1569.3	112-12-9	RI, MS	-	-	-	-	-	-	-	0.01±0.00	-
4-Methyldecane	1085.8	2847-72-5	RI	-	-	-	-	-	-	-	-	0.01±0.00
4-Methylundecanone	1136.2	2980-69-0	RI, MS	-	-	-	-	-	-	-	-	0.01±0.00
3-Methylheptadecane	1701.2	6418-44-6	RI, MS	-	-	-	-	-	-	-	-	0.01±0.00
<i>Olefin</i>												
1-Decene	1035.6	872-05-9	RI, MS	0.44±0.07 ^a	-	0.14±0.04 ^b	-	-	-	0.00±0.00 ^c	0.01±0.00 ^c	0.01±0.01 ^c
2-Octene	1283.4	111-67-1	RI, MS	0.62±0.58	-	-	-	-	-	-	-	-
5-Methyl-2-heptene	1760	22487-87-2	RI	0.02±0.00	-	-	-	-	-	-	-	-
3,5,5-Trimethyl-1-hexene	549.7	4316-65-8	RI	0.00±0.00	-	-	-	-	-	-	-	-
3-Methyl-3-heptene	578.5	7300-03-0	RI	-	-	0.01±0.01	-	-	-	-	-	-
1,3-Octadiene	1486.2	1002-33-1	RI, MS	-	-	0.03±0.01	-	-	-	-	-	-
Propylene	1150.7	115-07-1	RI	-	-	-	0.01±0.01	0.42±0.56	7.4±10.44	-	-	-

1-Heptene	535.9	592-76-7	RI, MS	-	-	-	0.06±0.06	-	-	-	-	-
1-Pentene	1324.6	109-67-1	RI	-	-	-	0.07±0.10	-	-	-	-	-
1-Hexene	1451.5	592-41-6	RI	-	-	-	0.05±0.07	-	-	-	-	-
7-Methyl-1-octene	1688.1	13151-06-9	RI	-	-	-	-	-	0.00±0.01	-	-	-
9-Octadecene	2436.8	5090-41-5	RI	-	-	-	-	-	0.15±0.09	-	-	-
1-Dodecene	1219.5	112-41-4	RI, MS	-	-	-	-	-	-	0.01±0.00	0.01±0.00	0.01±0.01
2-Methyl-1-hexene	1030.2	6094-2-6	RI	-	-	-	-	-	-	0.01±0.01	-	-
2,4-Dimethyl-2-pentene	693.3	625-65-0	RI	-	-	-	-	-	-	0.07±0.01	-	-
3-Heptane	1643.3	592-78-9	RI	-	-	-	-	-	-	0.07±0.03	-	-
1-Tetradecene	1509	1120-36-1	RI, MS	-	-	-	-	-	-	-	0.00±0.00	-
6-Methyl-1-octene	1181.8	13151-10-5	RI	-	-	-	-	-	-	-	0.01±0.01	-
5-Dodecene	1267.3	7206-28-2	RI	-	-	-	-	-	-	-	-	0.00±0.00
2,2-Dimethyl-3-octene	1000.4	86869-76-3	RI	-	-	-	-	-	-	-	-	0.02±0.01
<i>Pyrazine</i>												
Trimethyl pyrazine	1507.7	14667-55-1	RI, MS	1.29±0.89 ^b	-	2.68±0.81 ^a	0.01±0.01 ^c	0.44±0.32 ^c	0.23±0.14 ^c	0.23±0.14 ^c	0.13±0.05 ^c	0.68±0.20 ^{bc}
2,3,5,6-Tetramethylpyrazine	1532.2	1124-11-4	RI, MS	0.46±0.65	-	0.60±0.53	-	-	0.10±0.07	0.07±0.04	0.04±0.04	0.31±0.40
2,3-Dimethylpyrazine	1434.9	5910-89-4	RI, MS	0.43±0.25 ^a	-	0.21±0.06 ^b	-	-	0.03±0.03 ^b	0.04±0.02 ^b	0.01±0.01 ^b	0.04±0.02 ^b
2,6-Diethylpyrazine	1516.6	13067-27-1	RI, MS	-	-	0.09±0.08 ^a	-	0.01±0.01 ^b	0.01±0.01 ^b	0.00±0.00 ^b	0.00±0.00 ^b	0.00±0.00 ^b
2-Methylpyrazine	1314.6	109-08-0	RI, MS	0.29±0.26	-	-	-	-	0.03±0.04	-	-	-
2,6-Dimethylpyrazine	1417	108-50-9	RI, MS	0.61±0.58 ^{ab}	-	2.91±2.58 ^a	-	-	0.14±0.10 ^b	-	0.10±0.09 ^b	0.83±0.86 ^{ab}
2-Ethyl-6-methyl-pyrazine	1493.5	13925-03-6	RI, MS	-	-	0.20±0.03 ^a	-	-	-	0.01±0.01 ^b	0.01±0.00 ^b	0.04±0.02 ^b
Methyl pyrazine	1281.4	109-08-0	RI, MS	-	-	0.25±0.22 ^b	-	-	-	0.05±0.03 ^b	0.03±0.01 ^b	0.06±0.03 ^a
Vinyl pyrazine	1551.3	4177-16-6	RI, MS	-	-	-	-	-	0.13±0.18	-	-	-

