

Supplementary Data

Variations in physicochemical characteristics, antioxidant activity, phenolic and volatile profiles, and sensory attributes of tea-flavored Chardonnay wine during bottle aging

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Table S1 HPLC-DAD-ESI-QTOF-MS/MS identification of phenolic compounds from tea-macerated Chardonnay wines, tea-macerated model wines and tea ethanolic extracts

Class/subclass	Compound name	Peak No.	Compound No.	Molecular formula	Retention time (min)	UV λ_{\max} (nm)	Theoretical mass (m/z)	Experimental mass (m/z)	Error (ppm)	Major fragment ion MS/MS (m/z, abundance %)	Wine type ^a	ID ^b	Literature ^c
<i>Negative mode [M-H]⁻</i>													
Phenolic acids													
Hydroxybenzoic acid	Quinic acid 1	1	1	C ₇ H ₁₂ O ₆	6.171	208, 234, 282	191.0561	191.0560	-0.52	85 (100), 87 (23), 127 (22)	BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6	D, L	(2)
	Quinic acid 2	16	35	C ₇ H ₁₂ O ₆	13.807	194, 282	191.0561	191.0550	-5.76	191 (100), 93 (14)	BC-9	D, L	(2)
	Gallic acid	17	37	C ₇ H ₆ O ₅	14.146	218, 274	169.0143	169.0156	7.69	125 (100)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	S, D, L	(6), (10), (14), (17)
	Galloylquinic acid	18	39	C ₁₄ H ₁₆ O ₁₀	14.341	216, 276	343.0671	343.0684	3.79	191 (100), 192 (56), 193 (11)	GE, BE, GC-1, GM-1, GC-3, BC-3, GM-3, BM-3, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9,	D, L	(6), (17)

	3,4-Dihydroxybenzoic acid	30	66	C ₇ H ₆ O ₄	20.024	210, 268, 293	153.0193	153.0205	7.91	108 (100), 110 (69), 87 (15)	GM-9, BM-9, GE, BE, C-1, BC-1, C-3, BC-3, C-6, BC-6, C-9, GC-9, BC-9	S, D, L	(14)
	2,6-Bis- <i>O</i> -(3,4,5-trihydroxybenzoyl)-D-glucopyranose 1	34	76	C ₂₀ H ₂₀ O ₁₄	21.375	196, 226, 281, 328sh	483.0780	483.0818	7.87	97 (100), 387 (66), 438 (58)	GC-1, BC-1, GC-3, BM-3, GC-6, GC-9, GM-9, BM-9	D	
	2,6-Bis- <i>O</i> -(3,4,5-trihydroxybenzoyl)-D-glucopyranose 2	35	81	C ₂₀ H ₂₀ O ₁₄	21.850	220, 282	483.0780	483.0745	-7.25	305 (100)	GC-9, BM-9	D	
	3- <i>O</i> -Methylgallic acid	43	106	C ₈ H ₈ O ₅	24.428	194, 238, 272	183.0299	183.0290	-4.90	183 (100)	GC-9	D	
	Syringic acid	64	152	C ₉ H ₁₀ O ₅	34.439	218, 275	197.0455	197.0470	7.61	124 (100), 125 (32), 79 (6)	C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	S, D	
	Gallic acid 1	85	196	C ₇ H ₆ O ₅	44.887	194, 270	169.0143	169.0161	10.65	83 (100), 107 (72)	BC-1, BC-3, GC-6, GC-9	D, L	(6), (10), (14), (17)
Hydroxycinnamic acid	Tartaric acid 1	1	2	C ₄ H ₆ O ₆	6.171	208, 234, 282	149.0092	149.0089	-1.48	74 (100), 88 (40), 85 (21)	C-1, GM-1, BM-1, C-3, GM-3, BM-3, C-6, GM-6, BM-6	D	
	Caffeic acid 3-sulfate	24	52	C ₉ H ₈ O ₇ S	16.471	206, 275	258.9918	258.9880	-14.66	136 (100), 134 (64)	C-9, BC-9	D	
	Glutathionyl caftaric acid	27	62	C ₂₃ H ₂₇ N ₃ O ₁₅ S	18.638	212, 272	616.1090	616.1133	6.98	273 (100), 211 (41), 467 (31)	C-1, GC-1, BC-1, C-3, GC-3, BC-3, C-6, GC-6, BC-6, C-9, GC-9	D, L	(14), (17)

Chlorogenic acid 1	31	68	C ₁₆ H ₁₈ O ₉	20.153	327	353.0878	353.0894	4.53	192 (100), 180 (70), 135 (57)	9, BC-9 GE, GC-1, GM-1, GC-3, GM-3, BM-3, GC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9	D, L	(6), (10)
<i>p</i> -Coumaroylquinic acid 1	41	100	C ₁₆ H ₁₈ O ₈	23.825	312	337.0929	337.0933	1.19	163 (100), 93 (8)	GE, BE, GC-1, BC-1, BM-1, GC-3, BC-3, BM-3, GC-6, BC-6, BM-6, BC-9, GM-9, BM-9	D, L	(6), (9)
Tartaric acid 2	42	103	C ₄ H ₆ O ₆	24.075	194, 210, 298	149.0092	149.0110	12.54	73 (100)	C-1, BC-1, C-3, BC-3, C-6, C-9, BC-9	D	
Chlorogenic acid	44	108	C ₁₆ H ₁₈ O ₉	24.894	242, 327	353.0878	353.0881	0.85	173 (100)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	S, D, L	(6), (10)
Subaphyllin	47	118	C ₁₄ H ₂₀ N ₂ O ₃	25.832	216, 246, 344	263.1401	263.1434	12.59	97 (100), 92 (37), 190 (34)	C-1, GM-3	D	
<i>p</i> -Coumaroylquinic acid 2	53	129	C ₁₆ H ₁₈ O ₈	27.791	211, 278	337.0929	337.0948	5.64	173 (100), 155 (14), 164 (11)	GE, BE, GC-1, BC-1, BM-1, GC-3, BC-3,	D, L	(6), (9)

	Caffeic acid	56	133	C ₉ H ₈ O ₄	29.502	208, 290, 326	179.0350	179.0368	10.05	134 (100), 136 (95), 107 (22)	BM-3, BC-6, BM-6, BC-9, 3BM-9 GE, BE, C-1, C-3, C-6, C-9, BC-9	S, D, L	(14)
	<i>p</i> -Coumaroylquinic acid 3	56	134	C ₁₆ H ₁₈ O ₈	29.502	216, 276	337.0929	337.0953	7.12	173 (100), 174 (50)	GE, BE, GC-1, BC-1, BM-1, GC-3, BC-3, BM-3, GC-6, BC-6, BM-6, GC-9, BC-9, GM-9, BM-9	D, L	(6), (9)
	<i>p</i> -Coumaroylquinic acid 4	59	141	C ₁₆ H ₁₈ O ₈	31.447	208, 238, 298	337.0929	337.0967	11.33	191 (100), 173 (27)	BE, BC-1, BC-3, BM-3, GC-6, BC-6, BC-9, 3BM-9	D, L	(6), (9)
	Hydrocaffeic acid	96	216	C ₉ H ₁₀ O ₄	49.100	195, 226, 264, 300	181.0506	181.0519	6.73	109 (100)	C-1, C-6, C-9, BC-9	D	
	3,4-Dimethoxycinnamic acid	120	259	C ₁₁ H ₁₂ O ₄	60.747	224, 327	207.0663	207.0665	0.97	135 (100), 134 (64), 179 (53), 136 (18)	C-1, GC-1, BC-1, C-3, GC-3, BC-3, C-6, GC-6, BC-6, C-9, GC-9, BC-9	D, L	(14)
Flavonoids													
	3-Deoxyanthocyanidin	128	271	C ₁₇ H ₂₆ O ₃	66.719	200, 222, 254, 486	277.1809	277.1797	-4.32	277 (100)	GE, BE	D	
	Flavan-3-ol Theasinensin C 1	24	51	C ₃₀ H ₂₆ O ₁₄	16.471	206, 275	609.1250	609.1273	3.78	423 (100), 484 (86), 187 (48), 177 (38)	GM-1, GM-3, GM-6, GM-9	D, L	(4)
	Theasinensin C 2	25	54	C ₃₀ H ₂₆ O ₁₄	17.180	206, 275	609.1250	609.1271	3.45	406 (100)	GM-1, GM-3	D, L	(4)
	Gallocatechin	28	63	C ₁₅ H ₁₄ O ₇	19.054	200, 216, 273	305.0667	305.0716	16.06	126 (100), 165 (78), 139 (67), 219 (49), 177 (36)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3,	S, D, L	(2), (6), (17)

Theasinensin C 3	32	70	C ₃₀ H ₂₆ O ₁₄	20.509	206, 276	609.1250	609.1258	1.31	424 (100), 305 (72), 441 (37), 125 (32), 592 (10)	BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9, GC-1, GM-1, GC-3, GM-3, GC-6, GM-6	D, L	(4)
Theacitrin A	34	77	C ₃₇ H ₂₈ O ₁₈	21.375	206, 274	759.1203	759.1171	-4.16	759 (100), 742 (52), 607 (36), 425 (16)	BE	D, L	(4)
Theasinensin B 1	37	84	C ₃₇ H ₃₀ O ₁₈	22.193	230, 298	761.1359	761.1305	-7.09	591 (100), 594 (24), 467 (19), 753 (14)	BE, BC-1, GM-1, BM-1, BC-3, GM-3, BM-3, BC-6, GM-6, BM-6, GM-9, BM-9	D, L	(4)
Epigallocatechin	40	97	C ₁₅ H ₁₄ O ₇	23.385	196, 236, 272	305.0667	305.0657	-3.28	125 (100), 179 (28), 167 (23), 221 (11)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	S, D, L	(2), (17)
Unknown (epi)gallocatechin conjugate	40	98	C ₃₀ H ₂₈ O ₁₄	23.385	196, 236, 272	611.1406	611.1393	-2.13	306 (100), 305 (95), 307 (18)	GE, GC-1, GM-1, GC-3, GM-3, GC-6, BC-6, GM-6, GC-9, GM-9	D	

Theasinensin B 2	43	104	C ₃₇ H ₃₀ O ₁₈	24.428	212, 276	761.1359	761.1411	6.83	305 (100), 169 (83), 161 (47), 593 (40)	GE, BE, GC-1, GM-1, GC-3, GM-3, GC-6, GM-6, GC-9, GM-9	D, L	(4)
Catechin	46	112	C ₁₅ H ₁₄ O ₆	25.502	206, 224sh, 285	289.0718	289.0664	-18.68	289 (100), 245 (65), 232 (50), 165 (42), 137 (38), 93 (37)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	S, D, L	(4), (6), (14)
(Epi)gallocatechin-(epi)catechin gallate 1	51	127	C ₃₇ H ₃₀ O ₁₇	27.305	208, 281	745.1410	745.1456	6.17	424 (100), 560 (44), 407 (38), 593 (29), 575 (25)	BE, GC-1, GM-1, GC-3, GM-3, GC-6, GM-6, GC-9, GM-9	D, L	(9)
Theasinensin A	52	128	C ₄₄ H ₃₄ O ₂₂	27.498	210, 282	913.1469	913.1527	6.35	744 (100), 425 (65), 733 (54), 763 (46)	GE, BE, GC-1, BC-1, BM-1, GC-3, BC-3, BM-3, GM-9	D, L	(4)
(Epi)gallocatechin-(epi)catechin gallate 2	54	131	C ₃₇ H ₃₀ O ₁₇	28.239	204, 280	745.1410	745.1453	5.77	576 (100), 728 (42), 550 (28), 348 (20), 273 (17)	GC-1, GM-1, GC-3, GM-3, GC-6, GM-6, GC-9, GM-9	D, L	(9)
Epicatechin	56	135	C ₁₅ H ₁₄ O ₆	29.502	216, 276	289.0720	289.0747	9.34	109 (100), 123 (82), 97 (28), 125 (29)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3,	S, D, L	(4), (6), (14)

Epigallocatechin gallate	58	137	C ₂₂ H ₁₈ O ₁₁	30.555	196, 232, 276	457.0776	457.0795	4.16	169 (100), 170 (45)	BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	S, D, L	(4), (6)
(Epi)gallocatechin-cyanidin-3,5-diglucoside	58	138	C ₄₂ H ₄₄ O ₂₃	30.555	196, 234, 277	915.2201	915.1880	-35.04	169 (100), 170 (32), 458 (18), 459 (15), 457 (12), 305 (12)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, GM-9	D, L	(10)
Gallocatechin gallate dimer	58	140	C ₄₄ H ₃₆ O ₂₂	30.555	196, 232, 276	915.1625	915.1680	6.01	170 (100), 458 (79), 459 (63), 171 (49)	BE, GC-9, GM-9	D, L	(7)
Gallocatechin gallate	63	149	C ₂₂ H ₁₈ O ₁₁	33.413	210, 275, 334	457.0776	457.0808	7.00	170 (100)	GC-1, GM-1, GC-3, GM-3, GC-6, GM-6	S, D	
Theasinensin B 3	65	155	C ₃₇ H ₃₀ O ₁₈	35.315	208, 281	761.1359	761.1315	-5.78	761 (100), 591 (39), 743 (28)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GM-9	D, L	(4)
Epigallocatechin 3,3'-di-gallate	74	171	C ₂₉ H ₂₂ O ₁₅	38.451	228, 310	609.0886	609.0902	2.63	457 (100), 306 (71), 573 (39)	GM-9	D, L	(9)
(Epi)catechin gallate dimer	76	174	C ₄₄ H ₃₆ O ₂₀	39.622	196, 232, 280	883.1727	883.1759	3.62	442 (100), 441 (81), 443 (77), 289 (9)	GE, GC-1, GC-3	D, L	(20)

Epicatechin gallate	76	176	C ₂₂ H ₁₈ O ₁₀	39.622	196, 232, 280	441.0827	441.0844	3.85	169 (100), 290 (61), 170 (51)	GC-3, BC-3, GM-3, GC-6, GM-6, GC-9, GM-9, GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, GM-9, BM-9	S, D, L	(2), (6), (9)
Epicatechin gallate 1	79	182	C ₂₂ H ₁₈ O ₁₀	41.522	208, 276	441.0827	441.0844	3.85	290 (100), 246 (14), 272 (13)	GE, GC-1, GM-1, GC-3, GM-3, GC-6, GM-6, GC-9	D, L	(2), (6), (9)
Theacitrin C	89	203	C ₄₄ H ₃₂ O ₂₂	46.401	194, 277	911.1312	911.1406	10.22	875 (100), 396 (95), 850 (75)	BE	D, L	(13)
Methyl epicatechin gallate	90	204	C ₂₃ H ₂₀ O ₁₀	46.533	194, 280	455.0984	455.1007	5.05	79 (100), 181 (50), 283 (31)	GE, BE, GM-1, GC-3, GM-3	D, L	(2)
Epiafzelechin 3-gallate	92	209	C ₂₂ H ₁₈ O ₉	47.594	194, 206sh, 275	425.0878	425.0884	1.41	274 (100), 169 (10), 255 (9)	GE, BE, GC-1, GM-1, GC-3, GM-3, GM-6	D, L	(11)
Epicatechin-3,5-digallate	96	217	C ₂₉ H ₂₂ O ₁₄	49.100	208, 282	593.0937	593.0977	6.74	563 (100), 443 (36), 476 (27), 523 (27)	GE, GM-1, GM-3, GC-9, GM-9, BM-9	D, L	(18)
Epigallocatechin gallate-4',4''-O-β-D-glucopyranoside	98	219	C ₃₄ H ₃₈ O ₂₁	50.154	248, 364	781.1833	781.2199	46.85	607 (100)	GE, GC-1, GM-1, 4GM-9	D, L	(22)
Unknown (epi)catechin glucoside	100	222	C ₂₁ H ₂₄ O ₁₁	50.678	210, 226sh, 316	451.1246	451.1080	-36.70	161 (100), 417 (55), 286 (54)	GE	D, L	(21)
Unknown di(epi)catechin-dimethylglyoxal adduct	102	226	C ₃₆ H ₃₆ O ₁₆	51.259	214, 316	723.1931	723.5140	443.76	678 (100), 679 (17)	BE	D, L	(3)
Epigallocatechin 3-O-gallate-7-O-glucoside-	106	231	C ₂₂ H ₁₈ O ₁₁	54.599	210, 276	795.1625	795.1700	9.43	463 (100), 490 (61), 305 (48), 331 (39)	GC-9	D, L	(15)

	4"- <i>O</i> -glucuronide												
	Theasinensin F	109	235	C ₄₄ H ₃₄ O ₂₁	55.721	228, 380	897.1520	897.1536	1.80	441 (100), 285 (47), 855 (37)	GM-9	D	
	Epiphylloumarin- [4β→8]-(-)-epicatechin- [4β→8]-(-)-epicatechin	113	242	C ₄₈ H ₃₈ O ₁₉	58.117	320	917.1935	917.2600	67.85	887 (100)	BE	D	
	Epitheaflavic acid 3'- gallate	114	244	C ₂₈ H ₂₀ O ₁₄	58.315	208, 282, 398	579.0780	579.0900	22.19	150 (100), 505 (47), 193 (38), 339 (35)	BE	D	
	Theaflavic acid	115	246	C ₂₁ H ₁₆ O ₁₀	58.863	282, 396	427.0671	427.0727	13.16	227 (100), 137 (66)	BE	D	
	Theaflavin	115	247	C ₂₉ H ₂₄ O ₁₂	58.863	210, 268, 298, 365sh	563.1195	563.1227	5.68	546 (100), 523 (91), 138 (81), 231 (77), 486 (53)	GE, BE, BC-1, BC-3	D, L	(6)
	Theaflavin 3-gallate	118	255	C ₃₆ H ₂₈ O ₁₆	59.832	210, 270	715.1305	715.1359	7.55	564 (100)	GE, BE, BC-1, BC-3	D, L	(6)
	Epitheafagallin 3- <i>O</i> - gallate	119	256	C ₂₇ H ₂₀ O ₁₃	60.230	214, 314	551.0831	551.0922	16.52	125 (100), 234 (31), 295 (30), 187 (25)	BE	D, L	(8)
	Theaflavin-3,3'-gallate	119	257	C ₄₃ H ₃₂ O ₂₀	60.230	208, 277, 378	867.1414	867.1475	7.03	715 (100)	GE, BE, BC-1, BC-3	D, L	(6)
Flavanonol	Katuranin 2	102	225	C ₁₅ H ₁₂ O ₆	51.259	204, 268, 316	287.0561	287.0549	-4.15	125 (100), 260 (17)	GE, BE, C-9	D	
Flavone	Apigenin 6,8-di- <i>C</i> - glucoside	55	132	C ₂₇ H ₃₀ O ₁₅	29.077	272, 352	593.1512	593.1541	4.89	391 (100)	BE, BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, BC-6, GM-6, BM-6, GC-9	D, L	(4)
Flavonoid- <i>O</i> - glucuronide	6-[6-((6-[(6-Aminopyridin-3-yl)methyl]-2-(3,4-dihydroxyphenyl)-5-hydroxy-4-oxo-4 <i>H</i> -chromen-7-yl}oxy)-3,4,5-trihydroxyoxane-2-carboxyloxy]-3,4,5-trihydroxyoxane-2-carboxylic acid	104	229	C ₃₃ H ₃₂ N ₂ O ₁₈	52.761	195, 280, 330	743.1577	743.1546	-4.22	439 (100), 302 (94), 147 (83), 573 (47)	GC-9	D	
Flavonoid- <i>O</i> -glycoside	Tilioside 1	29	65	C ₃₀ H ₂₆ O ₁₃	19.574	204, 274	593.1301	593.1334	5.56	426 (100), 289 (34), 409 (30), 304 (17)	GC-1, GM-1, GC-3, GM-3, GM-6, GC-9, GM-9	D	
	Tilioside 2	33	74	C ₃₀ H ₂₆ O ₁₃	20.888	200, 268sh, 334	593.1301	593.1351	8.43	426 (100), 177 (17), 468 (23), 380 (10)	GC-1, GC-9, GM-9	D	
	Tilioside 3	39	94	C ₃₀ H ₂₆ O ₁₃	22.905	216, 260, 294sh	593.1301	593.1345	7.42	424 (100), 287 (11), 177 (10)	GC-1, GM-1	D	
	Miscanthoside	81	190	C ₂₁ H ₂₂ O ₁₁	42.668	204, 230sh, 295	449.1089	449.1124	7.65	286 (100), 355 (50), 381 (24), 324 (22), 284 (19)	C-1, C-6, C-9	D	

	Unknown flavonoid-7- <i>O</i> -glycoside 1 (myriophylloside B or myriophylloside type)	114	243	C ₄₁ H ₄₄ O ₂₂	58.315	210, 285	887.2251	887.2325	8.34	840 (100)	BC-1, GM-1, GC-3, GM-3, BM-3, GC-6, GM-6, BM-6, GC-9, BC-9	D	
	Unknown flavonoid-7- <i>O</i> -glycoside 2 (myriophylloside D or myriophylloside type)	118	254	C ₄₁ H ₄₄ O ₂₁	59.832	266, 318	871.2302	871.2324	2.53	828 (100), 585 (78)	BC-1, GM-1, GM-3, BM-3, BM-6, GC-9, BC-9, GM-9, BM-9	D	
	Tiliroside 4	120	258	C ₃₀ H ₂₆ O ₁₃	60.747	206, 268, 372	593.1301	593.1290	-1.85	563 (100), 483 (33), 413 (66), 284 (22), 531 (18)	GC-1, GM-1, GC-3, GM-3, BM-3, BM-6, GM-9, BM-9	D	
Flavonol	Quercetin 3- <i>O</i> -di-galactoside/glucoside	65	154	C ₂₇ H ₃₀ O ₁₇	35.315	208, 281	625.1410	625.1448	6.08	317 (100)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, GM-9	D, L	(5)
	Myricetin 3- <i>O</i> -glucoside	67	158	C ₂₁ H ₂₀ O ₁₃	36.197	254, 343	479.0831	479.0810	-4.43	316 (100), 271 (14)	BE	D, L	(10)
	Myricetin-3- <i>O</i> -galactoside 1	68	159	C ₂₁ H ₂₀ O ₁₃	36.379	212, 270, 350	479.0831	479.0858	5.64	318 (100), 328 (19), 152 (17)	GE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6,	D, L	(10)

