

**Supplementary Table S1** The volatile flavor compounds identified by GC/MS of apple juice and cider samples fermented with different non-*Saccharomyces* yeast

No.	Name	Formul a	FAJ	Cider (μg/L)				
				Sc	Sc+Zb	Sc+Rm	Sc+Km	Sc+Dh
C0 1	2,3-Butanediol	C <sub>4</sub> H <sub>10</sub> O <sub>2</sub>	nd	16.05 ± 2.84a	24.41 ± 28.74a	36.15 ± 12.52a	16.70 ± 3.38a	39.08 ± 6.83a
C0 2	2-Furanmethanol	C <sub>5</sub> H <sub>6</sub> O <sub>2</sub>	nd	nd	nd	3.91 ± 5.53	nd	nd
C0 3	1-Butanol, 3-methyl-	C <sub>5</sub> H <sub>12</sub> O	20.22 ± 0.57b	1954.68 ± 412.61a	1860.53 ± 464.12a	1492.47 ± 189.52a	1292.73 ± 35.12a	1510.91 ± 243.64a
C0 4	3-Hexen-1-ol, (Z)-	C <sub>6</sub> H <sub>12</sub> O	8.94 ± 2.04	nd	nd	nd	nd	nd
C0 5	Phenylethyl alcohol	C <sub>8</sub> H <sub>10</sub> O	nd	561.18 ± 40.14a	614.18 ± 213.97a	609.01 ± 187.79a	488.65 ± 277.71a	782.68 ± 165.47a
C0 6	Cyclohexanol, 5- methyl-2-(1- methylethyl)-	C <sub>10</sub> H <sub>20</sub> O	10.49 ± 1.24	nd	nd	nd	nd	nd
C0 7	1-Dodecanol	C <sub>12</sub> H <sub>26</sub> O	nd	nd	nd	nd	3.00 ± 4.25	nd
C0 8	5,9-Undecadien-2-ol, 6,10-dimethyl-	C <sub>13</sub> H <sub>24</sub> O	nd	nd	nd	29.21 ± 9.79	nd	nd
C0 9	2,6,10-Dodecatrien-1- ol, 3,7,11-trimethyl-	C <sub>15</sub> H <sub>26</sub> O	nd	nd	9.12 ± 12.89a	nd	nd	27.30 ± 20.36a
Alcohols			39.65 ± 1.88b	2531.91 ± 446.04a	2508.24 ± 491.84a	1801.08 ± 248.36a	2359.97 ± 304.51a	2170.75 ± 401.24a
C1	Hexanal	C <sub>6</sub> H <sub>12</sub> O	169.80 ± 1.39	nd	nd	nd	nd	nd



