

Supplementary table 1. Correlation matrix among instrumental flavor analysis, antioxidant acti

Variables	Carbon disulfide	Isobutyraldehyde	Benzene	3-methyl butanal
Carbon disulfide	<b>1</b>	0.016	<b>0.500</b>	<b>0.586</b>
Isobutyraldehyde		<b>1</b>	<b>-0.508</b>	<b>0.590</b>
Benzene			<b>1</b>	<b>0.228</b>
3-methyl butanal				<b>1</b>
Thiophene				
2-methyl butanal				
2-methyl-2-butenal				
2-pentanol				
Pentanal				
Dimethyl disulfide				
3-methyl 1-butanol				
2-Methyl pentanal				
2-methyl 1-butanol				
Pyridine				
Methyl benzene				
Isoamyl phenyl acetate				
2-Methylthiophene				
1-Pentanol				
Hexanal				
2-methyl pyrazine				
1,3,5,7-Cyclooctatetraene				
2-Heptanone				
2,5-dimethyl-Pyrazine				
2-Pentyl furan				
Benzaldehyde				
2,3,5-trimethyl pyrazine				
DPPH				
Total phenol (Free)				
Total phenol (Bound)				
Polyphenol				
TEAC (Free)				
TEAC (Bound)				
Total glucan				
α- glucan				
β- glucan				

α=0.05

bold text=significant

ivities, and functional activity (glucan contents) results

Thiophene	2-methyl butanal	2-methyl-2- butenal	2-pentanol	Pentanal	Dimethyl disulfide
<b>0.035</b>	<b>-0.302</b>	<b>0.663</b>	<b>-0.910</b>	<b>-0.575</b>	<b>-0.575</b>
<b>0.361</b>	<b>0.343</b>	<b>0.382</b>	<b>0.054</b>	<b>-0.131</b>	<b>-0.131</b>
<b>0.076</b>	<b>-0.630</b>	<b>-0.279</b>	<b>-0.736</b>	<b>-0.620</b>	<b>-0.620</b>
<b>0.553</b>	<b>-0.055</b>	<b>0.423</b>	<b>-0.620</b>	<b>-0.789</b>	<b>-0.789</b>
<b>1</b>	<b>-0.534</b>	<b>-0.210</b>	<b>-0.292</b>	<b>-0.160</b>	<b>-0.160</b>
	<b>1</b>	<b>0.378</b>	<b>0.606</b>	0.015	0.015
		<b>1</b>	<b>-0.316</b>	<b>-0.200</b>	<b>-0.200</b>
			<b>1</b>	<b>0.641</b>	<b>0.641</b>
				<b>1</b>	<b>1.000</b>
					<b>1</b>

3-methyl 1-butanol	2-Methyl pentanal	2-methyl 1-butanol	Pyridine	Methyl benzene	Isoamyl phenyl acetate
0.607	-0.867	0.701	-0.575	-0.615	0.252
-0.584	-0.011	-0.547	-0.131	0.085	-0.923
0.824	-0.760	0.766	-0.620	-0.291	0.657
-0.043	-0.742	0.016	-0.789	-0.620	-0.246
-0.427	-0.269	-0.450	-0.160	-0.410	-0.210
-0.380	0.438	-0.314	0.015	0.187	-0.365
0.041	-0.301	0.200	-0.200	-0.430	-0.200
-0.643	0.957	-0.678	0.641	0.578	-0.316
-0.407	0.837	-0.429	1.000	0.546	-0.200
-0.407	0.837	-0.429	1.000	0.546	-0.200
1	-0.613	0.985	-0.407	-0.097	0.645
	1	-0.646	0.837	0.620	-0.301
		1	-0.429	-0.201	0.642
			1	0.546	-0.200
				1	-0.430
					1

2-Methyl thiophene	1-Pentanol	Hexanal	2-methyl pyrazine	1,3,5,7- Cyclooctatetraen e	2-Heptanone
-0.575	-0.575	-0.317	-0.575	0.049	-0.575
-0.131	-0.131	-0.787	-0.131	-0.696	-0.131
-0.620	-0.620	-0.093	-0.620	0.535	-0.620
-0.789	-0.789	-0.730	-0.789	-0.321	-0.789
-0.160	-0.160	-0.509	-0.160	-0.693	-0.160
0.015	0.015	0.178	0.015	0.163	0.015
-0.200	-0.200	-0.162	-0.200	-0.194	-0.200
0.641	0.641	0.437	0.641	-0.027	0.641
1.000	1.000	0.423	1.000	-0.320	1.000
1.000	1.000	0.423	1.000	-0.320	1.000
-0.407	-0.407	0.106	-0.407	0.695	-0.407
0.837	0.837	0.472	0.837	-0.141	0.837
-0.429	-0.429	0.122	-0.429	0.665	-0.429
1.000	1.000	0.423	1.000	-0.320	1.000
0.546	0.546	-0.023	0.546	0.059	0.546
-0.200	-0.200	0.657	-0.200	0.691	-0.200
1	1.000	0.423	1.000	-0.320	1.000
	1	0.423	1.000	-0.320	1.000
		1	0.423	0.527	0.423
			1	-0.320	1.000
				1	-0.320
					1