Myricetin-3- <i>O</i> -galactoside 2	69	160	C ₂₁ H ₂₀ O ₁₃	36.799	210, 264, 360	479.0831	479.0862	6.47	318 (100), 289 (13)	GC-9, GM-9, GE, BE, GC-1, BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, GC-6, GM-6, GC-9, BC-9, GM-9, GE, BE, BC-1, BM-1, GC-3, BC-3, GM-3, BM-3, BC-6, BM-6, BC-9, BM-9	D, L	(10)
Quercetin 3- <i>O</i> -glucosylrutinoside 1	70	161	C ₃₃ H ₄₀ O ₂₁	37.066	206, 268, 350	771.1989	771.2065	9.85	675 (100)	GC-9, GE, BE, BC-1, BM-1, GC-3, BC-3, GM-3, BM-3, BC-6, BM-6, BC-9, BM-9	D, L	(4), (9)
Kaempferol 3- <i>O</i> -rutinoside 1	72	164	C ₂₇ H ₃₀ O ₁₅	37.274	216, 271, 347	593.1512	593.1553	6.91	440 (100), 294 (92)	GC-1, BC-1, GM-1, BM-1, BC-3, BM-3, GC-6, GM-6, GC-9, GM-9, BM-9	D, L	(6)
Quercetin 3- <i>O</i> -glucosylrutinoside 2	73	167	C ₃₃ H ₄₀ O ₂₁	38.075	208, 260, 360	771.1989	771.2077	11.41	302 (100), 730 (52)	GC-1, BC-1, GM-1, BM-1, BC-3, BM-3, GC-6, GM-6, GC-9, BC-9, GM-9, BM-9	D, L	(4), (9)
Kaempferol 3- <i>O</i> -glucosylrhamnosyl-galactoside 1	77	177	C ₃₃ H ₄₀ O ₂₀	40.518	200, 264, 350	755.2040	755.2013	-3.61	285 (100), 755 (69), 286 (18), 284 (15)	BE	D	
Rutin	77	178	C ₂₇ H ₃₀ O ₁₆	40.518	204, 258, 357	609.1461	609.1490	4.76	301 (100), 458 (27), 179 (10)	GE, BE, C-1, GC-	S, D, L	(2), (4)

Quercetin-3- <i>O</i> -glucoside 2	80	186	C ₂₁ H ₂₀ O ₁₂	42.306	208, 258, 356	463.0882	463.0890	1.73	301 (100)	1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9, GE, BE, GC-1, GM-1, GC-3, GM-3, GC-6, BC-6, GM-6, GC-9, GM-9, BM-9	D, L	(9)
Kaempferol 3- <i>O</i> -glucosylrhamnosyl-galactoside 2	80	187	C ₃₃ H ₄₀ O ₂₀	42.306	200, 268, 350	755.2040	755.2166	16.68	285 (100), 415 (9)	GE, BE, BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, BC-6, BM-6, BC-9, BM-9	D	
Quercetin 3- <i>O</i> -galactoside	82	192	C ₂₁ H ₂₀ O ₁₂	42.886	208, 258, 357	463.0882	463.0892	2.16	179 (100), 389 (77)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9	S, D, L	(6), (17)
Kaempferol 3- <i>O</i> -rutinoside 2	83	193	C ₂₇ H ₃₀ O ₁₅	43.739	206, 254, 342	593.1512	593.1516	0.67	498 (100)	GE, BE, C-1, GC-	D, L	(6)

Kaempferol 3- <i>O</i> -(6- <i>O</i> -rhamnosyl-hexoside)7- <i>O</i> -rhamnoside	84	194	C ₃₃ H ₄₀ O ₁₉	44.325	212, 240, 268, 342	739.2091	739.2080	-1.49	593 (100)	1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GE, BE, GC-1, GM-1, GC-3, GM-3, GC-9, GM-9	D	
Kaempferol 3- <i>O</i> -rutinoside 3	87	198	C ₂₇ H ₃₀ O ₁₅	45.751	198, 268, 349	593.1512	593.1510	-0.34	483 (100), 563 (48), 286 (44)	GE, BE, GC-1, BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, GC-6, GM-6, GC-9, BC-9, GM-9, BM-9	D, L	(6)
Kaempferol 3- <i>O</i> -glucoside 1	89	201	C ₂₁ H ₂₀ O ₁₁	46.401	194, 280	447.0933	447.0955	4.92	285 (100)	GE, BE, BC-1, BC-3, GC-6, BC-6, BM-6	D, L	(9)
Kaempferol 7-(6"-galloyl)glucoside)	89	202	C ₂₈ H ₂₄ O ₁₅	46.401	194, 210sh, 280	599.1042	599.1044	0.26	301 (100)	BE, GC-1, GM-9	D	
Kaempferol 3- <i>O</i> -glucoside	94	212	C ₂₁ H ₂₀ O ₁₁	48.399	198, 266, 350	447.0933	447.0942	2.01	284 (100), 447 (58)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9	S, D, L	(9)

Kaempferol 3- <i>O</i> -glucoside 3	101	223	C ₂₁ H ₂₀ O ₁₁	51.044	198, 266, 350	447.0933	447.0904	-6.49	300 (100), 258 (38), 374 (36), 179 (32)	GC-9, BC-9, GM-9, BM-9	D, L	(9)
Quercetagenin	108	234	C ₁₅ H ₁₀ O ₈	55.294	228, 380	317.0303	317.0279	-7.57	151 (100), 137 (92), 255 (29)	GC-9	D	
Quercetin 1	109	236	C ₁₅ H ₁₀ O ₇	55.721	228, 380	301.0354	301.0351	-1.00	301 (100)	GC-1, GM-1, GM-3, GC-9, GM-9	D, L	(17)
Unknown quercetin conjugate 1	113	241	C ₄₂ H ₄₆ O ₂₂	58.117	210, 270, 318	901.2408	901.2466	6.44	757 (100), 302 (17)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9	D, L	(6)
Kaempferol 3-(2G-xylosylrutinoside)-7-glucoside	114	245	C ₃₈ H ₄₈ O ₂₄	58.315	208, 282, 398	887.2463	887.2400	-4.54	858 (100)	BE, GM-9, BM-9	D	
Unknown kaempferol conjugate	115	248	C ₄₂ H ₄₆ O ₂₁	58.863	212, 270, 315	885.2459	885.2456	-0.34	740 (100), 787 (72), 792 (69)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9	D, L	(6)
Unknown quercetin conjugate 2	116	249	C ₄₂ H ₄₆ O ₂₂	59.059	206, 316	901.2408	901.2262	-16.20	755 (100)	GC-9, GM-9, BM-9	D, L	(6)
Kaempferol 3- <i>O</i> -glucosylrhamnosylgalactoside 3	116	250	C ₃₃ H ₄₀ O ₂₀	59.059	206, 316	755.2040	755.1815	-29.79	289 (100), 745 (59)	GE, GM-1, GC-3, GM-3, GC-9, GM-9	D	
Quercetin	121	261	C ₁₅ H ₁₀ O ₇	61.100	206, 258, 374	301.0354	301.0385	10.30	151 (100), 179 (51), 107 (13)	GE, BE, GC-1,	S, D, L	L data 9, 10

	Kaempferol	124	266	C ₁₅ H ₁₀ O ₆	63.686	228, 370	285.0405	285.0392	-4.56	257 (100), 227 (74), 210 (47), 151 (45)	GM-1, BM-1, GC-3, GM-3, BM-3, GC-6, GM-6, BM-6, GC-9, GM-9, GE, BE, GC-1, GM-1, GC-3, GM-3, GM-6, GC-9, GM-9, BM-9	S, D, L	(17)
Isoflavonoid	3-Oxo-3-[[3,4,5-trihydroxy-6-(3-hydroxy-14-methoxy-8,17-dioxatetracyclo[8.7.0.0 ² ,7.0 ¹¹ ,1 ⁶]heptadeca-2,4,6,11,13,15-hexaen-5-yl]oxy)oxan-2-yl]methoxy}propanoic acid	75	173	C ₂₅ H ₂₆ O ₁₃	39.040	220, 274, 338	533.1301	533.1362	11.44	441 (100), 97 (20)	BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, BC-6, BM-6, GC-9, BC-9, GM-9, BM-9	D	
Neoflavone	Dihydrogenistin	103	228	C ₂₁ H ₂₂ O ₁₀	52.494	195, 280, 330	433.1140	433.1099	-9.51	180 (100), 259 (62), 126 (12)	C-9	D	
Neoflavone	8-Methyl-2-oxo-4-phenyl-2 <i>H</i> -chromen-7-yl 2-(4-methylbenzenesulfonamido)-3-phenylpropanoate	99	220	C ₃₂ H ₂₇ NO ₆ S	50.467	224, 280	552.1486	552.1455	-5.67	401 (100), 382 (60), 236 (24), 508 (10)	GC-9, GM-9	D	
Proanthocyanidin	Procyanidin B3	38	89	C ₃₀ H ₂₆ O ₁₂	22.477	204, 213sh, 282	577.1351	577.1394	7.45	484 (100), 370 (18), 543 (14), 481 (12)	C-1, GC-1, BC-1, GM-1, C-3, GC-3, BC-3, GM-3, GM-6, GC-9, GM-9	D, L	(17)
Proanthocyanidin	Procyanidin B1	49	123	C ₃₀ H ₂₆ O ₁₂	26.435	204, 282	577.1351	577.1390	6.76	290 (100), 426 (34), 560 (19)	GE, BE, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-	S, D, L	(17)

											6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9		
	Procyanidin B2	50	124	C ₃₀ H ₂₆ O ₁₂	27.127	204, 325	577.1351	577.1393	7.28	408 (100), 261 (18), 485 (17), 287 (15)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6	S, D, L	(17)
	Procyanidin B1 3-O-gallate	59	142	C ₃₇ H ₃₀ O ₁₆	31.447	208, 282, 310	729.1461	729.1497	4.94	635 (100), 636 (40)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6	D	
	Procyanidin B2 3-O-gallate	61	147	C ₃₇ H ₃₀ O ₁₆	32.412	204, 280	729.1461	729.1530	9.46	636 (100)	GC-1, BC-1, GM-1, BC-3, GM-3, GC-6, GC-9, BM-9	D	
Pyranoflavonoid	Unknown pyranoflavonoid	90	205	C ₄₅ H ₅₀ N ₄ O ₇	46.533	194, 280	757.3607	757.3591	-2.08	271 (100), 485 (82)	GC-3	D	
Others: phenolic compounds													
Anthraquinone	Unknown anthraquinone	16	36	C ₂₈ H ₃₃ O ₁₈ P	13.807	194, 282	687.1332	687.1348	2.32	343 (100)	GC-9, BM-9	D	
Benzenediol	2-(3,4-Dihydroxyphenyl)-3,5,7-trihydroxy-4a,8a-dihydro-4H-chromen-4-one	79	183	C ₁₅ H ₁₂ O ₇	41.522	205, 290	303.0510	303.0522	3.71	217 (100)	C-1, C-6, C-9	D	
Benzenetriol	Pyrogallol	17	38	C ₆ H ₆ O ₃	14.146	214, 271	125.0244	125.0243	-1.02	125 (100), 81 (13), 97 (11)	BE	D	
Benzopyran	7,8-Dihydroxy-3-propyl-	124	265	C ₁₂ H ₁₆ O ₅	63.686	228, 370	239.0925	239.0897	-11.70	124 (100), 168 (29)	GC-9,	D	

	3,4,5,6,7,8-hexahydro-1 <i>H</i> -2-benzopyran-1,5-dione										BC-9		
	8-Hydroxycannabinol	128	270	C ₂₁ H ₂₆ O ₃	66.719	200, 222, 254, 486	325.1809	325.1828	5.79	325 (100)	C-9, GC-9, BC-9, GM-9, BM-9	D	
C-Glycosyl compound	Unknown C-glycosyl compound (5-ethyl-2-(hydroxymethyl)-1 <i>H</i> -indol-3-yl 3,4,5-trihydroxy-5-(hydroxymethyl)oxolane-2-carboxylate or C-glycosyl compound type)	60	143	C ₁₇ H ₂₁ NO ₈	32.078	202, 224, 282	366.1194	366.1219	6.69	142 (100), 322 (79), 300 (48)	C-1, C-9	D	
Coumarin glycoside	Aspergillazine C 1	86	197	C ₂₀ H ₂₂ N ₂ O ₈ S	45.255	194, 270	449.1024	449.1039	3.21	286 (100), 199 (30)	C-9	D	
	Aspergillazine C 2	94	213	C ₂₀ H ₂₂ N ₂ O ₈ S	48.399	198, 266, 350	449.1024	449.1057	7.37	217 (100), 147 (55), 156 (54)	C-9	D	
Gallotannin	Trigalloyl glucopyranose	60	145	C ₂₇ H ₂₄ O ₁₈	32.078	224, 282	635.0890	635.0940	7.87	565 (100), 424 (67)	BE, GC-1, BC-1, GM-1, GM-3, GM-6, GM-9	D, L	(4)
Galloyl ester	(3,4,5,6-Tetrahydroxyoxan-2-yl)methyl 3,4,5-trihydroxybenzoate	8	20	C ₁₃ H ₁₆ O ₁₀	10.415	196, 226sh, 275	331.0671	331.0636	-10.48	169 (100), 211 (44), 271 (25)	GC-9, GM-9, BM-9	D	
	2,3-Dihydroxypropyl 3,4,5-trihydroxybenzoate	26	57	C ₁₀ H ₁₂ O ₇	17.301	212, 270, 308	243.0510	243.0519	3.72	107 (100), 106 (20), 71 (20)	GC-1, BC-1, GC-3, BC-3, GC-6, BC-6, GC-9, BC-9	D	
Hydrolyzable tannin	Strictinin	45	109	C ₂₇ H ₂₂ O ₁₈	25.141	196, 222, 288, 326sh	633.0733	633.0790	9.00	302 (100), 631 (4)	BE, GC-1, BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, GC-6, GM-6, BM-6, GM-9	D, L	(16)
Hydroxycoumarin	Esculetin	66	157	C ₉ H ₆ O ₄	35.542	334	177.0193	177.0197	2.08	109 (100), 91 (96), 152 (39)	BC-1, BM-1, BC-3, BM-3, BM-6, BC-9	D	
O-Benzoquinone	5,6-Dioxocyclohexa-1,3-diene-1-carboxylic acid	107	232	C ₇ H ₄ O ₄	55.163	192, 262, 300sh	151.0037	151.0022	-9.81	65 (100), 83 (23), 107 (20)	GE, BE, BC-1,	D	

Phenol ether	Unknown phenol ether (4-[3-(4-amidinophenoxy)propoxy]benzamidine or other phenol ether type)	62	148	$C_{21}H_{32}N_4O_{10}S_2$	33.214	202, 236, 272sh, 336	563.1487	563.1472	-2.75	504 (100), 387 (83), 147 (76)	GC-3, BC-3, GC-6, GC-9, BC-9, BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9	D
Phenolic glycoside	Turgorin	21	43	$C_{13}H_{16}O_{13}S$	15.458	194, 340	411.0239	411.0237	-0.45	97 (100), 242 (32), 139 (11)	GC-1, BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9	D
Styrene	4-Vinylphenol	73	168	C_8H_8O	38.075	230, 313	119.0502	119.0515	10.92	119 (100), 93 (6)	C-1, BC-1, C-6, C-9	D
Tetracenequinone	Maggiemycin	97	218	$C_{22}H_{18}O_9$	49.699	208, 282	425.0878	425.0829	-11.54	255 (100), 133 (31), 205 (24)	GC-9, GM-9	D
Others: non-phenolic compounds												
Alpha amino acid	Pyroglutamic acid	4	9	$C_5H_7NO_3$	8.277	196, 262	128.0353	128.0336	-13.25	88 (100)	C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9,	D

Beta hydroxy acid	<i>L</i> -Malic acid	2	3	C ₄ H ₆ O ₅	6.849	194, 210sh, 266	133.0142	133.0130	-9.75	73 (100), 89 (9)	GM-9, BM-9 BC-1, GM-1, BM-1, GM-3, C- 6, GC-6, BC-6, GM-6, BM-6, C- 9, GC-9, BC-9	D
	3-Hydroxyglutaric acid	7	15	C ₅ H ₈ O ₅	9.797	194, 205sh, 263	147.0299	147.0301	1.24	87 (100)	C-1, GC- 1, BC-1, C-3, GC- 3, BC-3, C-6, GC- 6, BC-6, C-9, GC- 9, BC-9	D
Dicarboxylic acid	Succinic acid	6	13	C ₄ H ₆ O ₄	9.325	198, 270	117.0193	117.0169	-20.70	73 (100), 99 (6)	C-9, GC- 9	D
Dipeptide	gamma-Glutamylleucine	10	23	C ₁₁ H ₂₀ N ₂ O ₅	11.194	194, 280	259.1299	259.1303	1.52	215 (100), 171 (39), 102 (24)	C-1, BC- 1, C-3, C-9, BC- 9	D
	2-(2-Amino-4-methylpentanamido)-3-methylbutanoic acid 3	33	72	C ₁₁ H ₂₂ N ₂ O ₃	20.888	194, 222, 285, 333	229.1558	229.1574	7.31	130 (100), 115 (13)	C-1, BC- 1, C-3, BC-3, C- 6, BC-6, C-9, GC- 9, BC-9	D
	2-(2-Amino-4-methylpentanamido)-3-phenylpropanoic acid 1	61	146	C ₁₅ H ₂₂ N ₂ O ₃	32.412	204, 280	277.1558	277.1543	-5.29	164 (100), 153 (11)	C-1, BC- 1, C-3, C-9, GC- 9, BC-9	D
Estrogen	Estradiol	112	240	C ₁₈ H ₂₄ O ₂	57.697	317	271.1704	271.1678	-9.42	209 (100), 128 (71), 84 (47)	BC-1, C- 9, BC-9	D
Fatty aldehyde	Triptophenolide	127	269	C ₂₀ H ₂₄ O ₃	66.239	208, 280, 325	311.1653	311.1671	5.89	311 (100)	C-9, GC- 9, BC-9, GM-9, BM-9	D
Furoic acid	2-Furoic acid	26	60	C ₅ H ₄ O ₃	17.301	250, 308	111.0088	111.0072	-14.40	111 (100)	GC-9, BC-9, BM-9	D
Gamma butyrolactone	Parvistemoamide	47	117	C ₁₂ H ₁₉ NO ₄	25.832	270, 315	240.1241	240.1245	1.53	160 (100), 99 (19)	C-3, GM- 3, C-6, C-9	D
Gamma-keto acid	4,5-Dioxopentanoic acid	34	75	C ₅ H ₆ O ₄	21.375	206, 274	129.0193	129.0190	-2.57	85 (100), 101 (33)	C-9, GC- 9, BC-9	D
Gluconolactone	Gluconolactone	13	30	C ₆ H ₁₀ O ₆	12.689	204, 297	177.0405	177.0405	0.44	73 (100), 87 (9), 135 (5)	GC-1, BC-1, BM-1, C- 3, BC-3, C-6, GC- 6, BC-6,	D

Glutamic acid	Saccharopine	2	4	C ₁₁ H ₂₀ N ₂ O ₆	6.849	194, 210sh, 266	275.1249	275.1202	-16.79	115 (100)	GM-6, BM-6, C- 9, GC-9, BC-9 C-3, GC- 3, BC-3, GM-3	D
Glutamine	Theanine	3	6	C ₇ H ₁₄ N ₂ O ₃	7.270	200, 248, 280sh	173.0932	173.0926	- 3.1543 71	155 (100), 84 (49)	BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9	D
Hippuric acid	Herdmanine H	36	83	C ₁₃ H ₁₈ N ₄ O ₄	22.002	220, 282	293.1255	293.1253	-0.64	89 (100), 129 (51), 119 (39), 175 (23), 188 (16)	C-9	D
Histrionicotoxin	Histrionicotoxin	12	26	C ₁₉ H ₂₅ NO	12.358	204, 256	282.1863	282.1879	5.54	150 (100)	BM-3, GM-6, BM-6	D
Hydroxy fatty acid	Mevalonic acid	9	22	C ₆ H ₁₂ O ₄	10.823	194, 280	147.0663	147.0645	-12.12	85 (100)	GC-9	D
Iridoid <i>O</i> -glycoside	Caryoptosidic acid 1	13	29	C ₁₆ H ₂₄ O ₁₁	12.689	208, 327	391.1246	391.1226	-5.02	243 (100), 242 (87), 168 (34)	C-3, BC- 3, C-6, BC-6, C- 9	D
	Caryoptosidic acid 2	22	47	C ₁₆ H ₂₄ O ₁₁	15.965	264, 338	391.1246	391.1231	-3.69	277 (100), 242 (81), 263 (47), 228 (43), 358 (39)	C-3, BC- 3, C-6, C-9	D
Lineolic acid	Corchorifatty acid F	122	262	C ₁₈ H ₃₂ O ₅	63.212	210, 305	327.2177	327.2167	-3.05	222 (100), 311 (96), 165 (89), 209 (64)	BC-9, BM-9	D
Long-chain fatty acid	Myristic acid	29	64	C ₁₄ H ₂₈ O ₂	19.574	206, 275	227.2017	227.2007	-4.20	112 (100), 184 (54), 107 (49), 159 (35)	BC-3	D
Methyl-branched fatty acid	3-Methylpent-2-enedioic acid	23	50	C ₆ H ₈ O ₄	16.276	195, 205sh, 280, 320	143.0350	143.0350	4.74	72 (100), 98 (55)	C-1, GC- 1, BC-1, C-3, GC- 3, BC-3, C-6, GC- 6, BC-6, C-9, GC- 9, BC-9	D
Phenylalanine	L-Phenylalanine	20	40	C ₉ H ₁₁ NO ₂	14.758	236, 296	164.0717	164.0691	-15.86	103 (100), 147 (47), 72 (21)	C-9	D
Purine 2'-deoxyribonucleoside	9-[4-Hydroxy-5-(hydroxymethyl)oxolan-2-yl]-2-imino-3,9-dihydro-2 <i>H</i> -purine-6,8-diol	11	24	C ₁₀ H ₁₃ N ₅ O ₅	11.528	210, 260	282.0844	282.0846	0.84	151 (100), 133 (27)	GM-9, BM-9	D
Purine nucleoside	2-({9-[3,4-dihydroxy-5-(hydroxymethyl)oxolan-	26	58	C ₁₄ H ₁₇ N ₅ O ₈	17.301	196, 266, 328	382.1004	382.1054	12.94	74 (100)	C-1, GC- 1, BC-1,	D