C2 2	9-Decenoic acid, ethyl ester	C <sub>12</sub> H <sub>22</sub> O <sub>2</sub>	nd	51.49 ± 18.59a	43.40 ± 43.98a	30.80 ± 43.55a	10.85 ± 15.34a	78.08 ± 27.72a
C2 3	Decanoic acid, ethyl ester	C <sub>12</sub> H <sub>24</sub> O <sub>2</sub>	nd	683.19 ± 68.36a	665.49 ± 104.18a	690.20 ± 119.92a	579.15 ± 204.16a	845.10 ± 66.20a
C2 4	4-Ethylbenzoic acid, 2-butyl ester	C <sub>13</sub> H <sub>18</sub> O <sub>2</sub>	nd	42.93 ± 4.99b	46.41 ± 15.34b	51.73 ± 14.13a	34.67 ± 6.61b	80.06 ± 27.07a
C2 5	Acetic acid, [4-(1,1-dimethylethyl)phenoxy]-, methyl ester	C <sub>13</sub> H <sub>18</sub> O <sub>3</sub>	nd	nd	166.72 ± 235.78a	90.07 ± 127.38a	nd	nd
C2 6	Tridecanoic acid, 3-hydroxy-, ethyl ester	C <sub>15</sub> H <sub>30</sub> O <sub>3</sub>	nd	nd	nd	10.29 ± 2.63a	3.82 ± 5.40b	12.55 ± 1.83a
C2 7	Nerolidyl acetate	C <sub>17</sub> H <sub>28</sub> O <sub>2</sub>	nd	nd	nd	nd	1.97 ± 2.78	nd
C2 8	Hexadecanoic acid, ethyl ester	C <sub>18</sub> H <sub>36</sub> O <sub>2</sub>	nd	nd	nd	64.06 ± 90.60a	22.75 ± 32.18a	nd
C2 9	Succinic acid, hept-2-yl oct-1-en-3-yl ester	C <sub>19</sub> H <sub>34</sub> O <sub>4</sub>	nd	nd	nd	6.02 ± 4.49a	2.83 ± 4.00a	nd
C3 0	Isopropyl palmitate	C <sub>19</sub> H <sub>38</sub> O <sub>2</sub>	nd	8.68 ± 12.28a	131.29 ± 185.67a	8.08 ± 11.43a	nd	nd
C3 1	1,2-Benzenedicarboxylic acid, butyl octyl ester	C <sub>20</sub> H <sub>30</sub> O <sub>4</sub>	nd	nd	nd	nd	35.83 ± 50.67	nd
C3 2	Benzoic acid, 2-formyl-4,6-dimethoxy-, 8,8-dimethoxyoct-2-yl	C <sub>20</sub> H <sub>30</sub> O <sub>7</sub>	nd	4.46 ± 6.31	nd	nd	nd	nd

## ester

C3 3	Decanoic acid, tetradecyl ester	C <sub>24</sub> H <sub>48</sub> O <sub>2</sub>	nd	nd	nd	nd	nd	150.45 ± 212.76
	Esters		1257.76 ± 4.28d	1970.95 ± 448.70bc	2254.68 ± 321.71ab	1643.29 ± 411.02bc	2894.15 ± 346.26a	2120.68 ± 282.80b
C3 4	Acetic acid	C <sub>2</sub> H <sub>4</sub> O <sub>2</sub>	nd	36.33 ± 6.22b	64.99 ± 33.35ab	71.96 ± 28.79ab	44.19 ± 6.50b	93.69 ± 10.43a
C3 5	Hexanoic acid	C <sub>6</sub> H <sub>12</sub> O <sub>2</sub>	nd	158.21 ± 48.94b	204.63 ± 15.30ab	179.67 ± 20.69b	157.02 ± 68.56b	276.99 ± 43.38a
C3 6	Octanoic acid	C <sub>8</sub> H <sub>16</sub> O <sub>2</sub>	nd	633.88 ± 14.41b	863.37 ± 299.93ab	903.24 ± 231.02ab	838.51 ± 405.44ab	1216.82 ± 207.03a
C3 7	9-Decenoic acid	C <sub>10</sub> H <sub>18</sub> O <sub>2</sub>	nd	24.69 ± 17.49a	32.93 ± 31.28a	42.00 ± 18.40a	nd	62.72 ± 19.75a
C3 8	n-Decanoic acid	C <sub>10</sub> H <sub>20</sub> O <sub>2</sub>	0.42 ± 0.59c	756.94 ± 42.87b	967.74 ± 284.78ab	951.23 ± 229.16ab	883.20 ± 540.30ab	1485.78 ± 192.57a
C3 9	10-Hydroxydecanoic acid	C <sub>10</sub> H <sub>20</sub> O <sub>3</sub>	nd	nd	nd	nd	3.01 ± 4.26	nd
C4 0	Dodecanoic acid	C <sub>12</sub> H <sub>24</sub> O <sub>2</sub>	nd	194.62 ± 28.45a	242.76 ± 40.87a	228.68 ± 35.16a	149.53 ± 115.25a	235.03 ± 166.23a
C4 1	n-Hexadecanoic acid	C <sub>16</sub> H <sub>32</sub> O <sub>2</sub>	nd	nd	16.85 ± 23.83a	180.91 ± 237.56a	19.78 ± 14.71a	16.76 ± 12.02a
C4 2	Oleic Acid	C <sub>18</sub> H <sub>34</sub> O <sub>2</sub>	nd	nd	nd	nd	22.37 ± 18.61	nd
	Acids		0.42 ± 0.59c	1804.68 ±	2393.27 ±	2117.61 ±	3387.78 ±	2557.69 ±