Ethyl pyrazine	2172.9	13925-00-3	RI, MS	-	-	-	-	-	0.07±0.09	-	-	0.83±0.86
2,5-Dimethylpyrazine	1354.5	123-32-0	RI, MS	-	-	-	-	-	-	0.40±0.18	0.00±0.00	-
Pyrazine	1865.9	290-37-9	RI, MS	-	-	-	-	-	-	-	0.00±0.00	-
2-Acetyl-3-methylpyrazine	1623.8	23787-80-6	RI, MS	-	-	-	-	-	-	-	-	0.00±0.00
2,3-Dimethyl-5-ethyl pyrazine	1504	15707-34-3	RI, MS	-	-	-	-	-	-	-	-	0.02±0.01
2,3,5-Trimethyl-6-ethylpyrazine	1530.5	17398-16-2	RI, MS	-	-	-	-	-	-	-	-	0.01±0.01
<i>Furan</i>												
2-Pentylfuran	1232.8	3777-69-3	RI, MS	0.22±0.21 ^{ab}	0.34±0.18 ^{ab}	0.58±0.49 ^a	0.26±0.22 ^{ab}	0.5±0.26 ^{ab}	0.43±0.38 ^{ab}	0.04±0.03 ^{ab}	0.02±0.01 ^b	0.08±0.02 ^{ab}
2-Ethyl-furan	820.3	3208-16-0	RI, MS	-	0.04±0.05	-	-	-	-	-	-	-
5-Methyl-2-acetylfuran	1590.7	1193-79-9	RI, MS	-	0.01±0.01	-	-	-	-	-	-	-
2-Propyl-furan	1515.9	4229-91-8	RI, MS	-	-	-	0.01±0.01	0.01±0.01	-	0.01±0.01	0.01±0.01	-
2-Hexyl-furan	1581	3777-70-6	RI, MS	-	-	-	0.03±0.04	-	0.01±0.01	-	-	-
2-Methylfuran	2105.1	534-22-5	RI, MS	-	-	-	0.00±0.00	-	-	-	-	-
Furan	1086.8	110-00-9	RI, MS	-	-	-	-	-	0.05±0.06	0.00±0.00	-	-
2-Ethyl-5-methyl-furan	1584.9	1703-52-2	RI	-	-	-	-	-	0.01±0.01	-	-	-
2,5-Dibutylfuran	1161	72636-53-4	RI	-	-	-	-	-	-	0.00±0.00	0.00±0.00	0.00±0.00
3-Methyl-furan	1470.8	930-27-8	RI, MS	-	-	-	-	-	-	0.00±0.00	-	-
<i>Sulfide</i>												
Dimethyl trisulfide	1455.4	3658-80-8	RI, MS	0.02±0.02 ^b	0.01±0.00 ^b	0.01±0.00 ^b	1.21±0.99 ^a	0.65±0.10 ^{ab}	0.34±0.19 ^b	0.00±0.00 ^b	-	0.00±0.00 ^b
Dimethyl sulfide	557.9	75-18-3	RI, MS	-	-	0.05±0.00 ^a	-	0.01±0.01 ^b	-	0.00±0.00 ^b	-	-
Dimethyl disulfide	1048.4	624-92-0	RI, MS	-	-	-	0.95±1.34	1.71±2.39	-	-	-	-
Dimethyl tetrasulfide	1693.9	5756-24-1	RI, MS	-	-	-	-	0.03±0.02	-	-	-	-
Methyl mercaptan	510.6	74-93-1	RI, MS	-	-	-	0.38±0.28	0.16±0.12	-	-	-	-

Notes: “-” indicates that the substance is not detected whiles “0.00” indicates that the value is less than 0.01. All the value is expressed as the percentage of peak area (%). The relative percentage is the ratio of the volatile component to the total volatile aroma compounds in one aging time.

¹ Volatile component were identified by RI and MS

² Retention index, calculated according to peak time of series alkanes under the same gas phase conditions

^{a, b, c, d} Means with different letters with in a row differ significantly ($P < 0.05$)