	2-yl]-9H-purin-6-yl} amino)butanedioic acid										C-3, BC-6, C-9, GC-9, BC-9, BM-9	
Pyranone	4-Methoxy-6-(3-phenyloxiran-2-yl)-2H-pyran-2-one	5	11	C ₁₄ H ₁₂ O ₄	8.902	198, 270	243.0663	243.0623	-16.51	139 (100), 85 (64)		D
Pyrazine	2-(1-Hydroxy-1-methylpropyl)-3-methoxypyrazine	123	263	C ₉ H ₁₄ N ₂ O ₂	63.570	210, 305	181.0983	181.0959	-12.98	153 (100), 79 (50)	BC-3, GM-3, BM-3	D
Pyrimidone	5-Ethyl-5-(4-hydroxypentan-2-yl)-1,3-diazinane-2,4,6-trione	48	121	C ₁₁ H ₁₈ N ₂ O ₄	26.298	258	241.1194	241.1201	2.86	130 (100), 168 (66)	C-1, C-3, GM-3, C-6, C-9, BC-9	D
Pyrrole	2-Heptyl-1H-pyrrole	21	44	C ₁₁ H ₁₉ N	15.458	194, 210sh, 289	164.1445	164.1443	-1.05	120 (100)	C-3, C-6	D
Stilbene	3,5-Dihydroxy-4-(2,3-epoxy-3-methylbutyl)bibenzyl	126	268	C ₁₉ H ₂₂ O ₃	65.720	208, 280, 325	297.1496	297.1531	11.72	297 (100)	C-9, GC-9, BC-9, GM-9, BM-9	D
Sugar acid	1,3-Diethyl 2-hydroxypropanedioate	35	80	C ₇ H ₁₂ O ₅	21.850	220, 282	175.0612	175.0622	5.67	115 (100), 85 (60)	C-1, BC-1, GM-1, C-3, GC-3, BC-3, C-6, GC-6, BC-6, C-9, GC-9, BC-9, GM-9	D
Terpene glycoside	Rhodiolide C	117	251	C ₂₂ H ₃₈ O ₁₂	59.227	206, 316	493.2291	493.2305	2.94	293 (100), 89 (81), 119 (62), 404 (61)	GM-9	D
Terpene lactone	Heritol	45	110	C ₁₅ H ₁₆ O ₃	25.141	316	243.1027	243.1062	14.49	63 (100), 83 (68)	C-6, C-9	D
	Amberboin	110	238	C ₁₅ H ₂₀ O ₄	56.333	192, 260	263.1289	263.1294	1.96	147 (100), 71 (83)	BC-1, BM-6, BC-9, BM-9	D
Tricarboxylic acid	Citric acid	4	7	C ₆ H ₈ O ₇	8.277	198, 265	191.0197	191.0201	1.85	111 (100), 87 (47)	C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	D
	3-Hydroxy-3-carboxymethyl-adipic acid	27	61	C ₈ H ₁₂ O ₇	18.638	212, 272	219.0510	219.0485	-11.58	112 (100), 95 (19), 155 (13)	C-9, GC-9, BC-9	D
Tropone	N-{9-hydroxy-3,4,5,14-tetramethoxy-13-	75	172	C ₂₁ H ₂₃ NO ₇	39.040	340	400.1402	400.1398	-0.94	357 (100), 204 (39), 342 (39)	GC-1, BC-3,	D

Tyrosine	oxotricyclo[9.5.0.0 ^{2,7}]hexadeca-1(16),2(7),3,5,11,14-hexaen-10-yl}formamide <i>L</i> -Tyrosine	7	14	C ₉ H ₁₁ NO ₃	9.797	194, 205sh, 263	180.0666	180.0644	-12.42	93 (100)	GC-9, BC-9, GM-9, BM-9	D	
Vinylogous acid	3,4-Dihydroxy-5-(3-methylbut-2-en-1-yl)-2-(3-methylbutanoyl)cyclopent-2-en-1-one	125	267	C ₁₅ H ₂₂ O ₄	65.555	208, 280, 325	265.1445	265.1476	11.69	265 (100), 97 (42)	C-9, GC-9, BC-9, BM-9	D	
<i>Positive mode [M+H]⁺</i>													
Phenolic acids													
Hydroxybenzoic acid	Gallic acid	17	37	C ₇ H ₆ O ₅	14.146	218, 274	171.0288	171.0289	0.70	79 (100)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	S, D, L	(6), (10), (14), (17)
	Galloylquinic acid	18	39	C ₁₄ H ₁₆ O ₁₀	14.341	216, 276	345.0816	345.0841	7.18	154 (100), 155 (18), 111 (13),	GE, BE, GC-1, GM-1, GC-3, BC-3, GM-3, BM-3, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9	D, L	(6), (17)
	Syringic acid 1	37	85	C ₉ H ₁₀ O ₅	22.193	230, 298	199.0601	199.0602	0.70	128 (100)	BC-1	D	
	<i>p</i> -Hydroxybenzoic acid	46	113	C ₇ H ₆ O ₃	25.502	204, 226sh, 276	139.0390	139.0388	-1.44	139 (100), 67 (9)	GE, BE	S, D	
	Syringic acid	64	152	C ₉ H ₁₀ O ₅	34.439	218, 275	199.0601	199.0608	3.72	109 (100), 153 (57), 81 (22)	C-1, GC-1, BC-1, GM-1,	S, D	

											BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9		
Hydroxycinnamic acid	3-(4-Hydroxyphenyl)-N-{4-[3-(4-hydroxyphenyl)prop-2-enamido]butyl}prop-2-enamide	15	34	C ₂₂ H ₂₄ N ₂ O ₄	13.587	194, 282	381.1809	381.1788	-5.55	381 (100)	C-3, GC-3, BC-3, GM-3	D	
	Glutathionyl caftaric acid	27	62	C ₂₃ H ₂₇ N ₃ O ₁₅ S	18.638	196, 231, 255, 330	618.1236	618.1269	5.40	211 (100), 146 (60), 130 (48), 194 (47)	C-1, GC-1, BC-1, C-3, GC-3, BC-3, C-6, GC-6, BC-6, C-9, GC-9, BC-9	D, L	(14), (17)
	Feruloylagmatine 1	38	90	C ₁₅ H ₂₂ N ₄ O ₃	22.477	212, 234sh, 274	307.1765	307.1724	-13.24	139 (100), 163 (7)	GC-3, GM-3, GC-6, GM-9	D	
	Feruloylagmatine 2	44	107	C ₁₅ H ₂₂ N ₄ O ₃	24.894	194, 238, 272	307.1765	307.1723	-13.58	140 (100), 181 (69)	GC-3, GM-3	D	
	Chlorogenic acid	44	108	C ₁₆ H ₁₈ O ₉	24.894	242, 327	355.1024	355.1056	9.01	164 (100), 145 (65), 135 (11)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	S, D, L	(6), (10)
	Grandidentoside	45	111	C ₂₁ H ₂₈ O ₁₀	25.141	203, 232, 280	441.1755	441.1736	-4.41	441 (100)	BE	D	
	Subaphyllin	48	122	C ₁₄ H ₂₀ N ₂ O ₃	26.298	246, 356	265.1547	265.1550	1.47	120 (100)	C-1, GM-3, C-9, GC-9, BC-9	D	
	3,4-Dimethoxycinnamic acid	120	259	C ₁₁ H ₁₂ O ₄	60.747	224, 327	209.0808	209.0817	4.14	135 (100), 117 (91), 89 (32)	C-1, GC-1, BC-1,	D, L	(14)

Hydroxyphenylpropanoic acid	2-(2-Amino-4-methylpentanamido)-3-phenylpropanoic acid 1	61	146	C ₁₅ H ₂₂ N ₂ O ₃	32.412	204, 280	279.1703	279.1714	3.94	83 (100)	C-3, GC-3, BC-3, C-6, GC-6, BC-6 C-1, BC-1, C-3, C-9, GC-9, BC-9	D	
	2-(2-Amino-4-methylpentanamido)-3-phenylpropanoic acid 2	65	153	C ₁₅ H ₂₂ N ₂ O ₃	35.315	208, 281	279.1703	279.1648	-19.70	86 (100), 120 (14)	C-9	D	
Flavonoids													
Dihydrochalcone	2',6'-Dihydroxy-4'-methoxy-3'-methylidihydrochalcone	111	239	C ₁₇ H ₁₈ O ₄	56.810	228, 270	287.1278	287.1256	-7.61	287 (100), 187 (19)	C-1, GC-1, BC-1, BC-3	D	
Flavan-3-ol	Theasinensin C 1	24	51	C ₃₀ H ₂₆ O ₁₄	16.471	206, 275	611.1395	611.1475	12.97	317 (100), 425 (58), 308 (35), 275 (30)	GM-1, GM-3, GM-6	D, L	(4)
	Gallocatechin	28	63	C ₁₅ H ₁₄ O ₇	19.054	200, 216, 273	307.0812	307.0834	7.07	139 (100), 195 (7)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	S, D, L	(2), (6), (17)
	Theasinensin C 3	32	70	C ₃₀ H ₂₆ O ₁₄	20.509	206, 276	611.1395	611.1446	8.35	69 (100), 251 (66), 203 (63), 270 (31)	GC-1, GM-1, GC-3, GM-3, GC-6, GM-6	D, L	
	Theasinensin B 1	37	84	C ₃₇ H ₃₀ O ₁₈	22.193	230, 298	763.1505	763.1641	17.82	275 (100), 449 (63), 283 (58), 409 (25), 261 (20)	BE, BC-1, GM-1, BM-1, BC-3, GM-3, BM-3, BC-6, GM-6, BM-6	D, L	(4)
	Leucodelphinidin	37	87	C ₁₅ H ₁₄ O ₈	22.193	214, 250, 270	323.0761	323.0775	4.23	161 (100)	GM-1	D	
	Epigallocatechin	40	97	C ₁₅ H ₁₄ O ₇	23.385	196, 236, 272	307.0800	307.0843	14.00	163 (100), 140 (67), 205 (13)	GE, BE, C-1, GC-1, BC-1, GM-1,	S, D, L	(2), (17)

Theasinensin B 2	43	104	C ₃₇ H ₃₀ O ₁₈	24.428	212, 276	763.1505	763.1583	10.22	277 (100), 425 (87), 139 (76), 577 (27)	BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9 GE, BE, GC-1, GM-1, GC-3, GM-3, GC-6, GM-6, GC-9, GM-9	D, L	(4)
Catechin	46	112	C ₁₅ H ₁₄ O ₆	25.502	206, 224sh, 285	291.0900	291.0871	-9.96	123 (100), 111 (20)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9 BE, GC-1, GM-1, GC-3, GM-3, GC-6, GM-6, GC-9, GM-9	S, D, L	(4), (6), (14)
(Epi)gallocatechin-(epi)catechin gallate 1	51	127	C ₃₇ H ₃₀ O ₁₇	27.305	208, 281	747.1556	747.1642	11.51	139 (100), 290 (47), 595 (40), 275 (38)	BE, GC-1, GM-1, GC-3, GM-3, GC-6, GM-6, GC-9, GM-9	D, L	(9)
Theasinensin A	52	128	C ₄₄ H ₃₄ O ₂₂	27.498	210, 282	915.1614	915.1708	10.27	288 (100), 575 (18), 449 (14)	GE, BE, GC-1, BC-1, BM-1, GC-3, BC-3, BM-3	D, L	(4)
(Epi)gallocatechin-(epi)catechin gallate 2	54	131	C ₃₇ H ₃₀ O ₁₇	28.239	204, 280	747.1556	747.1641	11.38	123 (100), 273 (17), 315 (10)	GC-1, GM-1	D, L	(9)

Epicatechin	56	135	C ₁₅ H ₁₄ O ₆	29.502	216, 276	291.0863	291.0895	10.94	147 (100), 165 (83), 140 (54)	GC-3, GM-3, GC-6, GM-6, GC-9, GM-9 GE, BE, C-1, GC- 1, BC-1, GM-1, BM-1, C- 3, GC-3, BC-3, GM-3, BM-3, C- 6, GC-6, BC-6, GM-6, BM-6, C- 9, GC-9, BC-9, GM-9, BM-9	S, D, L	(4), (6), (14)
Epigallocatechin gallate	58	137	C ₂₂ H ₁₈ O ₁₁	30.555	196, 232, 276	459.0922	459.0983	13.29	139 (100), 151 (14), 289 (9)	GE, BE, C-1, GC- 1, BC-1, GM-1, BM-1, C- 3, GC-3, BC-3, GM-3, BM-3, C- 6, GC-6, BC-6, GM-6, BM-6, GC-9, GM-9, BM-9	S, D, L	(4), (6)
Galocatechin gallate dimer	58	140	C ₄₄ H ₃₆ O ₂₂	30.555	196, 232, 276	917.1771	917.1899	13.96	153 (100), 140 (78), 271 (28)	GC-1, GM-1, GC-3, GM-3, GC-6, GM-6	D, L	(7)
Galocatechin gallate	63	149	C ₂₂ H ₁₈ O ₁₁	33.413	210, 275, 334	459.0900	459.0976	16.55	140 (100), 120 (31), 247 (25)	GE, BE, C-1, GC- 1, BC-1, GM-1, BM-1, C- 3, GC-3, BC-3, GM-3, BM-3, C- 6, GC-6, BC-6, GM-6,	S, D	

(Epi)catechin gallate dimer	76	174	C ₄₄ H ₃₆ O ₂₀	39.622	196, 232, 280	885.1873	885.2010	15.51	257 (100)	BM-6, GM-9, GE, BE, GC-1, BC-1, GM-1, GC-3, BC-3, GM-3, GC-6, GM-6	D, L	(20)
Epicatechin gallate	76	176	C ₂₂ H ₁₈ O ₁₀	39.622	196, 232, 280	443.0973	443.1032	13.32	153 (100), 165 (51), 140 (35), 124 (30)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, GM-9, BM-9	S, D, L	(2), (6), (9)
Epicatechin gallate 1	79	182	C ₂₂ H ₁₈ O ₁₀	41.522	208, 276	443.0973	443.1013	9.09	123 (100)	GE, GC-1, GM-1, GC-3, GM-3, GC-6, GM-6	D, L	(2), (6), (9)
2,3-Dihydroxy-5-(3,5,7-trihydroxy-4-oxo-3,4-dihydro-2H-1-benzopyran-2-yl)phenyl 3,4-dihydroxy-5-methoxybenzoate	81	191	C ₂₃ H ₁₈ O ₁₂	42.668	210, 258, 357	487.0871	487.0903	6.57	325 (100)	BE, GC-1, BC-1, GM-1, BM-1, BC-3	D	
Epicatechin 3-O-gallate and dehydroascorbic acid conjugate	88	199	C ₂₈ H ₂₄ O ₁₆	46.074	198, 268, 348	617.1137	617.1548	66.60	550 (100)	GE, BE, GC-1, BC-1, GM-1, BM-1, GC-3	D, L	(23)
Methyl epicatechin gallate	90	204	C ₂₃ H ₂₀ O ₁₀	46.533	194, 280	457.1129	457.1130	0.17	123 (100)	GE, BE, GM-1, GC-3, GM-3	D, L	(2)
Epiafzelechin 3-gallate	92	209	C ₂₂ H ₁₈ O ₉	47.594	194, 206sh, 275	427.1024	427.1068	10.40	257 (100), 140 (94), 109 (42)	GE, BE, GC-1, GM-1, GC-3, GM-3	D, L	(11)
C-8 N-ethyl-2-	94	211	C ₂₈ H ₂₇ NO ₁₁	48.399	194, 210, 280	554.1657	554.1710	9.56	384 (100), 217 (18)	GE, GC-	D, L	(12)

	pyrrolidinone-substituted (-)-epicatechin-3- <i>O</i> - gallate									1, GC-9, GM-9		
	Epigallocatechin gallate- 4',4''- <i>O</i> -β-D-gluco- pyranoside	98	219	C ₃₄ H ₃₈ O ₂₁	50.154	248, 364	783.1978	783.2441	59.07	189 (100), 111 (40), 329 (19)	GE, GC- 1, GM-1	D, L (22)
	Theaflavate B	107	233	C ₃₆ H ₂₈ O ₁₅	55.163	326	701.1501	701.5045	505.46	701 (100)	GE, BE, GC-1, BC-1, C- 3, GC-3, BC-3	D, L (19)
	Theaflavin	115	247	C ₂₉ H ₂₄ O ₁₂	58.863	210, 268, 298, 365sh	565.1341	565.1412	12.65	139 (100)	GE, BE, BC-1, BC-3	D, L (6)
	Theaflavin 3-gallate	118	255	C ₃₆ H ₂₈ O ₁₆	59.832	210, 270	717.1450	717.1531	11.29	151 (100), 379 (84), 229 (42), 367 (38), 241 (31), 111 (30)	GE, BE, BC-1, BC-3	D, L (6)
	Theaflavin-3,3'-gallate	119	257	C ₄₃ H ₃₂ O ₂₀	60.230	208, 277, 378	869.1560	869.1664	12.00	391 (100), 153 (98), 333 (45), 593 (42), 441 (13)	GE, BE, BC-1, BC-3	D, L (6)
Flavanonol Flavone	Katuranin 1	58	139	C ₁₅ H ₁₂ O ₆	30.555	212, 275	289.0707	289.0707	-0.05	139 (100), 151 (14)	GE, BE	D
	Apigenin 6,8-di- <i>C</i> - glucoside	55	132	C ₂₇ H ₃₀ O ₁₅	29.077	272, 352	595.1657	595.1745	14.71	559 (100), 439 (85), 421 (57), 294 (31)	BE, BC- 1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, BC-6, GM-6, BM-6	D, L (4)
	Apigenin 7- <i>O</i> -apiosyl- glucoside	63	150	C ₂₆ H ₂₈ O ₁₄	33.413	208, 270, 338	565.1552	565.1553	0.16	565 (100), 547 (100), 529 (82), 427 (78), 511 (56)	GE, BE, GC-9, BC-9, GM-9, BM-9	D
Flavonol	3,5,7-Trihydroxy-2-(3- hydroxy-4- methoxyphenyl)-6,8- dimethoxy-4 <i>H</i> -chromen- 4-one	47	116	C ₁₈ H ₁₆ O ₉	25.832	194, 222, 286, 320sh	377.0867	377.0897	7.93	377 (100)	GM-1, GM-3	D
	Quercetin 3- <i>O</i> -di- galactoside/galucoside	65	154	C ₂₇ H ₃₀ O ₁₇	35.315	208, 281	627.1556	627.1631	12.00	319 (100), 320 (24)	GE, BE, C-1, GC- 1, BC-1, GM-1, BM-1, C- 3, GC-3, BC-3, GM-3, BM-3, C- 6, GC-6, BC-6, GM-6, BM-6	D, L (5)
	Myricetin-3- <i>O</i> - galactoside 1	68	159	C ₂₁ H ₂₀ O ₁₃	36.379	212, 270, 350	481.0977	481.1023	9.63	320 (100)	GE, C-1, GC-1,	D, L (10)

Myricetin-3- <i>O</i> -galactoside 2	69	160	C ₂₁ H ₂₀ O ₁₃	36.799	210, 264, 360	481.0977	481.1036	12.33	320 (100), 97 (15)	BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GE, BE, GC-1, BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, GC-6, GM-6, GE, BE, BC-1, BM-1, GC-3, BC-3, GM-3, BM-3, BC-6, BM-6, GC-1, GM-1, GC-9, GM-9	D, L	(10)
Quercetin 3- <i>O</i> -glucosylrutinoside 1	70	161	C ₃₃ H ₄₀ O ₂₁	37.066	206, 268, 350	773.2135	773.2271	17.61	304 (100), 449 (12)	BC-1, BM-1, GC-3, BC-3, GM-3, BM-3, BC-6, BM-6, GC-1, GM-1, GC-9, GM-9, GE, BE, GC-1, BC-1, GM-1, BM-1, BC-3, BM-3, GC-6, GM-6, GC-9, GM-9, BM-9	D, L	(4), (9)
Myricetin	71	162	C ₁₅ H ₁₀ O ₈	37.156	210, 264, 360	319.0448	319.0481	10.34	319 (100)	GC-1, GM-1, GC-9, GM-9	D	
Kaempferol 3- <i>O</i> -rutinoside 1	72	164	C ₂₇ H ₃₀ O ₁₅	37.274	216, 271, 347	595.1658	595.1732	12.43	433 (100), 379 (16), 283 (12)	GE, BE, GC-1, BC-1, GM-1, BM-1, BC-3, BM-3, GC-6, GM-6, GC-9, GM-9, BM-9	D, L	(6)
Quercetin 3- <i>O</i> -glucosylrutinoside 2	73	167	C ₃₃ H ₄₀ O ₂₁	38.075	208, 260, 360	773.2135	773.2219	10.88	303 (100)	GE, BE, GC-1, BC-1, GM-1, BM-1, BC-3, BM-3, GC-6, BM-6, GC-9,	D, L	(4), (9)

Quercetin-3- <i>O</i> -glucoside 1	76	175	C ₂₁ H ₂₀ O ₁₂	39.622	345	465.1028	465.0879	-32.02	303 (100)	BC-9, GM-9, BM-9, BC-1, BM-1, BC-3, BM-3, BM-6	D, L	(9)
Rutin	77	178	C ₂₇ H ₃₀ O ₁₆	40.518	204, 258, 357	611.1610	611.1676	10.80	304 (100)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9	S, D, L	(2), (4)
Myricetin-galloyl-hexoside 1	77	179	C ₂₈ H ₂₄ O ₁₇	40.518	202, 256, 354	633.1086	633.1430	54.31	633 (100)	GE	D, L	(16)
Myricetin-galloyl-hexoside 2	79	185	C ₂₈ H ₂₄ O ₁₇	41.522	208, 258, 356	633.1086	633.1578	77.64	633 (100), 331 (2)	BE	D, L	(16)
Quercetin-3- <i>O</i> -glucoside 2	80	186	C ₂₁ H ₂₀ O ₁₂	42.306	208, 258, 356	465.1028	465.1061	7.22	304 (100), 127 (20)	GE, BE, GC-1, GM-1, GC-3, GM-3, GC-6, BC-6, GM-6	D, L	(9)
Kaempferol 3- <i>O</i> -glucosylrhamnosyl-galactoside 2	80	187	C ₃₃ H ₄₀ O ₂₀	42.306	198, 268, 349	757.2186	757.2331	19.19	288 (100), 449 (79), 217 (6)	GE, BE, BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, BC-6, BM-6, BC-9, BM-9	D	
Kaempferol 1	80	188	C ₁₅ H ₁₀ O ₆	42.306	198, 268, 349	287.0550	287.0601	17.77	122 (100)	BE, BC-1, BC-3, BC-6, BM-6	D, L	(5)
Quercetin 3- <i>O</i> -[(2,4-diacetyl- α -L-rhamnopyranosyl)-	80	189	C ₃₅ H ₃₈ O ₂₀	42.306	198, 268, 349	779.2029	779.2159	16.68	493 (100)	BE, BC-1, GM-1, BM-1	D	

(1→6)]-3,4-diacetyl-beta-D-galactopyranoside 1 Quercetin 3-O-galactoside	82	192	C ₂₁ H ₂₀ O ₁₂	42.886	208, 258, 357	465.1028	465.1080	11.28	303 (100)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9	S, D, L	(6), (17)
Kaempferol 3-O-rutinoside 2	83	193	C ₂₇ H ₃₀ O ₁₅	43.739	206, 254, 342	595.1657	595.1694	6.14	288 (100), 449 (39), 147 (18)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6	D, L	(6)
Kaempferol 3-O-rutinoside 3	87	198	C ₂₇ H ₃₀ O ₁₅	45.751	198, 268, 349	595.1657	595.1726	11.59	287 (100), 288 (18)	GE, BE, GC-1, BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, GC-6, GM-6, GC-9, BC-9, GM-9, BM-9	D, L	(6)
Kaempferol 2	88	200	C ₁₅ H ₁₀ O ₆	46.074	196, 266, 348	287.0550	287.0551	0.10	287 (100)	GE, BE, BC-9, BM-9	D, L	(5)
Kaempferol 3-O-glucoside 1	89	201	C ₂₁ H ₂₀ O ₁₁	46.401	194, 280	449.1078	449.1139	13.50	91 (100)	GE, BE, BC-1, BC-3, GC-6, BC-6, BM-6	D, L	(9)