				123.56b	630.46ab	1162.74ab	377.85a	693.43ab
C4 3	2-Buten-1-one, 1-(2,6,6-trimethyl-1,3-cyclohexadien-1-yl)-, (E)-	C <sub>13</sub> H <sub>18</sub> O	5.97 ± 0.58	nd	nd	nd	nd	nd
C4 4	Ketone		5.97 ± 0.58	nd	nd	nd	nd	nd
C4 5	3-Allyl-6-methoxyphenol	C <sub>10</sub> H <sub>12</sub> O <sub>2</sub>	nd	nd	4.62 ± 6.54b	14.69 ± 4.10a	12.79 ± 5.61b	nd
C4 6	2,4-bis(1,1-dimethylethyl)-phenol	C <sub>14</sub> H <sub>22</sub> O	9.60 ± 0.43b	94.88 ± 29.39ab	139.97 ± 73.96a	116.68 ± 53.11ab	120.66 ± 26.71ab	144.21 ± 72.50a
C4 7	2,4-bis(1,1-dimethylethyl)-5-methylphenol	C <sub>15</sub> H <sub>24</sub> O	nd	nd	4.14 ± 5.85	nd	nd	nd
C4 8	Phenols		9.60 ± 0.43b	94.88 ± 29.39ab	148.73 ± 73.96a	133.45 ± 26.71ab	144.21 ± 72.50a	131.37 ± 53.11ab
C4 9	p-Xylene	C <sub>8</sub> H <sub>10</sub>	nd	nd	42.38 ± 13.51a	nd	14.23 ± 10.28b	nd
C5 0	o-Xylene	C <sub>8</sub> H <sub>10</sub>	nd	nd	nd	nd	4.69 ± 6.63	nd
C4 9	Benzene, 1,2,4-trimethyl-	C <sub>9</sub> H <sub>12</sub>	nd	30.81 ± 10.22ab	45.47 ± 9.47a	24.97 ± 9.25b	34.83 ± 7.00ab	22.41 ± 6.74b
C5 0	Benzenes		nd	30.81 ± 10.22bc	87.85 ± 19.05a	53.75 ± 10.87b	22.41 ± 6.74c	24.97 ± 9.25c
C5 0	1-Octene, 3-(methoxymethoxy)-	C <sub>10</sub> H <sub>20</sub> O <sub>2</sub>	nd	nd	5.28 ± 7.46a	4.25 ± 6.01a	nd	8.82 ± 6.24a
	Alkene		nd	nd	5.28 ± 7.46a	nd	8.82 ± 6.24a	4.25 ± 6.01a
	Total		1626.97 ±	6465.48 ±	7416.57 ±	5782.89 ±	8817.34 ±	7058.80 ±

8.80c	502.14ab	1102.77ab	1814.94b	156.60ab	1407.18ab
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Results represent the mean  $\pm$  SD for three independent experiments. Mean values in the same row with the same letters are not significantly different at the 95% confidence level (Duncan's test). ND = not detected. Different letters represent significantly difference at the 95% percent confidence level (Duncan's test). Sc, *S. cerevisiae* monoculture. Sc+Zb, *S. cerevisiae* and *Z. bailii* co-inoculation. Sc+Rm, *S. cerevisiae* and *R. mucilaginosa* co-inoculation. Sc+Km, *S. cerevisiae* and *K. marxianus* co-inoculation. Sc+Dh, *S. cerevisiae* and *D. hansenii* co-inoculation. FAJ, apple juice.