2,5-dimethyl- Pyrazine	2-Pentyl furan	Benzaldehyde	2,3,5-trimethyl pyrazine	DPPH	Total phenol (Free)
-0.575	-0.575	-0.575	-0.575	-0.393	-0.842
-0.131	-0.131	-0.131	-0.131	0.835	0.242
-0.620	-0.620	-0.620	-0.620	-0.616	-0.841
-0.789	-0.789	-0.789	-0.789	0.097	-0.392
-0.160	-0.160	-0.160	-0.160	0.133	-0.073
0.015	0.015	0.015	0.015	0.348	0.636
-0.200	-0.200	-0.200	-0.200	0.021	-0.177
0.641	0.641	0.641	0.641	0.414	0.949
1.000	1.000	1.000	1.000	0.268	0.560
1.000	1.000	1.000	1.000	0.268	0.560
-0.407	-0.407	-0.407	-0.407	-0.604	-0.833
0.837	0.837	0.837	0.837	0.397	0.889
-0.429	-0.429	-0.429	-0.429	-0.634	-0.838
1.000	1.000	1.000	1.000	0.268	0.560
0.546	0.546	0.546	0.546	0.604	0.365
-0.200	-0.200	-0.200	-0.200	-0.976	-0.423
1.000	1.000	1.000	1.000	0.268	0.560
1.000	1.000	1.000	1.000	0.268	0.560
0.423	0.423	0.423	0.423	-0.604	0.319
1.000	1.000	1.000	1.000	0.268	0.560
-0.320	-0.320	-0.320	-0.320	-0.589	-0.242
1.000	1.000	1.000	1.000	0.268	0.560
1	1.000	1.000	1.000	0.268	0.560
	1	1.000	1.000	0.268	0.560
		1	1.000	0.268	0.560
			1	0.268	0.560
				1	0.462
					1

Total phenol (Bound)	Polyphenol	TEAC (Free)	TEAC (Bound)	Total glucan	α- glucan
-0.543	-0.770	-0.771	0.157	0.152	0.557
-0.670	0.350	0.135	0.744	0.273	0.156
-0.110	-0.911	-0.645	-0.109	0.367	0.680
-0.976	-0.322	-0.203	0.852	0.702	0.445
-0.666	-0.034	0.243	0.399	0.914	0.053
0.124	0.645	0.415	0.377	-0.502	-0.407
-0.427	-0.057	-0.291	0.303	-0.259	0.011
0.585	0.908	0.778	-0.143	-0.446	-0.713
0.640	0.574	0.377	-0.667	-0.512	-0.582
0.640	0.574	0.377	-0.667	-0.512	-0.582
0.171	-0.880	-0.850	-0.369	-0.186	0.636
0.660	0.865	0.698	-0.355	-0.513	-0.729
0.111	-0.867	-0.856	-0.326	-0.222	0.586
0.640	0.574	0.377	-0.667	-0.512	-0.582
0.588	0.333	0.029	-0.353	-0.497	0.081
0.356	-0.517	-0.209	-0.480	-0.042	-0.055
0.640	0.574	0.377	-0.667	-0.512	-0.582
0.640	0.574	0.377	-0.667	-0.512	-0.582
0.777	0.238	0.334	-0.636	-0.549	-0.681
0.640	0.574	0.377	-0.667	-0.512	-0.582
0.517	-0.359	-0.275	-0.278	-0.391	0.150
0.640	0.574	0.377	-0.667	-0.512	-0.582
0.640	0.574	0.377	-0.667	-0.512	-0.582
0.640	0.574	0.377	-0.667	-0.512	-0.582
0.640	0.574	0.377	-0.667	-0.512	-0.582
0.640	0.574	0.377	-0.667	-0.512	-0.582
-0.197	0.530	0.217	0.375	-0.019	0.079
0.332	0.988	0.900	0.078	-0.257	-0.795
1	0.242	0.147	-0.811	-0.730	-0.382
	1	0.872	0.124	-0.253	-0.784
		1	0.179	0.099	-0.838
			1	0.548	0.187
				1	0.271
					1

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$\beta$ - glucan

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0.132  
0.269  
0.346  
0.690  
0.922  
-0.495  
-0.264  
-0.424  
-0.492  
-0.492  
-0.211  
-0.490  
-0.246  
-0.492  
-0.502  
-0.040  
-0.492  
-0.492  
-0.529  
-0.492  
-0.401  
-0.492  
-0.492  
-0.492  
-0.492  
-0.492  
-0.021  
-0.231  
-0.721  
-0.227  
0.130  
0.543  
0.999  
0.238