Kaempferol 3- <i>O</i> -glucoside 2	93	210	C ₂₁ H ₂₀ O ₁₁	47.827	280	449.1078	449.0938	-31.17	288 (100), 127 (15)	BE, GM-1	D	
Kaempferol 3- <i>O</i> -glucoside	94	212	C ₂₁ H ₂₀ O ₁₁	48.399	198, 266, 350	449.1078	449.1132	11.94	288 (100), 85 (14), 127 (11)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9	S, D, L	(9)
Kaempferol 3	95	215	C ₁₅ H ₁₀ O ₆	48.707	198, 268, 348	287.0550	287.0574	8.36	153 (100)	GC-1, BC-1, GM-1, BM-1, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9	D, L	(5)
Unknown quercetin conjugate 1	113	241	C ₄₂ H ₄₆ O ₂₂	58.117	210, 270, 318	903.2554	903.2670	12.90	147 (100), 303 (27), 309 (24), 291 (13)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9	D, L	(6)
Unknown kaempferol conjugate	115	248	C ₄₂ H ₄₆ O ₂₁	58.863	212, 270, 315	887.2604	887.2701	10.89	147 (100), 85 (13)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6	D, L	(6)

	Quercetin 3- <i>O</i> -[(2,4-diacetyl- α -L-rhamnopyranosyl)-(1 \rightarrow 6)]-3,4-diacetyl-beta-D-galactopyranoside 2	117	253	C ₃₅ H ₃₈ O ₂₀	59.227	210, 316	779.2029	779.1913	-14.89	779 (100)	GM-6, BM-6, GM-1	D	
	Quercetin	121	261	C ₁₅ H ₁₀ O ₇	61.100	206, 258, 374	303.0499	303.0524	8.25	247 (100), 177 (57)	GE, BE, GC-1, GM-1, BM-1, GC-3, GM-3, BM-3, GC-6, GM-6, BM-6, GC-9, GM-9	S, D, L	(17)
	Kaempferol	124	266	C ₁₅ H ₁₀ O ₆	63.686	228, 370	287.0550	287.0575	8.71	165 (100)	GE, BE, GC-1, GM-1, GC-3, GM-3, GM-6, GC-9, GM-9, BM-9	S, D, L	(5)
Hydroxyflavonoid	3,5,7-Tris(acetyloxy)-2-[4-(acetyloxy)-3-hydroxyphenyl]-4 <i>H</i> -1-benzopyran-4-one 1	91	207	C ₂₃ H ₁₈ O ₁₁	46.86	194, 210, 280	471.0922	471.0938	3.40	471 (100)	BE, GC-1, BC-1, GM-1, BC-3	D	
	3,5,7-Tris(acetyloxy)-2-[4-(acetyloxy)-3-hydroxyphenyl]-4 <i>H</i> -1-benzopyran-4-one 2	95	214	C ₂₃ H ₁₈ O ₁₁	48.707	198, 268, 348	471.0922	471.0955	7.00	309 (100)	BE, GC-1, BC-1, GM-1, BM-1	D	
Isoflavonoid	3-Oxo-3-[[3,4,5-trihydroxy-6-(3-hydroxy-14-methoxy-8,17-dioxatetracyclo[8.7.0.0 ² ,7 ⁰ ,11 ¹ ,6 ⁶]heptadeca-2,4,6,11,13,15-hexaen-5-yl]oxy)oxan-2-yl]methoxy}propanoic acid	75	173	C ₂₅ H ₂₆ O ₁₃	39.040	220, 274, 338	535.1446	535.1539	17.35	517 (100), 481 (87), 379 (64), 427 (43)	BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, BC-6, BM-6, GC-9, BC-9, GM-9, BM-9	D	
<i>O</i> -Methylated flavonoid	Centaureidin	42	101	C ₁₈ H ₁₆ O ₈	24.075	312	361.0918	361.0929	3.05	361 (100), 169 (58)	GE, GC-1, BC-1	D	
	Sideritiflavone	43	105	C ₁₈ H ₁₆ O ₈	24.428	206, 273	361.0918	361.0890	-7.85	361 (100), 147 (58), 187 (20), 321 (14)	BE	D	
<i>O</i> -Methylated isoflavonoid	2,5,7-Trihydroxy-3-(4-methoxyphenyl)-3,4-	73	165	C ₁₆ H ₁₄ O ₆	38.075	206, 258, 356	303.0863	303.0833	-9.95	91 (100), 120 (69), 175 (51), 153 (21), 221 (16)	BC-6, BM-6	D	

Proanthocyanidin	dihydro-2 <i>H</i> -1-benzopyran-4-one Procyanidin B3	38	89	C ₃₀ H ₂₆ O ₁₂	22.477	204, 213sh, 282	579.1497	579.1566	11.91	123 (100), 233 (33), 259 (26), 151 (22)	C-1, GC-1, BC-1, GM-1, C-3, GC-3, BC-3, GM-3, GM-6, GC-9, GM-9	D, L	(17)
	Procyanidin B1	49	123	C ₃₀ H ₂₆ O ₁₂	26.435	204, 282	579.1497	579.1566	11.91	123 (100), 233 (33), 259 (26), 151 (22)	GE, BE, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	S, D, L	(17)
	Procyanidin B2	50	124	C ₃₀ H ₂₆ O ₁₂	27.127	204, 325	579.1497	579.1558	10.53	139 (100), 123 (52), 283 (28), 275 (28)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6	S, D, L	(17)
	Procyanidin B1 3- <i>O</i> -gallate	59	142	C ₃₇ H ₃₀ O ₁₆	31.447	208, 282, 310	731.1607	731.1704	13.27	127 (100), 139 (46), 275 (44)	GE, BE, C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, C-6, GC-6, BC-6, GM-6, BM-6	D	
	Procyanidin B2 3- <i>O</i> -gallate	61	147	C ₃₇ H ₃₀ O ₁₆	32.412	204, 280	731.1607	731.1704	13.27	127 (100), 139 (46), 275 (44)	GC-1, BC-1,	D	

	Procyanidin dimer-trimethylglyoxal adduct 1	71	163	C ₃₉ H ₃₈ O ₁₈	37.156	208, 268, 350	795.2131	795.2084	-5.91	732 (100)	GM-1, BC-3, GM-3, GC-6, BE, BC-1	D
	Procyanidin dimer-trimethylglyoxal adduct 2	73	166	C ₃₉ H ₃₈ O ₁₈	38.075	206, 258, 355	795.2131	795.2098	-4.15	493 (100), 389 (51)	BE, BC-1, BM-1	D
	Procyanidin dimer-trimethylglyoxal adduct 3	74	169	C ₃₉ H ₃₈ O ₁₈	38.451	208, 258, 350	795.2131	795.2040	-11.44	582 (100)	GC-1, GM-1	D
Others: phenolic compounds												
Benzenediol	2-(3,4-Dihydroxyphenyl)-3,5,7-trihydroxy-4a,8a-dihydro-4 <i>H</i> -chromen-4-one	79	183	C ₁₅ H ₁₂ O ₇	41.522	205, 290	305.0656	305.0667	3.67	123 (100), 111 (66), 141 (33)	C-1, C-6	D
Benzoyl derivative	1-Phenylprop-2-yn-1-one	38	88	C ₉ H ₆ O	22.477	204, 213sh, 282	131.0491	131.0468	-17.79	104 (100), 78 (56)	4GM-9	D
Coumarin	Coumarin	42	102	C ₉ H ₆ O ₂	24.075	248, 340	147.0441	147.0442	0.68	65 (100), 105 (30)	GE, BE, GC-1, BC-1, GM-6, BM-6, GM-9, BM-9	D
	Trichoderamide A	78	181	C ₂₀ H ₂₀ N ₂ O ₉	41.340	208, 282	433.1242	433.1201	-9.37	281 (100)	GM-1, GM-9	D
Coumarin glycoside	3,4,5-Trihydroxy-6-[(8-hydroxy-2-oxo-2 <i>H</i> -chromen-7-yl)oxy]oxane-2-carboxylic acid	14	31	C ₁₅ H ₁₄ O ₁₀	13.046	204, 222, 300	355.0660	355.0678	5.17	355 (100), 185 (9)	GM-1	D
Hydrolyzable tannin	Unknown hydrolyzable tannin (gomisin R or gomisin type)	79	184	C ₂₂ H ₂₄ O ₇	41.522	204, 268, 354	401.1595	401.1573	-5.53	401 (100)	BE	D
Hydroxycoumarin	Skimmetin 1	32	71	C ₉ H ₆ O ₃	20.509	206, 276	163.0390	163.0390	0.00	77 (100)	GC-1, GM-1, GM-9	D
	Skimmetin 2	46	114	C ₉ H ₆ O ₃	25.502	194, 222, 287, 328	163.0390	163.0395	3.07	77 (100), 107 (16)	C-1, GM-1, GM-3, GM-6, GM-9	D
	Skimmetin 3	121	260	C ₉ H ₆ O ₃	61.100	224, 328	163.0390	163.0381	-5.52	77 (100)	C-1, BC-1, C-3	D
Naphthol	Solistatinol	129	272	C ₁₈ H ₂₀ O ₄	67.839	200, 232, 278	301.1434	301.1421	-4.37	301 (100), 116 (64)	C-1, GC-1, BC-1, GM-1, BM-1	D
Phenol ether	Unknown phenol ether (4-[3-(4-amidinophenoxy)propoxy]benzamidine or other	62	148	C ₂₁ H ₃₂ N ₄ O ₁₀ S ₂	33.214	202, 236, 272sh, 336	565.1633	565.1626	-1.17	529 (100), 499 (59), 391 (50), 475 (30)	BC-1, GM-1, BM-1, GC-3,	D

phenol ether type)												BC-3, GM-3, BM-3, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9	
Phenolic glycoside	Platensimycin B4 methyl ester	35	79	C ₃₁ H ₃₉ NO ₁₂	21.850	198, 218sh, 257, 331	618.2545	618.2560	2.42	193 (100), 146 (95), 208 (74), 369 (37), 471 (15)	C-3, GC-3, BC-3	D	
Pyranoxanthone	Unknown pyranoxanthone	102	224	C ₂₃ H ₂₂ O ₉ S	51.259	206, 236, 316	475.1057	475.1050	-1.53	313 (100)	GC-1, GM-1	D	
Others: non-phenolic compounds													
Alkaloid	Pilosine	91	206	C ₁₆ H ₁₈ N ₂ O ₃	46.860	208, 268, 349	287.1390	287.1424	11.88	165 (100)	GC-3, BC-3, GM-3, BC-6	D	
Alkyl-phenylketone	<i>O</i> -Acetylaniline	8	19	C ₈ H ₉ NO	10.415	196, 226sh, 275	136.0757	136.0752	-3.53	92 (100), 107 (96)	C-1, GC-1, BC-1, BC-6	D	
Alpha amino acid	Pyroglutamic acid	4	9	C ₅ H ₇ NO ₃	8.277	196, 262	130.0499	130.0501	1.61	84 (100)	C-1, GC-1, BC-1, GM-1, BM-1, C-3, GC-3, BC-3, GM-3, BM-3, GC-6, BC-6, GM-6, BM-6, C-9, GC-9, BC-9, GM-9, BM-9	D	
	Dichotocejpipin B	35	78	C ₁₅ H ₁₆ N ₂ O ₅	21.850	296	305.1132	305.1142	3.28	84 (100), 86 (34), 139 (25), 215 (17)	GM-1	D	
	2-Amino-6-octanamidoheptanoic acid 1	103	227	C ₁₄ H ₂₈ N ₂ O ₃	52.494	195, 280, 330	273.2173	273.2196	8.53	87 (100)	GC-1, BC-1, C-3, GC-3, BC-3, C-6, GC-6, BC-6, C-9, GC-9	D	
	2-Amino-6-octanamidoheptanoic acid 2	105	230	C ₁₄ H ₂₈ N ₂ O ₃	54.283	214, 330	273.2173	273.2190	6.34	86 (100), 160 (23)	GC-1, BC-1, C-3, C-6, C-9, BC-9	D	
Alpha-hydrogen aldehyde	4-Acetamidobutanal	13	28	C ₆ H ₁₁ NO ₂	12.689	208, 327	130.0863	130.0841	-16.95	85 (100), 115 (9)	BC-9, GM-9, BM-9	D	

Aminobenzenesulfonamide	Sulfisoxazole	8	18	C ₁₁ H ₁₃ N ₃ O ₃ S	10.415	196, 226sh, 275	268.0750	268.0737	-4.88	136 (100)	GM-9	D
Aminobenzoic acid	2-Aminobenzoic acid	3	5	C ₇ H ₇ NO ₂	7.270	200, 248, 280sh	138.0550	138.0547	-1.85	138 (100), 94 (36)	C-1, C-6, BC-6, C-9, BC-9	D
Aryl alkyl ketone	1-(Pyrazin-2-yl)ethan-1-one	7	16	C ₆ H ₆ N ₂ O	9.797	192, 205sh, 262	123.0553	123.0561	6.35	81 (100), 108 (33)	GM-1, GM-6, GM-9	D
Biguanide	Buformin 1	20	42	C ₆ H ₁₅ N ₅	14.758	198, 266	158.1400	158.1433	20.73	72 (100), 84 (38), 130 (17), 102 (12)	GC-3	D
Biguanide	1-Carbamimidamido-N,N-dimethylmethanimidamide	22	48	C ₄ H ₁₁ N ₅	15.965	264, 338	130.1087	130.1051	-28.07	84 (100)	C-3, GC-3, BC-3, GM-3, BM-3	D
Biguanide	Buformin 2	47	120	C ₆ H ₁₅ N ₅	25.832	223, 246, 329	158.1400	158.1439	24.27	84 (100)	BC-3, GM-3	D
Butenolide	Dihydrotetrodecamycin	4	8	C ₁₈ H ₂₄ O ₆	8.277	196, 267	337.1646	337.1673	8.11	253 (100), 301 (36), 111 (20)	BM-1, BC-6, BC-9	D
Cyclic peptide	Patellamide C 1	41	99	C ₃₇ H ₄₆ N ₈ O ₆ S ₂	23.825	211, 280	763.3054	763.3088	4.39	163 (100), 275 (32), 151 (19), 257 (15)	BM-3	D
	Patellamide C 2	47	115	C ₃₇ H ₄₆ N ₈ O ₆ S ₂	25.832	216, 280	763.3054	763.3067	1.64	287 (100), 306 (42), 595 (15)	GC-3, GM-3	D
Cycloheptathiophene	Ketotifen	15	33	C ₁₉ H ₁₉ NOS	13.587	194, 282	310.1260	310.1232	-9.16	84 (100), 70 (57)	BC-9, BM-9	D
Dihydrofuran	Phaeofuran B 1	37	86	C ₁₂ H ₁₈ O ₅	22.193	194, 280, 330	243.1227	243.1262	14.23	86 (100), 84 (29)	C-3, GC-3, BC-3	D
	Phaeofuran B 2	39	93	C ₁₂ H ₁₈ O ₅	22.905	200, 284	243.1227	243.1256	12.05	120 (100)	C-3, GC-3	D
Dipeptide	gamma-Glutamylleucine	10	23	C ₁₁ H ₂₀ N ₂ O ₅	11.194	194, 280	261.1445	261.1474	11.23	130 (100)	C-1, BC-1, C-3	D
Dipeptide	2-(2-Amino-4-methylpentanamido)-3-methylbutanoic acid 1	23	49	C ₁₁ H ₂₂ N ₂ O ₃	16.276	195, 205sh, 280, 320	231.1703	231.1666	-15.91	69 (100), 87 (77)	C-9, GC-9, BC-9	D
	2-(2-Amino-4-methylpentanamido)-3-methylbutanoic acid 2	26	59	C ₁₁ H ₂₂ N ₂ O ₃	17.301	195, 205sh, 280, 320	231.1703	231.1674	-12.50	86 (100)	C-9, GC-9, BC-9	D
	2-(2-Amino-4-methylpentanamido)-3-methylbutanoic acid 3	33	72	C ₁₁ H ₂₂ N ₂ O ₃	20.888	192, 220, 280, 332	231.1703	231.1709	2.64	73 (100)	C-1, BC-1, C-3, BC-3, C-6, BC-6, C-9, GC-9, BC-9	D
	2-[(2-Amino-1-hydroxy-3-methylpentylidene)amino]-4-methylpentanoic acid 1	39	91	C ₁₂ H ₂₄ N ₂ O ₃	22.905	196, 220, 282	245.1860	245.1816	-17.70	86 (100), 69 (10)	C-1, BC-1, C-3, GM-3, GC-9, BC-9	D
	1-(2-Amino-3-phenylpropanoyl)pyrrolidine-2-carboxylic acid	40	95	C ₁₄ H ₁₈ N ₂ O ₃	23.385	216, 260, 293sh	263.1390	263.1343	-17.93	70 (100)	BC-9	D
	2-[(2-Amino-1-hydroxy-3-methylpentylidene)amino]-4-methylpentanoic acid	47	119	C ₁₂ H ₂₄ N ₂ O ₃	25.832	192, 214, 320	245.1860	245.1869	3.63	87 (100)	C-1, BC-1, C-3, GM-3, C-9, BC-9	D

	acid 2											
	2-(2-Amino-4-methylpentanamido)-3-methylbutanoic acid 4	50	126	C ₁₁ H ₂₂ N ₂ O ₃	27.127	204, 325	231.1703	231.1705	0.61	86 (100)	C-1, BC-1, C-6, C-9	D
	2-[(2-Amino-1-hydroxy-3-methylpentylidene)amino]-4-methylpentanoic acid 3	53	130	C ₁₂ H ₂₄ N ₂ O ₃	27.791	220, 242, 328	245.1860	245.1866	2.37	86 (100)	C-1, BC-1, C-3, BC-3	D
	2-[(2-Amino-1-hydroxy-3-methylpentylidene)amino]-4-methylpentanoic acid 4	63	151	C ₁₂ H ₂₄ N ₂ O ₃	33.413	208, 270, 338	245.1860	245.1808	-21.12	72 (100), 146 (12)	C-9	D
Diphenylmethane	Darifenacin	123	264	C ₂₈ H ₃₀ N ₂ O ₂	63.570	209, 270, 370	427.2380	427.2379	-0.25	427 (100)	GM-1	D
Enoate ester	Stagonolide H	117	252	C ₁₀ H ₁₂ O ₄	59.227	206, 316	197.0808	197.0784	-12.36	99 (100), 83 (80)	C-1, GC-1, BC-1, C-3, BC-3	D
Fatty acid ester	Phomester C	78	180	C ₁₉ H ₃₅ NO ₅	41.340	208, 282	358.2588	358.2635	13.12	86 (100)	C-9	D
Gamma butyrolactone	N,N-Dimethyl-5-oxoxolane-2-carboxamide	36	82	C ₇ H ₁₁ NO ₃	22.002	220, 282	158.0812	158.0782	-18.79	85 (100)	C-9, GC-9, BC-9	D
Glutamine	Theanine	3	6	C ₇ H ₁₄ N ₂ O ₃	7.270	200, 248, 280sh	175.1077	175.1047	-17.47	84 (100), 158 (49), 130 (8)	GC-9, BC-9, GM-9, BM-9	D
Guanidine	3-Guanidino-N-(2-hydroxyheptyl)propionamide	50	125	C ₁₁ H ₂₄ N ₄ O ₂	27.127	238, 330	245.1972	245.1956	-6.54	91 (100), 87 (76)	C-6, BC-6, C-9, GC-9, BC-9	D
Indole	<i>trans</i> -3-Indoleacrylic acid	39	92	C ₁₁ H ₉ NO ₂	22.905	196, 220, 282	188.0706	188.0707	0.53	144 (100), 119 (34), 147 (34), 91 (21)	C-1, GC-1, BC-1, GM-1, BM-1, GM-6, BM-6, GC-9, GM-9, BM-9	D
Indoline	2,3-Dihydro-1 <i>H</i> -indole 1	20	41	C ₈ H ₉ N	14.758	236, 296	120.0808	120.0879	59.13	77 (100)	C-6, GC-6, BC-6, C-9, GC-9, BC-9, GM-9, BM-9	D
	2,3-Dihydro-1 <i>H</i> -indole 2	21	45	C ₈ H ₉ N	15.458	194, 212sh, 290	120.0808	120.0806	-1.47	77 (100), 103 (13)	C-1, GC-1, BC-1, GM-1, BM-1	D
	2,3-Dihydro-1 <i>H</i> -indole 3	60	144	C ₈ H ₉ N	32.078	194, 266	120.0808	120.0810	1.67	93 (100), 104 (89)	C-1, BC-1, C-9	D

Indolizidine	Isoelaecarpiline	7	17	C ₁₆ H ₂₁ NO ₂	9.797	200, 264	260.1645	260.1636	-3.48	86 (100), 132 (9)	C-1, GC-1	D	
Isoleucine	<i>L</i> -Isoleucine 1	5	10	C ₆ H ₁₃ NO ₂	8.902	198, 270	132.1019	132.0996	-17.30	70 (100)	GC-9, BC-9		
	<i>L</i> -Isoleucine 2	6	12	C ₆ H ₁₃ NO ₂	9.325	198, 270	132.1019	132.1000	-14.34	91 (100)	GC-9, BC-9		
Macrolactam	Cespitulactam D	33	73	C ₂₀ H ₂₉ NO ₂	20.888	194, 365	316.2271	316.2254	-5.49	187 (100)	C-1, GC-1, BC-1	D	
Morpholine	4-Amino-2-[(morpholin-4-yl)methyl]oxolan-3-ol	22	46	C ₉ H ₁₈ N ₂ O ₃	15.965	264, 338	203.1390	203.1357	-16.34	86 (100), 132 (26)	C-9, GC-9, BC-9	D	
Morpholine	Ciclosidomine	25	53	C ₁₃ H ₂₀ N ₄ O ₃	17.180	198, 268	281.1608	281.1576	-11.48	84 (100)	C-6, GC-6, BC-6, C-9	D	
Naphthopyran	Inumakilactone A	40	96	C ₁₈ H ₂₀ O ₈	23.385	216, 260, 293sh	365.1231	365.1244	3.56	334 (100)	GE, BE, GC-1, BC-1, GM-1, BM-1	D	
Oxasteroid	Wortmannin	92	208	C ₂₃ H ₂₄ O ₈	47.594	194, 210, 280	429.1544	429.1568	5.61	235 (100), 181 (44)	GC-1	D	
Peptide	Methyl 2-(2-amino-4-methylpentanamido)-3-phenylpropanoate	100	221	C ₁₆ H ₂₄ N ₂ O ₃	50.678	224, 280	293.1860	293.1865	1.81	103 (100), 147 (87)	C-1, GC-1, C-6, C-9	D	
Physalin	Physalin D	84	195	C ₂₈ H ₃₂ O ₁₁	44.325	202, 218, 266, 302, 348	545.2017	545.2050	6.05	375 (100)	GC-1, BC-1, GM-1	D	
Proline	Proline betaine	9	21	C ₇ H ₁₃ NO ₂	10.823	194, 280	144.1019	144.0998	-14.54	71 (100)	C-9, GC-9, BC-9	D	
Purine alkaloid	Theophyllin	26	56	C ₇ H ₈ N ₄ O ₂	17.301	206, 275	181.0720	181.0759	21.36	138 (100), 181 (76), 110 (56), 163 (36), 135 (23), 139 (22), 108 (21)	BE	D, L	(1)
	Theobromine	31	69	C ₇ H ₈ N ₄ O ₂	20.153	204, 274	181.0720	181.0724	2.21	138 (100), 110 (30)	GE, BE, GC-1, BC-1, GM-1, BM-1, GC-6, BC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9	D	
	Caffeine	57	136	C ₈ H ₁₀ N ₄ O ₂	29.773	216, 276	195.0877	195.0898	10.76	138 (100), 139 (9)	GE, BE, GC-1, BC-1, GM-1, BM-1, GC-3, BC-3, GM-3, BM-3, GC-6,	S, D, L	(6), (9)

Purine nucleoside	Adenosine	11	25	C ₁₀ H ₁₃ N ₅ O ₄	11.528	210, 260	268.1040	268.1081	15.26	136 (100)	BC-6, GM-6, BM-6, GC-9, BC-9, GM-9, BM-9, GM-1, GM-6	D
Purine nucleoside	2-(<i>9</i> -[3,4-Dihydroxy-5-(hydroxymethyl)oxolan-2-yl]-9 <i>H</i> -purin-6-yl}amino)butanedioic acid	26	58	C ₁₄ H ₁₇ N ₅ O ₈	17.301	202, 212, 270, 329	384.1150	384.1165	3.91	253 (100), 234 (22)	C-1, GC-1, BC-1, C-3, BC-6, C-9, GC-9, BC-9	D
Pyrrroloindole	Physostigmine	26	55	C ₁₅ H ₂₁ N ₃ O ₂	17.301	206, 275	276.1707	276.1667	-14.48	231 (100), 260 (96), 175 (83), 91 (47), 112 (46)	C-9, GC-9, BC-9	D
Sesquiterpenoid	Debromolaurinterol	12	27	C ₁₅ H ₂₀ O	12.358	212, 294, 328	217.1587	217.1613	11.87	165 (100), 86 (89), 147 (75)	C-6, GC-6, BC-6, C-9, GC-9, BC-9	D
Sugar alcohol	Ribitol	30	67	C ₅ H ₁₂ O ₅	20.024	216, 277	153.0758	153.0796	25.15	153 (100), 80 (14), 123 (12)	GC-3, BC-3, GM-3	D
Terpene glycoside	Monomelittoside	14	32	C ₁₅ H ₂₂ O ₁₀	13.046	206, 220sh, 302	363.1286	363.1293	2.00	231 (100), 156 (26)	GC-1, BC-1	D
Tricarboxylic acid	Cryptoporic acid B	110	237	C ₂₃ H ₃₆ O ₈	56.333	203, 218sh, 278	441.2483	441.2505	4.99	139 (100)	GC-1	D
Triterpenoid	Gypsogenic acid	74	170	C ₃₀ H ₄₆ O ₅	38.451	214, 270, 358	487.3418	487.3372	-9.40	487 (100)	GM-3	D
Xanthanolide	Xanthanol	66	156	C ₁₇ H ₂₄ O ₅	35.542	334	309.1697	309.1744	15.36	136 (100), 72 (77), 210 (56), 193 (54)	C-9	D

^a Wine type: GE, BE = detected in green tea ethanolic extract and black tea ethanolic extract, respectively; C-1, C-3, C-6 etc. refer to Table 1.