**Supplementary Table S2** Odor-activity values (OVA) of apple juice and cider samples fermented with different non-*Saccharomyces* yeast

No .	Name	Odor	Thresh old		FAJ	OVA (µg/L)			
			value (µg/L )	Sc		Sc+Zb	Sc+Rm	Sc+Km	Sc+Dh
Alcohols									
C0 1	2,3-Butanediol	Butter, cheese aroma	668	nf	0.02 ± 0.00a	0.04 ± 0.04a	0.05 ± 0.02a	0.03 ± 0.01a	0.06 ± 0.01a
C0 2	2-Furanmethanol	Mildly irritating	8000	nf	nf	nf	nf	nf	nf
C0 3	1-Butanol, 3-methyl-	Mild odor	1.7	11.89 ± 0.34b	1149.81 ± 242.71a	1094.43 ± 580.27a	3 ± 8a	760.43 ± 111.4	888.77 ± 20.66a
C0 4	3-Hexen-1-ol, (Z)-	Powerful grassy-green odor	0.013	687.87 ± 157.28	nf	nf	nf	nf	nf
C0 5	Phenylethyl Alcohol	floral odor of roses	100000	nf	0.01 ± 0.01a	0.01 ± 0.00a	0.01 ± 0.00a	0.00a	0.01 ± 0.00a
C0 6	Cyclohexanol, 5-methyl-2-(1-methylethyl)-	Peppermint	nf	nf	nf	nf	nf	nf	nf
C0 7	1-Dodecanol	Typical fatty alcohol sweet	1.2	nf	nf	nf	nf	2.50 ± 3.54	nf





C2 6	Tridecanoic acid, 3-hydroxy-, ethyl ester	nf	nf	nf	nf	nf	nf	nf	nf
C2 7	Nerolidyl acetate	nf	nf	nf	nf	nf	nf	nf	nf
C2 8	Hexadecanoic acid, ethyl ester	nf	nf	nf	nf	nf	nf	nf	nf
C2 9	Succinic acid, hept- 2-yl oct-1-en-3-yl ester	nf	nf	nf	nf	nf	nf	nf	nf
C3 0	Isopropyl palmitate	Almost odorless	nf	nf	nf	nf	nf	nf	nf
	1,2-								
C3 1	Benzenedicarboxyl ic acid, butyl octyl ester	nf	nf	nf	nf	nf	nf	nf	nf
	Benzoic acid, 2- formyl-4,6-								
C3 2	dimethoxy-, 8,8- dimethoxyoct-2-yl ester	nf	nf	nf	nf	nf	nf	nf	nf
C3 3	Decanoic acid, tetradecyl ester Acids	nf	nf	nf	nf	nf	nf	nf	nf
C3 4	Acetic acid	Sharp pungent sour vinegar	13	nf	2.79 ± 0.48ab	5.00 ± 2.57b	5.54 ±	3.40 ± 0.50b	7.21 ± 0.80a



Ketone

	2-Buten-1-one, 1-								
C4 3	(2,6,6-trimethyl- 1,3-cyclohexadien- 1-yl)-, (E)- Phenols	nf	nf	nf	nf	nf	nf	nf	nf
C4 4	3-Allyl-6- methoxyphenol	nf	nf	nf	nf	nf	nf	nf	nf
C4 5	2,4-bis(1,1- dimethylethyl)- phenol	nf	nf	nf	nf	nf	nf	nf	nf
C4 6	2,4-bis(1,1- dimethylethyl)-5- methylphenol	nf	nf	nf	nf	nf	nf	nf	nf
	Benzenes								
C4 7	p-Xylene	Like benzene; characteristic aromatic.	58	nf	nf	0.73 ± 0.23	nf	0.25 ± 0.18	nf
C4 8	o-Xylene	Benzene-like; characteristic aromatic.	380	nf	nf	nf	nf	0.01 ± 0.02	nf
C4 9	Benzene, 1,2,4- trimethyl-	Distinctive, aromatic odor	120	nf	0.26 ± 0.09ab	0.38 ± 0.08ab	0.21 ± 0.08a	0.29 ± 0.06ab	0.19 ± 0.06b