^b ID: identification method by S = standard, D = database and L = literature.

^c Literature reported: (1) = Akhtar et al.¹, (2) = Bastos et al.², (3) = Chen et al.³, (4) = Dai et al.⁴, (5) = Dai et al.⁵, (6) = Del Rio et al.⁶, (7) = Fracassetti et al.⁷, (8) = Itoh et al.⁸, (9) = Jiang et al.⁹, (10) = Lantzouraki et al.¹⁰, (11) = Lin et al.¹¹, (12) = Meng et al.¹², (13) = Qi¹³, (14) = Rossetti et al.¹⁴, (15) = Sang and Yang¹⁵, (16) = Scoparo et al.¹⁶, (17) = Topi et al.¹⁷, (18) = Wang et al.¹⁸, (19) = Wang et al.¹⁹, (20) = Xie et al.²⁰, (21) = Zhang et al.²¹, (22) = Zhang et al.²², (23) = Zhu et al.²³.

* = Semi-quantification: Quinic acid, Procyanidin B1 3-*O*-gallate, Quercetin 3-*O*-di-galactoside/glucoside, Myricetin 3-*O*-galactoside 1, Kaempferol 3-*O*-rutoside 2, Unknown quercetin conjugate and Unknown kaempferol conjugate were semi-quantified using the calibration curves of Chlorogenic acid, Procyanidin B1, Quercetin 3-*O*-galactoside, Quercetin 3-*O*-glucoside, Kaempferol 3-*O*-glucoside, Quercetin and Kaempferol, respectively.

ND = Not detected or value below the detection threshold.

C, 2GC, 4GC, 1BC, 3BC = Control Chardonnay wine, 2% (w/v) green tea-macerated Chardonnay wine, 4% (w/v) green tea-macerated Chardonnay wine, 1% (w/v) black tea-macerated Chardonnay wine, 3% (w/v) black tea-macerated Chardonnay wine.

tr = trace.

-- = Not available.

^{abcd} = Values with different superscripts in the same wine group are significantly different ($p < 0.05$).

Table S3 HPLC-DAD quantification of phenolic compounds in tea-macerated model wines at different aging times (µg/mL)

Compounds	M				2GM				4GM				1BM				3BM			
	Aging time (month)																			
	1	3	6	9	1	3	6	9	1	3	6	9	1	3	6	9	1	3	6	9
<i>Phenolic acids</i>																				
Quinic acid*	ND	ND	ND	ND	tr	11.14 ± 1.08 ^a	177.59 ± 9.21 ^a	42.05 ± 2.32 ^a	1.56 ± 0.90 ^a	29.39 ± 10.01 ^b	319.58 ± 10.68 ^a	67.65 ± 38.03 ^b	16.26 ± 0.41 ^c	22.29 ± 0.59 ^b	54.61 ± 13.93 ^a	55.05 ± 10.45 ^a	76.32 ± 31.63 ^a	89.46 ± 35.19 ^a	122.30 ± 10.74 ^a	174.28 ± 26.64 ^a
Galic acid	ND	ND	ND	ND	110.20 ± 2.96 ^b	124.48 ± 8.10 ^b	tr	204.31 ± 19.13 ^a	213.38 ± 66.35 ^b	218.84 ± 50.20 ^b	tr	401.94 ± 161.35 ^a	6.10 ± 0.26 ^d	10.18 ± 0.46 ^c	23.23 ± 0.00 ^b	95.98 ± 106.13 ^a	48.26 ± 2.69 ^c	45.46 ± 22.46 ^c	64.81 ± 11.32 ^a	70.58 ± 32.65 ^a
Chlorogenic acid	ND	ND	ND	ND	24.01 ± 1.26 ^a	25.44 ± 2.13 ^a	22.77 ± 1.68 ^a	24.94 ± 17.31 ^a	51.31 ± 3.90 ^a	46.52 ± 10.33 ^a	47.76 ± 3.40 ^a	49.64 ± 23.29 ^a	0.32 ± 0.22 ^b	tr	1.61 ± 1.77 ^b	11.46 ± 19.44 ^a	5.46 ± 4.69 ^a	5.78 ± 2.93 ^a	3.37 ± 0.19 ^a	4.55 ± 3.95 ^a
Total phenolic acids	--	--	--	--	134.21 ± 4.22^c	161.05 ± 11.31^c	200.36 ± 10.89^b	271.30 ± 25.35^a	266.25 ± 84.55^c	294.76 ± 70.54^c	367.35 ± 14.07^b	519.23 ± 222.67^a	22.67 ± 0.89^c	32.47 ± 1.05^c	79.45 ± 0.00^b	162.49 ± 136.03^a	130.04 ± 39.02^a	140.69 ± 60.58^a	190.47 ± 22.25^a	249.41 ± 63.24^a
<i>Flavonoids</i>																				
Gallic acid	ND	ND	ND	ND	579.20 ± 23.48 ^d	780.93 ± 70.79 ^c	1001.59 ± 49.73 ^b	1836.30 ± 885.25 ^a	1070.10 ± 285.55 ^c	1349.94 ± 242.15 ^c	1891.93 ± 114.97 ^b	2364.61 ± 914.95 ^a	197.05 ± 4.27 ^d	246.22 ± 4.61 ^c	409.97 ± 39.15 ^b	833.81 ± 539.86 ^a	690.58 ± 318.56 ^b	785.98 ± 270.94 ^b	845.87 ± 60.22 ^b	965.80 ± 420.74 ^a
Epigallocatechin	ND	ND	ND	ND	626.08 ± 68.43 ^c	815.40 ± 48.55 ^b	1029.22 ± 52.71 ^a	895.27 ± 467.55 ^b	1187.07 ± 364.01 ^c	1383.97 ± 299.14 ^c	1795.85 ± 91.07 ^b	2095.17 ± 793.68 ^a	27.83 ± 5.47 ^d	43.63 ± 5.92 ^c	63.67 ± 17.07 ^b	374.53 ± 561.17 ^a	122.06 ± 62.42 ^c	165.58 ± 20.27 ^b	188.26 ± 13.91 ^a	109.08 ± 53.36 ^a
Procyanidin B1	ND	ND	ND	ND	26.85 ± 1.64 ^b	33.93 ± 1.23 ^a	27.88 ± 0.68 ^b	35.39 ± 0.97 ^a	42.23 ± 9.42 ^b	47.33 ± 5.37 ^b	42.73 ± 3.24 ^b	55.83 ± 18.42 ^a	8.27 ± 0.68 ^c	10.21 ± 0.83 ^b	10.38 ± 2.74 ^b	17.21 ± 14.17 ^a	21.52 ± 6.89 ^a	20.07 ± 3.64 ^a	20.39 ± 3.48 ^a	16.62 ± 8.73 ^a
Catechin	ND	ND	ND	ND	3165.80 ± 47.32 ^c	5592.70 ± 404.97 ^b	4952.16 ± 191.60 ^b	8335.28 ± 4249.99 ^a	5576.38 ± 1405.23 ^c	9405.58 ± 1430.41 ^c	8247.79 ± 485.22 ^b	9461.17 ± 3095.64 ^a	1175.54 ± 42.85 ^c	1438.55 ± 52.53 ^c	2207.71 ± 198.34 ^b	3751.95 ± 1811.69 ^a	3396.73 ± 1203.18 ^b	3890.03 ± 1056.94 ^b	4582.78 ± 289.20 ^b	4841.91 ± 2063.66 ^a
Procyanidin B2	ND	ND	ND	ND	99.38 ± 9.42 ^b	19.30 ± 1.63 ^a	140.25 ± 6.36 ^a	116.78 ± 45.06 ^b	182.94 ± 54.55 ^b	33.25 ± 5.85 ^c	252.13 ± 86.07 ^b	243.58 ± 0.44 ^b	3.22 ± 0.34 ^d	2.22 ± 0.34 ^d	2.91 ± 3.72 ^b	54.00 ± 78.50 ^a	19.10 ± 9.80 ^a	8.28 ± 2.43 ^b	11.93 ± 1.78 ^a	21.34 ± 11.87 ^a
Epicatechin	ND	ND	ND	ND	tr	24.31 ± 1.77 ^b	52.69 ± 5.82 ^a	45.19 ± 2.98 ^b	46.18 ± 33.96 ^b	56.20 ± 14.19 ^b	137.69 ± 35.37 ^b	102.05 ± 51.65 ^a	tr	tr	4.84 ± 0.00 ^b	68.99 ± 0.00 ^a	35.32 ± 20.95 ^a	14.02 ± 6.57 ^b	10.61 ± 3.98 ^b	15.09 ± 6.86 ^b
Epigallocatechin gallate	ND	ND	ND	ND	12.93 ± 0.43 ^d	22.07 ± 0.38 ^c	68.94 ± 5.02 ^b	137.73 ± 188.74 ^a	26.46 ± 11.05 ^c	35.12 ± 37.74 ^b	157.53 ± 19.51 ^b	64.75 ± 37.74 ^b	5.82 ± 0.46 ^d	8.71 ± 3.73 ^b	16.71 ± 20.83 ^a	33.60 ± 31.68 ^a	24.51 ± 3.43 ^b	28.40 ± 3.43 ^b	45.34 ± 11.64 ^b	72.61 ± 25.27 ^a
Procyanidin B1 3-O-gallate*	ND	ND	ND	ND	1.13 ± 0.37 ^c	18.63 ± 3.95 ^b	18.04 ± 5.89 ^b	79.13 ± 24.58 ^a	74.51 ± 54.82 ^b	56.44 ± 16.68 ^b	128.50 ± 11.51 ^b	119.71 ± 88.43 ^a	3.44 ± 2.21 ^c	2.61 ± 2.87 ^b	8.80 ± 81.97 ^a	48.65 ± 18.22 ^b	40.45 ± 4.52 ^b	24.95 ± 9.18 ^b	39.83 ± 9.18 ^b	64.16 ± 23.06 ^a
Gallic acid	ND	ND	ND	ND	365.63 ± 96.87 ^b	500.54 ± 56.93 ^b	618.32 ± 365.84 ^a	460.22 ± 166.06 ^b	745.65 ± 93.86 ^b	828.61 ± 172.21 ^a	1013.78 ± 294.80 ^b	986.09 ± 0.85 ^a	26.17 ± 4.68 ^c	35.95 ± 48.55 ^b	267.88 ± 5.14 ^b	66.02 ± 25.99 ^b	79.14 ± 18.02 ^b	60.10 ± 14.91 ^b	55.32 ± 22.62 ^b	
Quercetin 3-O-di-galactoside/glucoside*	ND	ND	ND	ND	6.25 ± 1.19 ^b	0.85 ± 0.00 ^c	0.85 ± 0.00 ^c	15.20 ± 5.20 ^a	15.96 ± 5.54 ^b	0.85 ± 0.00 ^c	23.41 ± 7.18 ^b	32.36 ± 13.53 ^a	tr	0.82 ± 0.83 ^c	2.76 ± 2.12 ^b	7.73 ± 12.81 ^a	4.18 ± 0.52 ^b	0.34 ± 0.00 ^c	1.30 ± 1.65 ^b	2.84 ± 2.23 ^b
Myricetin 3-O-galactoside 1*	ND	ND	ND	ND	tr	17.56 ± 2.38 ^a	tr	9.28 ± 0.00 ^c	tr	34.07 ± 6.09 ^b	tr	11.63 ± 0.00 ^c	tr	1.41 ± 0.33 ^d	tr	1.55 ± 0.31 ^d	tr	8.66 ± 2.82 ^a	1.65 ± 1.65 ^b	5.68 ± 4.10 ^a
Epicatechin gallate	ND	ND	ND	ND	24.01 ± 0.92 ^b	26.24 ± 1.95 ^b	6.57 ± 0.00 ^c	40.43 ± 3.50 ^b	42.62 ± 12.91 ^b	42.89 ± 7.33 ^b	59.92 ± 6.82 ^b	58.00 ± 19.17 ^a	7.43 ± 0.22 ^d	8.13 ± 0.73 ^b	4.46 ± 3.17 ^b	20.35 ± 16.22 ^a	30.38 ± 12.14 ^a	25.06 ± 7.61 ^b	18.95 ± 5.06 ^b	23.45 ± 11.42 ^b
Rutin	ND	ND	ND	ND	3.72 ± 0.50 ^d	6.88 ± 1.03 ^c	16.54 ± 1.31 ^a	11.32 ± 3.09 ^b	10.50 ± 4.62 ^b	14.33 ± 3.07 ^b	43.35 ± 3.32 ^a	16.36 ± 7.07 ^b	1.57 ± 0.19 ^d	2.09 ± 0.25 ^c	6.30 ± 0.49 ^a	6.84 ± 3.81 ^b	9.13 ± 2.39 ^b	9.17 ± 2.26 ^b	7.77 ± 4.23 ^b	
Quercetin 3-O-galactoside	ND	ND	ND	ND	3.09 ± 0.00 ^b	3.09 ± 0.61 ^b	0.85 ± ±0.00 ^c	6.47 ± 3.52 ^a	6.47 ± 2.35 ^b	9.06 ± 1.33 ^b	5.67 ± 0.96 ^c	8.74 ± 1.86 ^b	tr	0.34 ± 0.00 ^c	5.97 ± 0.16 ^b	2.28 ± 3.36 ^b	tr	3.58 ± 1.03 ^b	1.60 ± 0.66 ^b	0.34 ± 0.00 ^c
Kaempferol 3-O-rutinoside 2*	ND	ND	ND	ND	5.16 ± 0.57 ^b	7.39 ± 0.87 ^b	0.07 ± 0.00 ^c	14.38 ± 2.34 ^a	13.35 ± 4.86 ^b	18.26 ± 3.55 ^b	17.78 ± 3.70 ^b	17.78 ± 4.68 ^b	2.31 ± 0.09 ^d	2.79 ± 0.07 ^c	0.03 ± 0.00 ^c	3.58 ± 4.38 ^a	11.07 ± 4.38 ^a	10.81 ± 3.20 ^a	0.03 ± 0.00 ^c	10.73 ± 5.11 ^a
Kaempferol 3-O-glucoside	ND	ND	ND	ND	tr	2.51 ± 0.01 ^b	7.95 ± 0.56 ^a	7.64 ± 6.03 ^a	5.47 ± 0.73 ^b	4.61 ± 1.05 ^c	25.83 ± 4.62 ^a	6.53 ± 2.26 ^b	1.52 ± 0.05 ^d	1.86 ± 0.16 ^b	3.81 ± 0.25 ^a	3.13 ± 0.28 ^a	5.89 ± 2.19 ^b	7.11 ± 1.37 ^b	10.02 ± 1.34 ^a	5.79 ± 2.51 ^b
Unknown quercetin conjugate*	ND	ND	ND	ND	5.80 ± 0.20 ^b	6.36 ± 0.09 ^b	3.23 ± 0.00 ^c	6.53 ± 1.08 ^a	8.38 ± 1.45 ^a	8.88 ± 1.26 ^a	3.23 ± 0.00 ^c	9.84 ± 2.39 ^a	tr	1.83 ± 0.09 ^b	1.29 ± 0.00 ^b	3.22 ± 1.83 ^a	2.67 ± 0.80 ^b	4.24 ± 0.55 ^a	1.29 ± 0.00 ^c	3.19 ± 1.28 ^a
Unknown kaempferol conjugate*	ND	ND	ND	ND	3.66 ± 0.06 ^b	4.02 ± 0.11 ^a	5.31 ± 0.11 ^a	4.04 ± 0.35 ^b	4.79 ± 0.52 ^b	5.29 ± 0.56 ^b	7.21 ± 0.49 ^a	5.11 ± 0.87 ^b	tr	1.40 ± 0.12 ^c	1.14 ± 0.00 ^b	1.85 ± 0.64 ^a	2.20 ± 0.70 ^b	2.52 ± 0.13 ^b	2.03 ± 0.30 ^b	2.09 ± 0.98 ^a
Total flavonoids	--	--	--	--	4928.70 ± 251.40^c	7882.71 ± 579.80^b	7950.45 ± 376.72^b	12055.02 ± 6256.07^a	9059.06 ± 2417.61^c	13327.84 ± 2136.18^b	13858.20 ± 985.88^b	15659.30 ± 5433.21^a	1460.18 ± 56.73^d	1808.77 ± 76.39^c	2799.28 ± 278.94^b	5504.58 ± 3542.14^a	4481.82 ± 1701.00^b	5087.95 ± 1447.89^b	5874.55 ± 419.58^b	6223.79 ± 2668.05^a
<i>Purine alkaloid</i>																				
Caffeine	ND	ND	ND	ND	918.66 ± 175.72 ^b	419.07 ± 20.35 ^c	1427.96 ± 183.42 ^a	1126.41 ± 735.27 ^b	1911.92 ± 433.91 ^b	690.71 ± 145.68 ^b	2434.06 ± 464.72 ^a	2552.21 ± 926.54 ^a	30.63 ± 3.57 ^b	24.33 ± 2.80 ^b	33.62 ± 8.61 ^b	68.64 ± 11.43 ^a	114.91 ± 50.03 ^a	75.87 ± 25.07 ^b	65.74 ± 1.84 ^b	91.32 ± 41.49 ^b
Total purine alkaloid	--	--	--	--	918.66 ± 175.72^b	419.07 ± 20.35^c	1427.96 ± 183.42^a	1126.41 ± 735.27^b	1911.92 ± 433.91^b	690.71 ± 145.68^b	2434.06 ± 464.72^a	2552.21 ± 926.54^a	30.63 ± 3.57^b	24.33 ± 2.80^b	33.62 ± 8.61^b	68.64 ± 11.43^a	114.91 ± 50.03^a	75.87 ± 25.07^b	65.74 ± 1.84^b	91.32 ± 41.49^b

* = Semi-quantification: Quinic acid, Procyanidin B1 3-*O*-gallate, Quercetin 3-*O*-di-galactoside/glucoside, Myricetin 3-*O*-galactoside 1, Kaempferol 3-*O*-rutoside 2, Unknown quercetin conjugate and Unknown kaempferol conjugate were semi-quantified using the calibration curves of Chlorogenic acid, Procyanidin B1, Quercetin 3-*O*-galactoside, Quercetin 3-*O*-galactoside, Kaempferol 3-*O*-glucoside, Quercetin and Kaempferol, respectively.

ND = Not detected or value below the detection threshold.

M, 2GM, 4GM, 1BM, 3BM = model wine, 2% (w/v) green tea-macerated model wine, 4% (w/v) green tea-macerated model wine, 1% (w/v) black tea-macerated model wine, 3% (w/v) black tea-macerated model wine.

tr = trace.

-- = Not available.

^{abcd} = Values with different superscripts in the same wine group are significantly different ($p < 0.05$).

8	Limonene	117 0	ND	ND	ND	ND	0.02 ± 0.00 ^a	0.02 ± 0.00 ^b	0.02 ± 0.00 ^a	0.02 ± 0.00 ^a	0.03 ± 0.00 ^a	0.03 ± 0.00 ^a	0.03 ± 0.00 ^a	0.03 ± 0.00 ^a	0.02 ± 0.00 ^a	0.02 ± 0.00 ^a	0.02 ± 0.00 ^a	0.02 ± 0.00 ^a	0.02 ± 0.00 ^a	0.02 ± 0.00 ^b	0.02 ± 0.00 ^b	RI, MS, ES	Citrus, orange, fresh, sweet	
15	Terpinolene****	125 5	ND	ND	ND	ND	0.01 ± 0.00	tr	tr	tr	0.01 ± 0.00 ^a	tr	0.01 ± 0.00 ^b	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS, ES	Sweet, fresh pine, citrus, woody, lemon peel
	Total terpene		ND	ND	ND	ND	0.03 ± 0.00 ^a	0.02 ± 0.00 ^b	0.02 ± 0.00 ^a	0.02 ± 0.00 ^b	0.04 ± 0.00 ^a	0.04 ± 0.00 ^a	0.04 ± 0.00 ^b	0.03 ± 0.00 ^a	0.02 ± 0.00 ^b	0.02 ± 0.00 ^b	0.02 ± 0.00 ^a	0.02 ± 0.00 ^a	0.02 ± 0.00 ^{ab}	0.02 ± 0.00 ^{ab}	0.02 ± 0.00 ^b			
	Total	160.45 ± 14.40 ^a	166.68 ± 13.12 ^a	166.81 ± 32.22 ^a	118.39 ± 3.91 ^b	143.29 ± 10.77 ^b	160.2 3 ± 8.32 ^b	145.49 ± 8.32 ^b	123.61 ± 16.53 ^c	146.87 ± 12.65 ^b	172.84 ± 14.52 ^a	158.53 ± 17.92 ^{ab}	120.08 ± 13.98 ^c	136.57 ± 5.90 ^b	182.26 ± 43.71 ^a	149.45 ± 12.25 ^b	115.43 ± 6.98 ^c	151.87 ± 8.01 ^b	160.79 ± 11.86 ^a	151.66 ± 19.78 ^b	127.87 ± 16.28 ^c			

*RI = retention index

**ID = Identification by comparison of mass spectra (MS) with in-house NIST library 11.0, MS and retention index (RI) with NIST Webbook (NIST²⁴), and MS and retention time with external standards

*** = Odor description were obtained from TGSC²⁵.

****= Semi-quantification: Ethoxyacetic acid, acetic acid, pterin-6-carboxylic acid, 2-methyl-1-propanol, 2-pentanol, 2-hexanol, 1-(2-aminoethylamino)-2-propanol, 2-octen-4-ol, 3-methyl-1-pentanol, 2-methyl-4-heptanol, 2,3-butanediol, phenylethyl alcohol, 3-hydroxybutanal, myristicin, 3-hydroxymandelic acid ethyl ester, ethyl (*E*)-3-hexenoate, ethyl lactate, propyl hexanoate, ethyl 3-hydroxybutanoate, isoamyl lactate, ethyl 2-furancarboxylate, ethyl 9-decenoate and 3-hydroxy-2-butanone were semi-quantified using the calibration curves of ethyl hexanoate, ethyl octanoate, 2-ethyl-1-hexanol, 3-methylbutyl acetate, 1-butanol, 1-butanol, ethyl hexanoate, 1-hexanol, 1-hexanol, 1-hexanol, 1-octanol, benzyl alcohol, ethyl hexanoate, benzaldehyde, 1-butanol, 1-hexanol, 2-heptanol, 2-heptanol, ethyl nonanoate, 1-octanol, benzeneacetaldehyde, diethyl succinate and ethyl hexanoate, respectively.

Samples of C, 2GC, 4GC etc. refer to Table S2.

ND = Not detected or value below the detection threshold.

tr = trace

-- = Not available.

a-d = Values with different superscripts in the same wine group are significantly different ($p \leq 0.05$)

Table S5 Effect of aging time on volatile compounds of tea-macerated model wines ($\mu\text{g/mL}$)

Peak No.	Compound name	RI*	Aging time (month)																				ID**	Odor description***	
			M				2GM				4GM				1BM				3BM						
			1	3	6	9	1	3	6	9	1	3	6	9	1	3	6	9	1	3	6	9			
17	Ethoxyacetic acid****	1268	0.01 ± 0.00	tr	tr	tr	0.01 ± 0.00 ^a	0.01 ± 0.00 ^{ab}	0.01 ± 0.00 ^b	tr	0.01 ± 0.00 ^a	0.01 ± 0.01 ^{ab}	0.01 ± 0.00 ^b	0.01 ± 0.00 ^b	0.03 ± 0.02 ^a	0.02 ± 0.01 ^b	0.01 ± 0.00 ^c	0.01 ± 0.00 ^c	0.02 ± 0.02 ^a	0.02 ± 0.01 ^a	0.01 ± 0.00 ^b	0.01 ± 0.01 ^b	RI, MS	--	
28	Acetic acid****	1433	ND	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS	Pungent, acidic, cheesy, vinegar
29	Pterin-6-carboxylic acid****	1461	ND	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS	--
49	Hexanoic acid	1829	ND	ND	ND	0.58 ± 0.02	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	0.61 ± 0.01 ^a	tr	0.62 ± 0.01 ^a	0.60 ± 0.00 ^a	RI, MS, ES	Sour, fatty, sweaty, cheesy	
53	Octanoic acid	2046	ND	ND	ND	0.66 ± 0.03	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS, ES	Fatty, waxy, rancid, oily, vegetable, cheesy
54	Nonanoic acid	2269	ND	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS, ES	Waxy, dirty, cheese, cultured, dairy
	Total acid		0.01 ± 0.00	tr	tr	tr	1.25 ± 0.05^a	0.01 ± 0.00^a	0.01 ± 0.00^b	0.01 ± 0.00^b	0.01 ± 0.01^a	0.01 ± 0.01^a	0.01 ± 0.00^a	0.01 ± 0.01^a	0.03 ± 0.02^a	0.03 ± 0.01^a	0.01 ± 0.01^c	0.01 ± 0.01^c	0.64 ± 0.02^a	0.02 ± 0.01^a	0.63 ± 0.02^a	0.61 ± 0.01^a			
3	2-Methyl-1-propanol****	1079	ND	ND	ND	0.01 ± 0.01 ^a	0.01 ± 0.00 ^a	0.01 ± 0.00 ^a	0.01 ± 0.00 ^a	0.02 ± 0.01 ^a	0.01 ± 0.01 ^b	0.01 ± 0.00 ^b	0.01 ± 0.00 ^b	tr	0.01 ± 0.00	tr	tr	0.01 ± 0.00 ^a	0.01 ± 0.00 ^a	0.01 ± 0.00 ^a	0.01 ± 0.00 ^a	RI, MS	Ethereal, winery, cortex		
5	2-Pentanol****	1127	ND	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS	Mild, green, fusel, oil, fermented
6	1-Butanol	1132	ND	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS, ES	Fusel, oil, sweet, balsam, whiskey
7	2-Hexanol****	1142	ND	ND	ND	tr	tr	tr	tr	0.26 ± 0.05 ^b	0.93 ± 0.17 ^a	tr	tr	tr	tr	tr	tr	tr	0.67 ± 0.16	tr	tr	tr	tr	RI, MS	Chemical, winery, fruity, fatty, terpenic, cauliflower
10	3-Methyl-1-butanol	1196	ND	ND	ND	0.01 ± 0.00	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS, ES	Fusel, oily, alcoholic, whiskey, fruity, banana
12	1-(2-Aminoethylamino)-2-propanol****	1228	ND	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS	--
20	2-Octen-4-ol****	1304	ND	ND	ND	0.11 ± 0.02 ^a	0.03 ± 0.01 ^a	0.07 ± 0.02 ^b	0.08 ± 0.03 ^b	0.10 ± 0.02 ^b	0.07 ± 0.05 ^a	0.10 ± 0.07 ^b	0.23 ± 0.09 ^a	0.06 ± 0.00	tr	tr	tr	tr	0.03 ± 0.01 ^a	0.01 ± 0.00 ^b	0.01 ± 0.00 ^b	0.01 ± 0.01 ^b	RI, MS	Fruity, berry	
21	3-Methyl-1-pentanol****	1315	ND	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS	Fusel, cognac, wine, cocoa, green, fruity
24	(E)-3-Hexen-1-ol	1350	ND	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	0.03 ± 0.00	tr	tr	RI, MS, ES	Green, leafy
25	1-Hexanol	1343	ND	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	0.02 ± 0.00	tr	tr	tr	tr	RI, MS, ES	Pungen, ethereal, fusel, oily, fruity, alcoholic, sweet, green
26	2-Methyl-4-heptanol****	1385	ND	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	0.74 ± 0.00	tr	tr	tr	tr	RI, MS	--
30	2-Ethyl-1-hexanol	1480	ND	ND	0.01 ± 0.00	tr	tr	tr	0.01 ± 0.00	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	0.01 ± 0.00	tr	RI, MS, ES	Citrus, fresh, floral, oily, sweet
35	Linalool	1537	ND	ND	ND	tr	0.08 ± 0.00 ^b	tr	0.14 ± 0.00 ^a	0.07 ± 0.06 ^d	0.15 ± 0.00 ^b	0.35 ± 0.00 ^b	0.55 ± 0.00 ^a	tr	tr	tr	tr	tr	0.11 ± 0.01 ^b	tr	0.11 ± 0.00 ^b	0.14 ± 0.03 ^b	RI, MS, ES	Citrus, floral, sweet, rose, woody, green, blueberry	
36	1-Octanol	1550	ND	ND	ND	tr	tr	0.01 ± 0.00	tr	tr	0.01 ± 0.00 ^a	0.02 ± 0.00 ^b	0.04 ± 0.00 ^a	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS, ES	Waxy, green, orange, aldehydic, rose, mushroom
38	2,3-Butanediol****	1571	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS	Fruity, creamy, buttery
52	Phenylethyl alcohol****	1891	ND	ND	ND	0.36 ± 0.33	tr	tr	tr	tr	tr	tr	tr	tr	tr	0.01 ± 0.00 ^b	0.44 ± 0.07 ^a	1.27 ± 0.01 ^b	1.53 ± 0.00 ^b	2.23 ± 0.10 ^a	2.20 ± 0.41 ^a	tr	tr	RI, MS	Sweet, floral, fresh, breadly, rose, honey
	Total alcohol		ND	ND	0.01 ± 0.00	0.49 ± 0.36^a	0.12 ± 0.02^c	0.09 ± 0.02^c	0.23 ± 0.03^b	0.45 ± 0.14^c	1.18 ± 0.23^a	0.48 ± 0.07^c	0.84 ± 0.09^b	0.07 ± 0.00^b	0.02 ± 0.00^b	0.02 ± 0.00^b	0.44 ± 0.07^a	2.85 ± 0.18^a	1.56 ± 0.01^b	2.40 ± 0.11^a	2.37 ± 0.45^a				
13	3-Hydroxybutanal****	1243	ND	tr	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS	--
19	Octanal	1289	1.11 ± 0.15 ^a	0.89 ± 0.10 ^b	0.62 ± 0.51 ^c	0.76 ± 0.08 ^b	2.06 ± 0.52 ^a	1.40 ± 0.17 ^b	1.52 ± 0.31 ^{ab}	1.48 ± 0.26 ^{ab}	1.69 ± 0.98 ^a	1.02 ± 0.82 ^b	0.39 ± 0.60 ^c	1.04 ± 1.00 ^b	3.14 ± 1.13 ^a	2.55 ± 1.05 ^b	1.01 ± 0.79 ^c	2.25 ± 0.74 ^b	2.22 ± 1.23 ^a	2.33 ± 0.89 ^b	0.57 ± 0.88 ^c	1.26 ± 1.13 ^b	RI, MS	Aldehydic, waxy, citrus, orange, green, peely	
	Total aldehyde		1.11 ± 0.15^a	0.89 ± 0.10^b	0.62 ± 0.51^c	0.76 ± 0.08^b	2.06 ± 0.52^a	1.40 ± 0.17^b	1.52 ± 0.31^{ab}	1.48 ± 0.26^{ab}	1.69 ± 0.98^a	1.02 ± 0.82^b	0.39 ± 0.60^c	1.04 ± 1.00^b	3.14 ± 1.13^a	2.55 ± 1.05^b	1.01 ± 0.79^c	2.25 ± 0.74^b	2.22 ± 1.23^a	2.33 ± 0.89^b	0.57 ± 0.88^c	1.26 ± 1.13^b			

15	Terpinolene****	125 5	ND	ND	ND	ND	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	tr	RI, MS, ES	Sweet, fresh pine, citrus, woody, lemon peel
	Total terpene		ND	ND	ND	ND	0.02 ± 0.00 ^b	0.02 ± 0.00 ^b	0.02 ± 0.00 ^b	0.02 ± 0.00 ^b	0.03 ± 0.00 ^b	0.03 ± 0.00 ^b	0.03 ± 0.00 ^b	0.03 ± 0.00 ^b	0.02 ± 0.00 ^a	0.02 ± 0.00 ^a	0.02 ± 0.00 ^a	0.02 ± 0.00 ^a	0.02 ± 0.00 ^b	0.02 ± 0.00 ^b	0.02 ± 0.00 ^{ab}	0.02 ± 0.00 ^{ab}			
	Total		42.60 ± 29.24 ^a	33.76 ± 2.89 ^b	15.51 ± 1.88 ^c	8.55 ± 1.85 ^d	31.78 ± 8.86 ^a	34.58 ± 13.31 ^a	25.26 ± 5.84 ^b	17.24 ± 5.11 ^c	22.36 ± 13.22 ^b	38.84 ± 17.86 ^a	16.38 ± 14.20 ^c	17.66 ± 8.70 ^b	38.54 ± 12.76 ^c	55.51 ± 20.78 ^a	44.85 ± 25.35 ^b	20.90 ± 7.13 ^d	34.89 ± 16.27 ^c	41.05 ± 19.27 ^b	51.18 ± 14.98 ^a	21.57 ± 10.35 ^d			

*RI = retention index

**ID = Identification by comparison of mass spectra (MS) with in-house NIST library 11.0, MS and retention index (RI) with NIST Webbook (NIST²⁴), and MS and retention time with external standards

*** = Odor description were obtained from TGSC²⁵.

****= Semi-quantification: Ethoxyacetic acid, acetic acid, pterin-6-carboxylic acid, 2-methyl-1-propanol, 2-pentanol, 2-hexanol, 1-(2-aminoethylamino)-2-propanol, 2-octen-4-ol, 3-methyl-1-pentanol, 2-methyl-4-heptanol, 2,3-butanediol, phenylethyl alcohol, 3-hydroxybutanal, myristicin, 3-hydroxymandelic acid ethyl ester, ethyl (*E*)-3-hexenoate, ethyl lactate, propyl hexanoate, ethyl 3-hydroxybutanoate, isoamyl lactate, ethyl 2-furancarboxylate, ethyl 9-decenoate and 3-hydroxy-2-butanone were semi-quantified using the calibration curves of ethyl hexanoate, ethyl octanoate, 2-ethyl-1-hexanol, 3-methylbutyl acetate, 1-butanol, 1-butanol, ethyl hexanoate, 1-hexanol, 1-hexanol, 1-hexanol, 1-octanol, benzyl alcohol, ethyl hexanoate, benzaldehyde, 1-butanol, 1-hexanol, 2-heptanol, 2-heptanol, ethyl nonanoate, 1-octanol, benzeneacetaldehyde, diethyl succinate and ethyl hexanoate, respectively.

Samples of M, 2GM, 4GM etc. refer to Table S3.

ND = Not detected or value below the detection threshold.

tr = trace

-- = Not available.

a-d = Values with different superscripts in the same wine group are significantly different ($p \leq 0.05$)

Table S6 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated Chardonnay wines

(continued 1/7)

Variables	TA	pH	TPC	TFC	DPPH	FRAP	ABTS	L*	a*	b*	Quinic acid	Gallic acid	Chlorogenic acid	Total phenolic acids	Gallocatechin
TA	1	-0.967	-0.729	-0.748	-0.802	-0.743	-0.723	0.451	-0.375	-0.572	-0.602	-0.590	-0.739	-0.701	-0.806
pH	-0.967	1	0.786	0.801	0.857	0.790	0.777	-0.495	0.409	0.601	0.617	0.664	0.794	0.743	0.869
TPC	-0.729	0.786	1	0.940	0.975	0.844	0.910	-0.434	0.448	0.563	0.792	0.875	0.949	0.954	0.945
TFC	-0.748	0.801	0.940	1	0.932	0.916	0.961	-0.295	0.227	0.422	0.635	0.890	0.970	0.852	0.961
DPPH	-0.802	0.857	0.975	0.932	1	0.821	0.919	-0.494	0.462	0.618	0.758	0.854	0.931	0.921	0.970
FRAP	-0.743	0.790	0.844	0.916	0.821	1	0.917	-0.131	0.080	0.257	0.501	0.832	0.940	0.736	0.870
ABTS	-0.723	0.777	0.910	0.961	0.919	0.917	1	-0.168	0.112	0.301	0.504	0.926	0.966	0.773	0.938
L*	0.451	-0.495	-0.434	-0.295	-0.494	-0.131	-0.168	1	-0.893	-0.971	-0.747	-0.102	-0.232	-0.580	-0.452
a*	-0.375	0.409	0.448	0.227	0.462	0.080	0.112	-0.893	1	0.883	0.822	0.149	0.233	0.650	0.376
b*	-0.572	0.601	0.563	0.422	0.618	0.257	0.301	-0.971	0.883	1	0.814	0.250	0.359	0.692	0.575
Quinic acid	-0.602	0.617	0.792	0.635	0.758	0.501	0.504	-0.747	0.822	0.814	1	0.466	0.639	0.925	0.676
Gallic acid	-0.590	0.664	0.875	0.890	0.854	0.832	0.926	-0.102	0.149	0.250	0.466	1	0.899	0.765	0.877
Chlorogenic acid	-0.739	0.794	0.949	0.970	0.931	0.940	0.966	-0.232	0.233	0.359	0.639	0.899	1	0.861	0.934
Total phenolic acids	-0.701	0.743	0.954	0.852	0.921	0.736	0.773	-0.580	0.650	0.692	0.925	0.765	0.861	1	0.873
Gallocatechin	-0.806	0.869	0.945	0.961	0.970	0.870	0.938	-0.452	0.376	0.575	0.676	0.877	0.934	0.873	1
Epigallocatechin	-0.684	0.749	0.909	0.962	0.900	0.928	0.992	-0.128	0.088	0.259	0.496	0.931	0.972	0.770	0.929
Procyanidin B1	-0.713	0.766	0.905	0.981	0.905	0.886	0.955	-0.270	0.179	0.405	0.575	0.883	0.941	0.805	0.959
Catechin	-0.743	0.804	0.958	0.924	0.975	0.765	0.894	-0.524	0.476	0.630	0.758	0.834	0.894	0.910	0.965
Procyanidin B2	-0.587	0.643	0.621	0.713	0.603	0.879	0.722	-0.020	-0.017	0.137	0.285	0.698	0.742	0.518	0.664
Epicatechin	-0.674	0.740	0.729	0.805	0.779	0.697	0.807	-0.312	0.159	0.418	0.368	0.731	0.766	0.590	0.851
Epigallocatechin gallate	-0.646	0.684	0.910	0.780	0.862	0.663	0.679	-0.635	0.711	0.730	0.963	0.661	0.796	0.983	0.801
Procyanidin B1 3-O-gallate	-0.469	0.576	0.551	0.670	0.569	0.665	0.636	-0.261	0.068	0.269	0.281	0.523	0.636	0.443	0.684
Gallocatechin gallate	-0.625	0.680	0.763	0.874	0.772	0.905	0.955	0.053	-0.149	0.067	0.254	0.866	0.884	0.569	0.831
Quercetin 3-O-di-galactoside/glucoside	-0.692	0.767	0.792	0.872	0.837	0.842	0.923	-0.199	0.054	0.306	0.351	0.828	0.870	0.623	0.895
Myricetin 3-O-galactoside 1	-0.347	0.380	0.484	0.587	0.527	0.402	0.595	-0.142	0.013	0.206	0.232	0.540	0.473	0.398	0.603
Epicatechin gallate	-0.671	0.745	0.737	0.816	0.787	0.711	0.837	-0.308	0.153	0.403	0.345	0.778	0.759	0.589	0.870
Rutin	-0.704	0.760	0.776	0.818	0.825	0.671	0.801	-0.460	0.310	0.519	0.513	0.682	0.768	0.674	0.878
Quercetin 3-O-galactoside	-0.558	0.600	0.417	0.494	0.527	0.397	0.520	-0.330	0.121	0.359	0.186	0.372	0.412	0.301	0.608

Kaempferol 3-O-rutinoside 2	-0.482	0.527	0.491	0.636	0.562	0.506	0.665	-0.198	0.000	0.275	0.141	0.604	0.533	0.363	0.688
Kaempferol 3-O-glucoside	-0.341	0.416	0.221	0.293	0.331	0.139	0.249	-0.486	0.328	0.466	0.164	0.229	0.189	0.213	0.432
Unknown quercetin conjugate	-0.611	0.666	0.817	0.921	0.826	0.809	0.910	-0.200	0.086	0.333	0.450	0.872	0.849	0.704	0.899
Unknown kaempferol conjugate	-0.624	0.668	0.840	0.870	0.852	0.658	0.831	-0.406	0.284	0.517	0.570	0.766	0.789	0.745	0.890
Total flavonoids	-0.760	0.824	0.964	0.958	0.979	0.835	0.941	-0.445	0.388	0.560	0.703	0.877	0.935	0.892	0.986
Caffeine	-0.668	0.729	0.752	0.860	0.756	0.962	0.911	0.015	-0.085	0.095	0.314	0.819	0.902	0.597	0.807
Ethoxyacetic acid	-0.226	0.220	0.167	0.333	0.186	0.260	0.352	0.257	-0.383	-0.229	-0.207	0.303	0.302	-0.008	0.272
Acetic acid	-0.054	0.059	0.276	0.288	0.256	0.068	0.225	-0.300	0.218	0.253	0.309	0.233	0.198	0.317	0.279
Pterin-6-carboxylic acid	-0.567	0.614	0.856	0.724	0.827	0.569	0.673	-0.597	0.643	0.644	0.866	0.618	0.739	0.895	0.762
Hexanoic acid	-0.005	-0.004	-0.316	-0.395	-0.275	-0.354	-0.359	-0.009	0.038	-0.065	-0.159	-0.400	-0.327	-0.284	-0.343
Octanoic acid	0.507	-0.540	-0.668	-0.606	-0.687	-0.531	-0.533	0.696	-0.591	-0.756	-0.726	-0.458	-0.543	-0.721	-0.685
Nonanoic acid	0.459	-0.467	-0.539	-0.442	-0.549	-0.389	-0.350	0.708	-0.629	-0.779	-0.702	-0.332	-0.377	-0.644	-0.510
Total acid	0.433	-0.459	-0.632	-0.589	-0.638	-0.518	-0.513	0.611	-0.517	-0.683	-0.669	-0.469	-0.517	-0.683	-0.643
2-Methyl-1-propanol	-0.361	0.400	0.379	0.329	0.400	0.188	0.342	-0.304	0.262	0.281	0.304	0.267	0.340	0.340	0.367
2-Pentanol	-0.381	0.389	0.246	0.394	0.281	0.555	0.450	0.304	-0.472	-0.200	-0.145	0.338	0.406	0.058	0.314
1-Butanol	-0.183	0.179	0.368	0.418	0.394	0.198	0.474	-0.079	-0.039	0.095	0.144	0.372	0.337	0.264	0.379
2-Hexanol	-0.184	0.198	0.308	0.388	0.347	0.158	0.394	-0.178	-0.007	0.156	0.127	0.273	0.290	0.214	0.339
3-Methyl-1-butanol	-0.145	0.180	0.105	0.011	0.120	-0.065	0.023	-0.307	0.288	0.230	0.187	-0.019	0.042	0.128	0.070
1-(2-Aminoethylamino)-2-propanol	-0.680	0.748	0.873	0.902	0.866	0.891	0.925	-0.207	0.124	0.308	0.514	0.802	0.921	0.732	0.882
2-Octen-4-ol	-0.718	0.776	0.916	0.950	0.910	0.898	0.974	-0.201	0.138	0.315	0.539	0.875	0.956	0.778	0.924
3-Methyl-1-pentanol	-0.133	0.154	-0.091	-0.128	-0.031	-0.198	-0.147	-0.293	0.219	0.185	0.058	-0.254	-0.123	-0.060	-0.055
(E)-3-Hexen-1-ol	-0.176	0.171	0.089	0.271	0.051	0.307	0.171	0.264	-0.280	-0.226	-0.051	0.143	0.195	0.034	0.102
1-Hexanol	-0.418	0.399	0.084	0.251	0.144	0.331	0.256	0.189	-0.393	-0.086	-0.226	0.150	0.209	-0.084	0.213
2-Methyl-4-heptanol	-0.201	0.286	-0.003	0.166	0.063	0.183	0.131	-0.039	-0.100	0.037	-0.183	0.118	0.107	-0.076	0.211
2-Ethyl-1-hexanol	-0.201	0.225	0.372	0.157	0.341	0.035	0.076	-0.691	0.775	0.645	0.731	0.057	0.171	0.547	0.234
Linalool	-0.687	0.745	0.944	0.829	0.937	0.670	0.786	-0.577	0.615	0.670	0.851	0.759	0.841	0.944	0.877
1-Octanol	0.551	-0.541	-0.598	-0.607	-0.621	-0.592	-0.575	0.338	-0.286	-0.474	-0.475	-0.564	-0.560	-0.585	-0.642
2,3-Butanediol	-0.147	0.113	0.133	0.130	0.139	0.007	0.075	-0.325	0.234	0.279	0.240	0.079	0.057	0.201	0.177
Phenylethyl alcohol	0.220	-0.214	-0.305	-0.419	-0.303	-0.442	-0.400	-0.102	0.172	-0.027	-0.074	-0.391	-0.357	-0.224	-0.366
Total alcohol	-0.141	0.167	0.114	0.051	0.137	-0.060	0.077	-0.248	0.197	0.172	0.137	0.008	0.064	0.105	0.088
3-Hydroxybutanal	-0.134	0.178	0.192	0.142	0.243	0.097	0.175	-0.286	0.169	0.275	0.157	0.056	0.149	0.145	0.222
Octanal	0.101	-0.089	0.232	0.321	0.143	0.250	0.284	0.287	-0.226	-0.288	0.015	0.371	0.320	0.172	0.195
Total aldehyde	0.099	-0.087	0.233	0.323	0.145	0.251	0.285	0.285	-0.224	-0.287	0.017	0.372	0.322	0.174	0.197