Alkene

C5	1-Octene, 3-(methoxymethoxy)	nf						
0	-							

Results represent the mean ± SD for three independent experiments. Mean values in the same row with the same letters are not significantly different at the 95% confidence level (Duncan's test). Thresholds refer to thresholds in water and air. nf = no threshold was found. Different letters represent significant difference at the 95% percent confidence level (Duncan's test). Sc, *S. cerevisiae* monoculture. Sc+Zb, *S. cerevisiae* and *Z. bailii* co-inoculation. Sc+Rm, *S. cerevisiae* and *R. mucilaginosa* co-inoculation. Sc+Km, *S. cerevisiae* and *K. marxianus* co-inoculation. Sc+Dh, *S. cerevisiae* and *D. hansenii* co-inoculation. FAJ, apple juice.

**Supplementary Table S3** GC-IMS global area set integration parameters obtained from apple juice and cider samples fermented with different non-*Saccharomyces* yeast

Count	Compound	CAS#	Formula	MW	RI	Rt [sec]	Dt [a.u.]	Comment
1	Acetic acid	C64197	C <sub>2</sub> H <sub>4</sub> O <sub>2</sub>	60.1	1496.3	1182.294	1.15026	
2	Octanoic acid ethyl ester	C106321	C <sub>10</sub> H <sub>20</sub> O <sub>2</sub>	172.3	1450.0	1078.527	1.48342	Monomer
3	Octanoic acid ethyl ester	C106321	C <sub>10</sub> H <sub>20</sub> O <sub>2</sub>	172.3	1448.7	1075.747	2.03352	Dimer
4	Ethyl lactate	C97643	C <sub>5</sub> H <sub>10</sub> O <sub>3</sub>	118.1	1351.0	886.122	1.54168	
5	Ethyl hexanoate	C123660	C <sub>8</sub> H <sub>16</sub> O <sub>2</sub>	144.2	1235.0	706.102	1.80106	
6	3-Methyl-1-butanol	C123513	C <sub>5</sub> H <sub>12</sub> O	88.1	1210.7	673.855	1.50511	
7	Isoamyl acetate	C123922	C <sub>7</sub> H <sub>14</sub> O <sub>2</sub>	130.2	1123.5	521.833	1.75012	
8	2-Methyl-1-propanol	C78831	C <sub>4</sub> H <sub>10</sub> O	74.1	1097.9	480.847	1.37251	
9	Ethyl 3-methylbutanoate	C108645	C <sub>7</sub> H <sub>14</sub> O <sub>2</sub>	130.2	1065.5	440.546	1.26456	Monomer
10	Ethyl 3-methylbutanoate	C108645	C <sub>7</sub> H <sub>14</sub> O <sub>2</sub>	130.2	1064.3	439.125	1.65941	Dimer
11	Ethyl 2-methylbutanoate	C7452791	C <sub>7</sub> H <sub>14</sub> O <sub>2</sub>	130.2	1048.0	420.413	1.656	
12	Ethyl butanoate	C105544	C <sub>6</sub> H <sub>12</sub> O <sub>2</sub>	116.2	1036.7	407.86	1.56134	
13	2-Methylpropyl acetate	C110190	C <sub>6</sub> H <sub>12</sub> O <sub>2</sub>	116.2	1013.7	383.464	1.61933	
14	2-Pentanone	C107879	C <sub>5</sub> H <sub>10</sub> O	86.1	981.2	354.659	1.3786	
15	Ethyl acetate	C141786	C <sub>4</sub> H <sub>8</sub> O <sub>2</sub>	88.1	872.9	283.419	1.34468	
16	Ethanol	C64175	C <sub>2</sub> H <sub>6</sub> O	46.1	918.0	311.168	1.1329	
17	Acetone	C67641	C <sub>3</sub> H <sub>6</sub> O	58.1	814.5	251.17	1.11831	
18	Furfural	C98011	C <sub>5</sub> H <sub>4</sub> O <sub>2</sub>	96.1	1485.2	1156.53	1.08772	Monomer
19	Furfural	C98011	C <sub>5</sub> H <sub>4</sub> O <sub>2</sub>	96.1	1488.7	1164.774	1.34056	Dimer
20	(Z)-3-hexenyl acetate	C3681718	C <sub>8</sub> H <sub>14</sub> O <sub>2</sub>	142.2	1323.6	839.039	1.82042	
21	Ethyl propanoate	C105373	C <sub>5</sub> H <sub>10</sub> O <sub>2</sub>	102.1	959.0	338.739	1.45712	
22	Hexanal	C66251	C <sub>6</sub> H <sub>12</sub> O	100.2	1087.9	467.8	1.56455	