Benzaldehyde	0.253	-0.216	-0.240	-0.371	-0.199	-0.338	-0.324	-0.276	0.229	0.198	0.004	-0.362	-0.366	-0.161	-0.262
Myristicin	0.426	-0.438	-0.354	-0.557	-0.400	-0.502	-0.513	-0.002	0.225	-0.088	-0.001	-0.455	-0.432	-0.203	-0.539
Benzeneacetaldehyde	-0.330	0.416	0.396	0.368	0.482	0.214	0.374	-0.606	0.423	0.561	0.324	0.297	0.277	0.358	0.454
Naphthalene	-0.815	0.869	0.928	0.887	0.964	0.772	0.842	-0.636	0.547	0.734	0.803	0.753	0.852	0.909	0.939
1,2-Dihydro-1,1,6-trimethyl-naphthalene	-0.815	0.869	0.928	0.887	0.964	0.772	0.842	-0.636	0.547	0.734	0.803	0.753	0.852	0.909	0.939
Benzyl alcohol	-0.298	0.273	0.021	0.134	0.065	0.231	0.160	0.187	-0.290	-0.071	-0.229	0.127	0.108	-0.105	0.134
Total benzenoid	-0.289	0.276	0.002	0.079	0.063	0.178	0.121	0.086	-0.203	0.016	-0.208	0.076	0.054	-0.113	0.109
Ethyl butanoate	-0.387	0.380	0.176	0.342	0.226	0.299	0.361	0.166	-0.340	-0.110	-0.168	0.236	0.309	-0.003	0.292
Ethyl 3-methylbutanoate	-0.195	0.139	-0.105	0.010	-0.058	0.001	0.038	0.252	-0.316	-0.262	-0.272	-0.048	0.029	-0.204	-0.028
3-Methylbutyl acetate	-0.342	0.333	0.117	0.320	0.169	0.326	0.342	0.281	-0.453	-0.196	-0.282	0.275	0.264	-0.074	0.272
3-Hydroxymandelic acid ethyl ester	0.016	-0.006	0.004	0.215	0.008	0.104	0.167	0.219	-0.340	-0.257	-0.255	0.199	0.110	-0.097	0.149
Ethyl hexanoate	0.116	-0.130	-0.188	-0.087	-0.204	-0.107	-0.049	0.442	-0.469	-0.449	-0.388	-0.109	-0.049	-0.313	-0.192
Hexyl acetate	-0.209	0.186	-0.013	0.180	0.020	0.204	0.206	0.424	-0.529	-0.351	-0.379	0.175	0.141	-0.189	0.109
Ethyl (E)-3-hexenoate	-0.464	0.502	0.170	0.318	0.253	0.245	0.289	-0.202	-0.023	0.174	-0.009	0.231	0.211	0.096	0.350
Ethyl lactate	0.256	-0.258	-0.137	-0.214	-0.171	-0.275	-0.202	-0.075	0.102	-0.040	0.006	-0.189	-0.186	-0.081	-0.211
Propyl hexanoate	0.105	-0.089	-0.036	0.020	-0.064	-0.078	0.024	0.253	-0.255	-0.305	-0.167	-0.035	0.051	-0.124	-0.060
Ethyl octanoate	0.571	-0.607	-0.514	-0.420	-0.578	-0.403	-0.373	0.724	-0.629	-0.777	-0.648	-0.285	-0.380	-0.590	-0.546
Ethyl 3-hydroxybutanoate	0.660	-0.621	-0.574	-0.619	-0.590	-0.526	-0.531	0.350	-0.274	-0.455	-0.487	-0.483	-0.526	-0.561	-0.649
Ethyl nonanoate	0.112	-0.139	-0.264	-0.125	-0.244	-0.185	-0.128	0.335	-0.464	-0.338	-0.453	-0.162	-0.167	-0.390	-0.184
Isoamyl lactate	-0.174	0.178	0.226	0.183	0.230	0.021	0.105	-0.491	0.405	0.402	0.419	0.032	0.119	0.316	0.227
Methyl decanoate	0.253	-0.210	-0.278	-0.147	-0.236	-0.157	-0.081	0.341	-0.519	-0.434	-0.494	-0.248	-0.142	-0.446	-0.174
Ethyl 2-furancarboxylate	0.470	-0.453	-0.367	-0.461	-0.410	-0.421	-0.431	0.082	0.012	-0.195	-0.144	-0.437	-0.394	-0.293	-0.473
Ethyl decanoate	0.439	-0.499	-0.470	-0.387	-0.524	-0.325	-0.319	0.783	-0.578	-0.799	-0.630	-0.173	-0.329	-0.533	-0.491
Diethyl succinate	0.391	-0.383	-0.253	-0.479	-0.291	-0.449	-0.421	-0.127	0.308	0.046	0.057	-0.346	-0.368	-0.119	-0.403
Ethyl 9-decenoate	0.413	-0.481	-0.457	-0.444	-0.496	-0.374	-0.349	0.656	-0.409	-0.678	-0.549	-0.165	-0.346	-0.476	-0.486
Ethyl benzeneacetate	-0.470	0.459	0.373	0.165	0.412	0.105	0.105	-0.785	0.843	0.785	0.716	0.092	0.168	0.549	0.307
2-Phenethyl acetate	0.126	-0.129	-0.153	-0.245	-0.094	-0.283	-0.136	-0.083	0.045	0.051	-0.157	-0.116	-0.193	-0.170	-0.101
Ethyl dodecanoate	0.173	-0.078	-0.124	-0.076	-0.162	-0.041	-0.057	0.222	-0.286	-0.269	-0.247	-0.144	-0.071	-0.229	-0.106
Total ester	0.049	-0.042	-0.029	0.179	-0.028	0.072	0.138	0.255	-0.364	-0.295	-0.288	0.175	0.082	-0.132	0.110
3-Hydroxy-2-butanone	0.195	-0.185	-0.247	-0.239	-0.280	-0.203	-0.183	0.405	-0.227	-0.417	-0.305	-0.094	-0.187	-0.265	-0.259
Limonene	-0.768	0.823	0.847	0.811	0.893	0.692	0.746	-0.700	0.549	0.777	0.780	0.629	0.751	0.841	0.865
Terpinolene	-0.626	0.693	0.746	0.794	0.742	0.873	0.838	-0.083	0.001	0.171	0.369	0.704	0.843	0.589	0.755

Total terpene	-0.809	0.868	0.895	0.878	0.932	0.794	0.830	-0.607	0.465	0.694	0.749	0.704	0.839	0.854	0.912
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Values in **bold** indicate significant ($p < 0.05$) correlation

Table S6 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated Chardonnay wines

(continued 2/7)

Variables	Epigallocatechin	Procyanidin B1	Catechin	Procyanidin B2	Epicatechin	Epigallocatechin gallate	Procyanidin B1 3-O-gallate	Gallocatechin gallate	Quercetin 3-O-di-galactoside/glucoside	Myricetin 3-O-galactoside 1	Epicatechin gallate	Rutin	Quercetin 3-O-galactoside	Kaempferol 3-O-rutinoside 2	Kaempferol 3-O-glucoside
TA	-0.684	-0.713	-0.743	-0.587	-0.674	-0.646	-0.469	-0.625	-0.692	-0.347	-0.671	-	-0.558	-0.482	-0.341
pH	0.749	0.766	0.804	0.643	0.740	0.684	0.576	0.680	0.767	0.380	0.745	0.704	0.600	0.527	0.416
TPC	0.909	0.905	0.958	0.621	0.729	0.910	0.551	0.763	0.792	0.484	0.737	0.776	0.417	0.491	0.221
TFC	0.962	0.981	0.924	0.713	0.805	0.780	0.670	0.874	0.872	0.587	0.816	0.818	0.494	0.636	0.293
DPPH	0.900	0.905	0.975	0.603	0.779	0.862	0.569	0.772	0.837	0.527	0.787	0.825	0.527	0.562	0.331
FRAP	0.928	0.886	0.765	0.879	0.697	0.663	0.665	0.905	0.842	0.402	0.711	0.671	0.397	0.506	0.139
ABTS	0.992	0.955	0.894	0.722	0.807	0.679	0.636	0.955	0.923	0.595	0.837	0.801	0.520	0.665	0.249
L*	-0.128	-0.270	-0.524	-0.020	-0.312	-0.635	-0.261	0.053	-0.199	-0.142	-0.308	-	-0.330	-0.198	-0.486
a*	0.088	0.179	0.476	-0.017	0.159	0.711	0.068	-0.149	0.054	0.013	0.153	0.460	0.310	0.121	0.000
b*	0.259	0.405	0.630	0.137	0.418	0.730	0.269	0.067	0.306	0.206	0.403	0.519	0.359	0.275	0.466
Quinic acid	0.496	0.575	0.758	0.285	0.368	0.963	0.281	0.254	0.351	0.232	0.345	0.513	0.186	0.141	0.164
Gallic acid	0.931	0.883	0.834	0.698	0.731	0.661	0.523	0.866	0.828	0.540	0.778	0.682	0.372	0.604	0.229
Chlorogenic acid	0.972	0.941	0.894	0.742	0.766	0.796	0.636	0.884	0.870	0.473	0.759	0.768	0.412	0.533	0.189
Total phenolic acids	0.770	0.805	0.910	0.518	0.590	0.983	0.443	0.569	0.623	0.398	0.589	0.674	0.301	0.363	0.213
Gallocatechin	0.929	0.959	0.965	0.664	0.851	0.801	0.684	0.831	0.895	0.603	0.870	0.878	0.608	0.688	0.432
Epigallocatechin	1	0.955	0.884	0.722	0.800	0.679	0.675	0.956	0.919	0.583	0.831	0.798	0.499	0.645	0.236
Procyanidin B1	0.955	1	0.912	0.676	0.845	0.729	0.680	0.882	0.893	0.661	0.849	0.834	0.550	0.728	0.348
Catechin	0.884	0.912	1	0.487	0.780	0.853	0.615	0.748	0.814	0.644	0.820	0.882	0.600	0.634	0.411
Procyanidin B2	0.722	0.676	0.487	1	0.534	0.449	0.475	0.744	0.669	0.165	0.526	0.386	0.206	0.362	0.084
Epicatechin	0.800	0.845	0.780	0.534	1	0.532	0.719	0.757	0.926	0.450	0.944	0.888	0.583	0.763	0.507
Epigallocatechin gallate	0.679	0.729	0.853	0.449	0.532	1	0.407	0.458	0.538	0.298	0.509	0.622	0.221	0.273	0.178
Procyanidin B1 3-O-gallate	0.675	0.680	0.615	0.475	0.719	0.407	1	0.664	0.790	0.459	0.782	0.804	0.554	0.686	0.619
Gallocatechin gallate	0.956	0.882	0.748	0.744	0.757	0.458	0.664	1	0.915	0.582	0.809	0.729	0.530	0.677	0.202
Quercetin 3-O-di-galactoside/glucoside	0.919	0.893	0.814	0.669	0.926	0.538	0.790	0.915	1	0.508	0.927	0.870	0.581	0.755	0.425
Myricetin 3-O-galactoside 1	0.583	0.661	0.644	0.165	0.450	0.298	0.459	0.582	0.508	1	0.613	0.639	0.765	0.794	0.545
Epicatechin gallate	0.831	0.849	0.820	0.526	0.944	0.509	0.782	0.809	0.927	0.613	1	0.929	0.690	0.870	0.623

Rutin	0.798	0.834	0.882	0.386	0.888	0.622	0.804	0.729	0.870	0.639	0.929	1	0.754	0.782	0.624
Quercetin 3-O-galactoside	0.499	0.550	0.600	0.206	0.583	0.221	0.554	0.530	0.581	0.765	0.690	0.754	1	0.710	0.664
Kaempferol 3-O-rutinoside 2	0.645	0.728	0.634	0.362	0.763	0.273	0.686	0.677	0.755	0.794	0.870	0.782	0.710	1	0.749
Kaempferol 3-O-glucoside	0.236	0.348	0.411	0.084	0.507	0.178	0.619	0.202	0.425	0.545	0.623	0.624	0.664	0.749	1
Unknown quercetin conjugate	0.902	0.959	0.850	0.646	0.775	0.611	0.649	0.862	0.842	0.784	0.836	0.777	0.582	0.826	0.441
Unknown kaempferol conjugate	0.823	0.903	0.912	0.393	0.869	0.698	0.640	0.721	0.818	0.704	0.889	0.910	0.628	0.781	0.495
Total flavonoids	0.934	0.950	0.991	0.579	0.818	0.824	0.661	0.823	0.871	0.639	0.854	0.888	0.599	0.666	0.396
Caffeine	0.919	0.836	0.685	0.865	0.708	0.504	0.680	0.946	0.884	0.368	0.715	0.642	0.382	0.523	0.135
Ethoxyacetic acid	0.360	0.377	0.211	0.170	0.551	-0.044	0.385	0.433	0.475	0.283	0.487	0.467	0.394	0.438	0.211
Acetic acid	0.228	0.312	0.414	-0.245	0.165	0.309	0.272	0.164	0.156	0.565	0.281	0.411	0.296	0.366	0.251
Pterin-6-carboxylic acid	0.678	0.682	0.881	0.217	0.501	0.895	0.491	0.491	0.549	0.471	0.561	0.727	0.392	0.356	0.255
Hexanoic acid	-0.370	-0.420	-0.286	-0.407	-0.198	-0.237	-0.299	-0.343	-0.303	-0.378	-0.266	-	-0.047	-0.393	-0.209
Octanoic acid	-0.515	-0.588	-0.706	-0.392	-0.396	-0.700	-0.446	-0.390	-0.461	-0.435	-0.481	-	-0.381	-0.415	-0.376
Nonanoic acid	-0.314	-0.404	-0.524	-0.364	-0.219	-0.646	-0.186	-0.182	-0.252	-0.198	-0.279	-	-0.119	-0.218	-0.216
Total acid	-0.496	-0.575	-0.647	-0.435	-0.355	-0.657	-0.405	-0.378	-0.431	-0.419	-0.441	-	-0.289	-0.414	-0.345
2-Methyl-1-propanol	0.333	0.330	0.470	-0.094	0.427	0.345	0.273	0.283	0.343	0.314	0.440	0.554	0.460	0.286	0.206
2-Pentanol	0.445	0.378	0.152	0.660	0.342	-0.003	0.394	0.562	0.515	0.099	0.319	0.193	0.192	0.264	0.028
1-Butanol	0.429	0.458	0.486	-0.016	0.318	0.208	0.181	0.452	0.356	0.754	0.442	0.497	0.495	0.594	0.261
2-Hexanol	0.344	0.412	0.418	0.020	0.363	0.182	0.204	0.362	0.338	0.601	0.413	0.496	0.469	0.519	0.287
3-Methyl-1-butanol	0.019	-0.006	0.187	-0.293	0.124	0.158	0.073	-0.017	0.044	0.026	0.136	0.262	0.221	-0.019	0.038
1-(2-Aminoethylamino)-2-propanol	0.938	0.876	0.852	0.650	0.791	0.674	0.718	0.891	0.893	0.449	0.804	0.809	0.481	0.540	0.199
2-Octen-4-ol	0.980	0.932	0.908	0.635	0.807	0.703	0.699	0.929	0.911	0.577	0.845	0.855	0.548	0.628	0.246
3-Methyl-1-pentanol	-0.149	-0.137	0.042	-0.391	0.043	-0.022	0.088	-0.161	-0.071	0.032	0.050	0.212	0.354	-0.041	0.189
(E)-3-Hexen-1-ol	0.207	0.184	0.034	0.391	0.067	-0.010	0.146	0.224	0.063	0.069	0.116	0.029	0.008	0.041	-0.023
1-Hexanol	0.239	0.264	0.040	0.458	0.495	-0.116	0.225	0.337	0.412	-0.093	0.359	0.205	0.200	0.240	0.065
2-Methyl-4-heptanol	0.144	0.217	0.062	0.319	0.438	-0.097	0.572	0.192	0.365	0.179	0.454	0.338	0.355	0.550	0.730
2-Ethyl-1-hexanol	0.067	0.105	0.411	-0.203	-0.080	0.612	0.018	-0.117	-0.081	0.158	0.004	0.207	0.105	-0.089	0.082
Linalool	0.785	0.787	0.950	0.371	0.676	0.918	0.509	0.598	0.696	0.440	0.684	0.791	0.426	0.412	0.275
1-Octanol	-0.545	-0.595	-0.555	-0.619	-0.461	-0.532	-0.343	-0.476	-0.540	-0.279	-0.482	-	-0.197	-0.439	-0.284
												0.409			

2,3-Butanediol	0.080	0.146	0.278	-0.299	0.086	0.200	0.288	0.042	0.086	0.379	0.220	0.333	0.266	0.282	0.283
Phenylethyl alcohol	-0.396	-0.432	-0.240	-0.556	-0.299	-0.162	-0.249	-0.404	-0.384	-0.224	-0.296	-	-0.060	-0.345	-0.124
Total alcohol	0.062	0.044	0.213	-0.291	0.150	0.124	0.072	0.049	0.078	0.157	0.177	0.306	0.297	0.081	0.076
3-Hydroxybutanal	0.160	0.156	0.250	-0.007	0.316	0.176	0.186	0.136	0.259	-0.006	0.267	0.267	0.151	0.118	0.121
Octanal	0.342	0.319	0.216	0.104	0.176	0.145	0.312	0.323	0.230	0.308	0.204	0.243	0.069	0.171	-0.022
Total aldehyde	0.344	0.321	0.217	0.105	0.178	0.147	0.314	0.324	0.232	0.308	0.206	0.245	0.071	0.172	-0.021
Benzaldehyde	-0.351	-0.372	-0.179	-0.329	-0.385	-0.122	-0.279	-0.344	-0.344	-0.155	-0.306	-	-0.131	-0.267	-0.092
Myristicin	-0.513	-0.595	-0.408	-0.500	-0.626	-0.125	-0.571	-0.558	-0.616	-0.492	-0.635	0.311	-0.540	-0.655	-0.475
Benzeneacetaldehyde	0.322	0.331	0.520	0.167	0.361	0.320	0.352	0.294	0.382	0.441	0.521	0.538	0.532	0.489	0.564
Naphthalene	0.815	0.852	0.954	0.553	0.710	0.865	0.575	0.684	0.764	0.526	0.751	0.799	0.534	0.556	0.390
1,2-Dihydro-1,1,6-trimethylnaphthalene	0.815	0.852	0.954	0.553	0.710	0.865	0.575	0.684	0.764	0.526	0.751	0.799	0.534	0.556	0.390
Benzyl alcohol	0.139	0.156	-0.047	0.432	0.382	-0.140	0.072	0.217	0.303	-0.198	0.258	0.069	0.033	0.181	0.034
Total benzenoid	0.087	0.094	-0.055	0.396	0.342	-0.143	0.019	0.165	0.265	-0.242	0.236	0.038	0.025	0.149	0.043
Ethyl butanoate	0.349	0.376	0.199	0.249	0.636	-0.023	0.389	0.422	0.530	0.123	0.532	0.459	0.329	0.446	0.226
Ethyl 3-methylbutanoate	0.038	0.022	-0.059	-0.099	0.269	-0.203	0.151	0.115	0.178	-0.055	0.198	0.196	0.181	0.121	0.055
3-Methylbutyl acetate	0.331	0.359	0.120	0.391	0.566	-0.126	0.351	0.435	0.503	0.143	0.487	0.349	0.304	0.459	0.239
3-Hydroxymandelic acid ethyl ester	0.194	0.247	0.107	0.039	0.290	-0.153	0.396	0.257	0.256	0.419	0.333	0.353	0.361	0.428	0.354
Ethyl hexanoate	-0.036	-0.054	-0.219	-0.123	0.203	-0.279	0.005	0.036	0.076	-0.207	0.053	0.013	-0.108	0.000	-0.167
Hexyl acetate	0.203	0.206	-0.032	0.301	0.380	-0.236	0.207	0.312	0.336	0.055	0.309	0.177	0.174	0.304	0.123
Ethyl (E)-3-hexenoate	0.267	0.321	0.311	0.126	0.439	0.034	0.416	0.328	0.387	0.378	0.498	0.508	0.565	0.474	0.434
Ethyl lactate	-0.190	-0.217	-0.068	-0.470	-0.228	-0.044	-0.131	-0.196	-0.248	0.011	-0.181	-	-0.024	-0.197	-0.176
Propyl hexanoate	0.060	0.041	0.002	-0.266	0.200	-0.091	0.153	0.070	0.093	0.053	0.120	0.043	0.130	0.037	-0.038
Ethyl octanoate	-0.336	-0.387	-0.552	-0.334	-0.268	-0.572	-0.280	-0.233	-0.322	-0.249	-0.340	-	-0.337	-0.262	-0.349
Ethyl 3-hydroxybutanoate	-0.517	-0.608	-0.610	-0.390	-0.506	-0.524	-0.464	-0.437	-0.505	-0.478	-0.561	0.379	-0.465	-0.533	-0.452
Ethyl nonanoate	-0.129	-0.073	-0.245	-0.132	0.234	-0.368	-0.024	-0.046	0.051	-0.153	0.072	0.015	0.020	0.054	-0.041
Isoamyl lactate	0.106	0.172	0.381	-0.328	0.048	0.336	0.271	0.026	0.048	0.432	0.172	0.398	0.371	0.190	0.263
Methyl decanoate	-0.048	-0.088	-0.171	-0.281	0.148	-0.443	0.264	0.078	0.114	0.080	0.085	0.162	0.306	0.090	0.066
Ethyl 2-furancarboxylate	-0.406	-0.473	-0.356	-0.467	-0.490	-0.244	-0.336	-0.397	-0.499	-0.238	-0.466	-	-0.253	-0.453	-0.330
Ethyl decanoate	-0.285	-0.382	-0.526	-0.202	-0.305	-0.553	-0.327	-0.185	-0.317	-0.240	-0.317	0.378	-	-0.248	-0.287
Diethyl succinate	-0.422	-0.482	-0.271	-0.481	-0.491	-0.055	-0.468	-0.470	-0.493	-0.334	-0.490	0.400	-	-0.503	-0.344
												0.399	-0.350		