23	Butan-1-ol	C71363	C <sub>4</sub> H <sub>10</sub> O	74.1	1148.7	565.785	1.38459	
24	Methyl hexanoate	C106707	C <sub>7</sub> H <sub>14</sub> O <sub>2</sub>	130.2	1183.6	632.58	1.68889	
25	(Z)-3-hexenol	C928961	C <sub>6</sub> H <sub>12</sub> O	100.2	1395.2	967.324	1.22876	Monomer
26	(Z)-3-hexenol	C928961	C <sub>6</sub> H <sub>12</sub> O	100.2	1396.5	969.772	1.51253	Dimer
27	1-Hexanol	C111273	C <sub>6</sub> H <sub>14</sub> O	102.2	1362.7	906.929	1.64043	Monomer
28	1-Hexanol	C111273	C <sub>6</sub> H <sub>14</sub> O	102.2	1361.8	905.297	1.99814	Dimer
29	3-Pentanone	C96220	C <sub>5</sub> H <sub>10</sub> O	86.1	978.6	352.752	1.35687	

**Supplementary Table S4** Physicochemical parameters of apple juice and cider samples fermented with different non-*Saccharomyces* yeast

Samples	FAJ	Ciders				
		Sc	Sc+Zb	Sc+Rm	Sc+Km	Sc+Dh
Tartaric acid (g/L)	1.40 ± 0.02a	0.52 ± 0.02b	0.52 ± 0.04b	0.33 ± 0.04c	0.34 ± 0.02c	0.28 ± 0.11c
Malic acid (g/L)	1.14 ± 0.02a	0.68 ± 0.01b	0.46 ± 0.06c	0.11 ± 0.00d	0.17 ± 0.02e	0.18 ± 0.01e
Acetic acid (g/L)	2.18 ± 0.04a	2.32 ± 0.05a	2.66 ± 0.05ab	3.04 ± 0.13b	3.03 ± 0.08b	2.44 ± 0.56a
Citric acid (g/L)	0.38 ± 0.00a	0.46 ± 0.03b	ND	0.49 ± 0.02b	ND	0.43 ± 0.06ab
Succinic acid (g/L)	ND	0.38 ± 0.13a	0.27 ± 0.09a	0.17 ± 0.02ab	0.23 ± 0.05ab	0.21 ± 0.01ab

Results represent the mean ± SD for three independent experiments. Different letters represent significantly difference at the 95% confidence level (Duncan's test). ND = not detected. TA = total acid (expressed as percentage of tartaric acid). Sc, *S. cerevisiae* monoculture. Sc+Zb, *S. cerevisiae* and *Z. bailii* co-inoculation. Sc+Rm, *S. cerevisiae* and *R. mucilaginosa* co-inoculation. Sc+Km, *S. cerevisiae* and *K. marxianus* co-inoculation. Sc+Dh, *S. cerevisiae* and *D. hansenii* co-inoculation. FAJ, apple juice.