Ethyl 9-decenoate	-0.324	-0.427	-0.500	-0.276	-0.319	-0.488	-0.393	-0.241	-0.344	-0.285	-0.337	-	-0.343	-0.284	-0.283
Ethyl benzeneacetate	0.063	0.090	0.408	-0.024	-0.013	0.576	-0.022	-0.096	-0.002	0.029	0.073	0.406	0.133	-0.059	0.175
2-Phenethyl acetate	-0.173	-0.147	-0.083	-0.334	0.062	-0.146	-0.062	-0.135	0.040	-0.032	-0.006	0.043	0.101	0.080	0.106
Ethyl dodecanoate	-0.026	-0.036	-0.124	-0.048	0.057	-0.200	0.117	0.032	0.012	0.020	0.006	0.051	0.159	0.005	-0.027
Total ester	0.166	0.213	0.070	0.011	0.267	-0.184	0.365	0.234	0.229	0.389	0.305	0.322	0.332	0.399	0.323
3-Hydroxy-2-butanone	-0.170	-0.224	-0.279	-0.155	-0.173	-0.269	-0.218	-0.133	-0.183	-0.150	-0.191	-	-0.185	-0.137	-0.140
Limonene	0.711	0.774	0.884	0.497	0.638	0.813	0.543	0.588	0.684	0.486	0.678	0.733	0.514	0.507	0.390
Terpinolene	0.853	0.747	0.681	0.727	0.707	0.537	0.666	0.849	0.836	0.228	0.699	0.658	0.330	0.403	0.086
Total terpene	0.805	0.835	0.910	0.594	0.712	0.817	0.618	0.698	0.777	0.469	0.742	0.782	0.521	0.528	0.355

Values in **bold** indicate significant ($p < 0.05$) correlation

Table S6 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated Chardonnay wines

(continued 3/7)

Variables	Unknown quercetin conjugate	Unknown kaempferol conjugate	Total flavonoids	Caffeine	Ethoxyacetic acid	Acetic acid	Pterin-6-carboxylic acid	Hexanoic acid	Octanoic acid	Nonanoic acid	Total acid	2-Methyl-1-propanol	2-Pentanol	1-Butanol	2-Hexanol
TA	-0.611	-0.624	-0.760	-0.668	-0.226	-0.054	-0.567	-0.005	0.507	0.459	0.433	-0.361	-0.381	-0.183	-0.184
pH	0.666	0.668	0.824	0.729	0.220	0.059	0.614	-0.004	-0.540	-0.467	-0.459	0.400	0.389	0.179	0.198
TPC	0.817	0.840	0.964	0.752	0.167	0.276	0.856	-0.316	-0.668	-0.539	-0.632	0.379	0.246	0.368	0.308
TFC	0.921	0.870	0.958	0.860	0.333	0.288	0.724	-0.395	-0.606	-0.442	-0.589	0.329	0.394	0.418	0.388
DPPH	0.826	0.852	0.979	0.756	0.186	0.256	0.827	-0.275	-0.687	-0.549	-0.638	0.400	0.281	0.394	0.347
FRAP	0.809	0.658	0.835	0.962	0.260	0.068	0.569	-0.354	-0.531	-0.389	-0.518	0.188	0.555	0.198	0.158
ABTS	0.910	0.831	0.941	0.911	0.352	0.225	0.673	-0.359	-0.533	-0.350	-0.513	0.342	0.450	0.474	0.394
L*	-0.200	-0.406	-0.445	0.015	0.257	-0.300	-0.597	-0.009	0.696	0.708	0.611	-0.304	0.304	-0.079	-0.178
a*	0.086	0.284	0.388	-0.085	-0.383	0.218	0.643	0.038	-0.591	-0.629	-0.517	0.262	-0.472	-0.039	-0.007
b*	0.333	0.517	0.560	0.095	-0.229	0.253	0.644	-0.065	-0.756	-0.779	-0.683	0.281	-0.200	0.095	0.156
Quinic acid	0.450	0.570	0.703	0.314	-0.207	0.309	0.866	-0.159	-0.726	-0.702	-0.669	0.304	-0.145	0.144	0.127
Gallic acid	0.872	0.766	0.877	0.819	0.303	0.233	0.618	-0.400	-0.458	-0.332	-0.469	0.267	0.338	0.372	0.273
Chlorogenic acid	0.849	0.789	0.935	0.902	0.302	0.198	0.739	-0.327	-0.543	-0.377	-0.517	0.340	0.406	0.337	0.290
Total phenolic acids	0.704	0.745	0.892	0.597	-0.008	0.317	0.895	-0.284	-0.721	-0.644	-0.683	0.340	0.058	0.264	0.214
Gallocatechin	0.899	0.890	0.986	0.807	0.272	0.279	0.762	-0.343	-0.685	-0.510	-0.643	0.367	0.314	0.379	0.339
Epigallocatechin	0.902	0.823	0.934	0.919	0.360	0.228	0.678	-0.370	-0.515	-0.314	-0.496	0.333	0.445	0.429	0.344
Procyanidin B1	0.959	0.903	0.950	0.836	0.377	0.312	0.682	-0.420	-0.588	-0.404	-0.575	0.330	0.378	0.458	0.412
Catechin	0.850	0.912	0.991	0.685	0.211	0.414	0.881	-0.286	-0.706	-0.524	-0.647	0.470	0.152	0.486	0.418
Procyanidin B2	0.646	0.393	0.579	0.865	0.170	-0.245	0.217	-0.407	-0.392	-0.364	-0.435	-0.094	0.660	-0.016	0.020
Epicatechin	0.775	0.869	0.818	0.708	0.551	0.165	0.501	-0.198	-0.396	-0.219	-0.355	0.427	0.342	0.318	0.363
Epigallocatechin gallate	0.611	0.698	0.824	0.504	-0.044	0.309	0.895	-0.237	-0.700	-0.646	-0.657	0.345	-0.003	0.208	0.182
Procyanidin B1 3-O-gallate	0.649	0.640	0.661	0.680	0.385	0.272	0.491	-0.299	-0.446	-0.186	-0.405	0.273	0.394	0.181	0.204
Gallocatechin gallate	0.862	0.721	0.823	0.946	0.433	0.164	0.491	-0.343	-0.390	-0.182	-0.378	0.283	0.562	0.452	0.362
Quercetin 3-O-di-galactoside/glucoside	0.842	0.818	0.871	0.884	0.475	0.156	0.549	-0.303	-0.461	-0.252	-0.431	0.343	0.515	0.356	0.338
Myricetin 3-O-galactoside 1	0.784	0.704	0.639	0.368	0.283	0.565	0.471	-0.378	-0.435	-0.198	-0.419	0.314	0.099	0.754	0.601
Epicatechin gallate	0.836	0.889	0.854	0.715	0.487	0.281	0.561	-0.266	-0.481	-0.279	-0.441	0.440	0.319	0.442	0.413
Rutin	0.777	0.910	0.888	0.642	0.467	0.411	0.727	-0.190	-0.536	-0.279	-0.458	0.554	0.193	0.497	0.496
Quercetin 3-O-galactoside	0.582	0.628	0.599	0.382	0.394	0.296	0.392	-0.047	-0.381	-0.119	-0.289	0.460	0.192	0.495	0.469

Kaempferol 3-O-rutinoside 2	0.826	0.781	0.666	0.523	0.438	0.366	0.356	-0.393	-0.415	-0.218	-0.414	0.286	0.264	0.594	0.519
Kaempferol 3-O-glucoside	0.441	0.495	0.396	0.135	0.211	0.251	0.255	-0.209	-0.376	-0.216	-0.345	0.206	0.028	0.261	0.287
Unknown quercetin conjugate	1	0.882	0.889	0.771	0.369	0.359	0.582	-0.530	-0.574	-0.400	-0.591	0.244	0.400	0.560	0.484
Unknown kaempferol conjugate	0.882	1	0.911	0.591	0.446	0.450	0.725	-0.346	-0.555	-0.357	-0.525	0.465	0.178	0.591	0.553
Total flavonoids	0.889	0.911	1	0.771	0.261	0.363	0.836	-0.319	-0.680	-0.491	-0.631	0.436	0.247	0.467	0.402
Caffeine	0.771	0.591	0.771	1	0.345	0.003	0.448	-0.327	-0.397	-0.232	-0.390	0.166	0.642	0.204	0.171
Ethoxyacetic acid	0.369	0.446	0.261	0.345	1	0.254	0.014	0.121	0.379	0.508	0.390	0.492	0.357	0.317	0.460
Acetic acid	0.359	0.450	0.363	0.003	0.254	1	0.557	0.015	-0.242	-0.151	-0.187	0.605	-0.382	0.582	0.532
Pterin-6-carboxylic acid	0.582	0.725	0.836	0.448	0.014	0.557	1	-0.117	-0.662	-0.508	-0.574	0.562	-0.140	0.417	0.318
Hexanoic acid	-0.530	-0.346	-0.319	-0.327	0.121	0.015	-0.117	1	0.528	0.369	0.666	0.567	-0.233	-0.147	-0.046
Octanoic acid	-0.574	-0.555	-0.680	-0.397	0.379	-0.242	-0.662	0.528	1	0.896	0.980	0.017	-0.020	-0.196	-0.098
Nonanoic acid	-0.400	-0.357	-0.491	-0.232	0.508	-0.151	-0.508	0.369	0.896	1	0.896	0.069	-0.014	-0.067	-0.016
Total acid	-0.591	-0.525	-0.631	-0.390	0.390	-0.187	-0.574	0.666	0.980	0.896	1	0.163	-0.072	-0.178	-0.078
2-Methyl-1-propanol	0.244	0.465	0.436	0.166	0.492	0.605	0.562	0.567	0.017	0.069	0.163	1	-0.194	0.480	0.524
2-Pentanol	0.400	0.178	0.247	0.642	0.357	-0.382	-0.140	-0.233	-0.020	-0.014	-0.072	-0.194	1	-0.028	0.016
1-Butanol	0.560	0.591	0.467	0.204	0.317	0.582	0.417	-0.147	-0.196	-0.067	-0.178	0.480	-0.028	1	0.913
2-Hexanol	0.484	0.553	0.402	0.171	0.460	0.532	0.318	-0.046	-0.098	-0.016	-0.078	0.524	0.016	0.913	1
3-Methyl-1-butanol	-0.107	0.114	0.138	-0.084	0.220	0.546	0.393	0.754	0.100	0.061	0.260	0.899	-0.384	0.234	0.293
1-(2-Aminoethylamino)-2-propanol	0.774	0.791	0.895	0.878	0.307	0.183	0.703	-0.260	-0.532	-0.305	-0.479	0.355	0.419	0.335	0.264
2-Octen-4-ol	0.857	0.849	0.946	0.878	0.380	0.296	0.752	-0.261	-0.520	-0.312	-0.472	0.456	0.395	0.472	0.394
3-Methyl-1-pentanol	-0.225	0.005	-0.010	-0.206	0.268	0.419	0.216	0.805	0.194	0.203	0.364	0.809	-0.315	0.123	0.222
(E)-3-Hexen-1-ol	0.186	0.067	0.082	0.282	0.081	-0.221	-0.119	-0.198	-0.020	-0.016	-0.064	-0.196	0.335	-0.040	-0.015
1-Hexanol	0.219	0.204	0.116	0.392	0.535	-0.459	-0.312	0.082	0.191	0.133	0.174	-0.023	0.725	-0.143	0.022
2-Methyl-4-heptanol	0.286	0.228	0.116	0.238	0.381	-0.239	-0.190	-0.208	0.037	0.135	0.004	-0.108	0.398	-0.098	0.021
2-Ethyl-1-hexanol	0.038	0.198	0.317	-0.133	-0.396	0.558	0.732	0.220	-0.500	-0.530	-0.392	0.491	-0.534	0.207	0.131
Linalool	0.672	0.806	0.922	0.573	0.112	0.388	0.942	-0.154	-0.654	-0.516	-0.580	0.509	0.020	0.343	0.287
1-Octanol	-0.613	-0.502	-0.580	-0.526	0.161	0.190	-0.321	0.737	0.705	0.649	0.782	0.396	-0.389	-0.053	0.007
2,3-Butanediol	0.170	0.272	0.228	-0.062	0.159	0.879	0.456	0.165	-0.222	-0.179	-0.145	0.530	-0.335	0.282	0.205
Phenylethyl alcohol	-0.497	-0.305	-0.299	-0.440	0.046	0.388	0.061	0.843	0.391	0.302	0.525	0.652	-0.541	0.022	0.075
Total alcohol	-0.020	0.187	0.167	-0.064	0.320	0.598	0.379	0.710	0.138	0.120	0.287	0.931	-0.327	0.433	0.493
3-Hydroxybutanal	0.082	0.290	0.234	0.119	-0.023	0.013	0.229	0.042	-0.234	-0.093	-0.165	0.115	-0.042	0.034	-0.014
Octanal	0.337	0.293	0.241	0.270	0.597	0.562	0.249	-0.216	0.133	0.293	0.102	0.298	0.048	0.211	0.217
Total aldehyde	0.338	0.295	0.243	0.271	0.598	0.562	0.251	-0.214	0.133	0.293	0.102	0.301	0.049	0.211	0.217

Benzaldehyde	-0.349	-0.296	-0.232	-0.355	-0.671	-0.004	-0.013	0.129	-0.240	-0.280	-0.189	-0.157	-0.358	-0.115	-0.255
Myristicin	-0.655	-0.591	-0.466	-0.515	-0.537	0.041	-0.032	0.544	0.232	0.044	0.292	0.148	-0.573	-0.198	-0.271
Benzeneacetaldehyde	0.396	0.467	0.485	0.196	-0.073	0.289	0.443	-0.170	-0.584	-0.522	-0.539	0.310	-0.038	0.546	0.575
Naphthalene	0.786	0.819	0.944	0.677	0.069	0.321	0.838	-0.261	-0.788	-0.683	-0.729	0.379	0.224	0.379	0.338
1,2-Dihydro-1,1,6-trimethyl-naphthalene	0.786	0.819	0.944	0.677	0.069	0.321	0.838	-0.261	-0.788	-0.683	-0.729	0.379	0.224	0.379	0.338
Benzyl alcohol	0.132	0.094	0.024	0.280	0.237	-0.694	-0.396	-0.177	0.109	0.058	0.042	-0.373	0.555	-0.260	-0.155
Total benzenoid	0.072	0.054	0.004	0.226	0.093	-0.727	-0.384	-0.135	0.036	-0.043	-0.019	-0.375	0.492	-0.253	-0.161
Ethyl butanoate	0.330	0.435	0.256	0.385	0.855	-0.056	-0.049	0.155	0.314	0.375	0.326	0.369	0.498	0.203	0.354
Ethyl 3-methylbutanoate	-0.043	0.088	-0.032	0.095	0.791	0.145	-0.097	0.499	0.539	0.561	0.597	0.535	0.201	0.030	0.138
3-Methylbutyl acetate	0.372	0.361	0.197	0.417	0.758	-0.256	-0.235	-0.114	0.272	0.338	0.226	0.011	0.643	0.086	0.220
3-Hydroxymandelic acid ethyl ester	0.329	0.324	0.138	0.162	0.755	0.398	-0.051	-0.220	0.242	0.445	0.205	0.157	0.120	0.297	0.418
Ethyl hexanoate	-0.116	0.014	-0.179	0.006	0.737	-0.048	-0.258	0.424	0.714	0.690	0.723	0.393	0.184	0.076	0.224
Hexyl acetate	0.233	0.204	0.043	0.303	0.724	-0.320	-0.344	-0.089	0.431	0.477	0.367	-0.084	0.609	0.023	0.147
Ethyl (E)-3-hexenoate	0.333	0.376	0.322	0.262	0.572	0.402	0.152	0.326	0.060	0.091	0.138	0.552	0.137	0.330	0.482
Ethyl lactate	-0.249	-0.116	-0.119	-0.286	0.095	0.694	0.230	0.546	0.183	0.164	0.287	0.656	-0.547	0.233	0.230
Propyl hexanoate	-0.034	0.161	0.004	-0.010	0.744	0.364	0.088	0.481	0.544	0.647	0.615	0.687	-0.064	0.233	0.336
Ethyl octanoate	-0.356	-0.338	-0.520	-0.283	0.554	0.016	-0.505	0.233	0.877	0.874	0.825	0.056	-0.044	-0.048	0.031
Ethyl 3-hydroxybutanoate	-0.615	-0.653	-0.612	-0.414	-0.191	-0.097	-0.436	0.455	0.485	0.348	0.504	0.055	-0.273	-0.237	-0.211
Ethyl nonanoate	-0.091	0.059	-0.213	-0.091	0.795	-0.076	-0.406	0.264	0.651	0.644	0.633	0.198	0.213	-0.006	0.209
Isoamyl lactate	0.153	0.311	0.308	-0.077	0.051	0.878	0.618	0.246	-0.302	-0.200	-0.183	0.640	-0.458	0.426	0.391
Methyl decanoate	-0.121	0.016	-0.146	-0.007	0.576	0.059	-0.211	0.265	0.457	0.716	0.520	0.277	0.120	0.132	0.217
Ethyl 2-furancarboxylate	-0.512	-0.464	-0.396	-0.420	-0.367	0.178	-0.061	0.497	0.184	0.123	0.266	0.256	-0.519	-0.032	-0.086
Ethyl decanoate	-0.331	-0.376	-0.485	-0.210	0.408	-0.198	-0.528	0.082	0.816	0.834	0.740	-0.173	-0.006	-0.148	-0.150
Diethyl succinate	-0.529	-0.447	-0.334	-0.477	-0.507	0.164	0.069	0.493	0.091	-0.044	0.168	0.234	-0.627	-0.105	-0.183
Ethyl 9-decenoate	-0.393	-0.414	-0.476	-0.266	0.340	-0.037	-0.428	0.295	0.780	0.747	0.747	0.028	-0.170	-0.171	-0.202
Ethyl benzeneacetate	0.022	0.139	0.327	-0.047	-0.545	0.211	0.592	0.150	-0.620	-0.714	-0.528	0.219	-0.345	-0.008	-0.077
2-Phenethyl acetate	-0.130	-0.007	-0.104	-0.208	0.285	0.157	-0.045	0.222	0.266	0.292	0.292	0.221	-0.130	0.065	0.111
Ethyl dodecanoate	-0.063	0.018	-0.101	-0.015	0.273	-0.233	-0.164	0.181	0.390	0.519	0.413	0.077	0.081	0.115	0.216
Total ester	0.293	0.293	0.102	0.137	0.771	0.393	-0.077	-0.182	0.295	0.491	0.259	0.171	0.105	0.288	0.411
3-Hydroxy-2-butanone	-0.214	-0.206	-0.262	-0.147	0.153	-0.303	-0.273	0.100	0.575	0.613	0.536	-0.167	-0.096	-0.065	-0.069
Limonene	0.720	0.765	0.866	0.589	0.022	0.325	0.780	-0.236	-0.798	-0.727	-0.740	0.342	0.235	0.357	0.348
Terpinolene	0.631	0.606	0.746	0.898	0.259	-0.041	0.519	-0.218	-0.420	-0.234	-0.380	0.211	0.542	0.158	0.113
Total terpene	0.761	0.792	0.910	0.711	0.108	0.274	0.783	-0.237	-0.757	-0.658	-0.698	0.362	0.328	0.342	0.333

Values in **bold** indicate significant ($p < 0.05$) correlation

Table S6 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated Chardonnay wines

(continued 4/7)

Variables	3-Methyl-1-butanol	1-(2-Aminoethylamino)-2-propanol	2-Octen-4-ol	3-Methyl-1-pentanol	(E)-3-Hexen-1-ol	1-Hexanol	2-Methyl-4-heptanol	2-Ethyl-1-hexanol	Linalool	1-Octanol	2,3-Butanedio l	Phenylethy l alcohol	Total alcohol	3-Hydroxybutana l	Octanal
TA	-0.145	-0.680	-0.718	-0.133	-0.176	-0.418	-0.201	-0.201	-0.687	0.551	-0.147	0.220	-0.141	-0.134	0.101
pH	0.180	0.748	0.776	0.154	0.171	0.399	0.286	0.225	0.745	-0.541	0.113	-0.214	0.167	0.178	-0.089
TPC	0.105	0.873	0.916	-0.091	0.089	0.084	-0.003	0.372	0.944	-0.598	0.133	-0.305	0.114	0.192	0.232
TFC	0.011	0.902	0.950	-0.128	0.271	0.251	0.166	0.157	0.829	-0.607	0.130	-0.419	0.051	0.142	0.321
DPPH	0.120	0.866	0.910	-0.031	0.051	0.144	0.063	0.341	0.937	-0.621	0.139	-0.303	0.137	0.243	0.143
FRAP	-0.065	0.891	0.898	-0.198	0.307	0.331	0.183	0.035	0.670	-0.592	0.007	-0.442	-0.060	0.097	0.250
ABTS	0.023	0.925	0.974	-0.147	0.171	0.256	0.131	0.076	0.786	-0.575	0.075	-0.400	0.077	0.175	0.284
L*	-0.307	-0.207	-0.201	-0.293	0.264	0.189	-0.039	-0.691	-0.577	0.338	-0.325	-0.102	-0.248	-0.286	0.287
a*	0.288	0.124	0.138	0.219	-0.280	-0.393	-0.100	0.775	0.615	-0.286	0.234	0.172	0.197	0.169	-0.226
b*	0.230	0.308	0.315	0.185	-0.226	-0.086	0.037	0.645	0.670	-0.474	0.279	-0.027	0.172	0.275	-0.288
Quinic acid	0.187	0.514	0.539	0.058	-0.051	-0.226	-0.183	0.731	0.851	-0.475	0.240	-0.074	0.137	0.157	0.015
Gallic acid	-0.019	0.802	0.875	-0.254	0.143	0.150	0.118	0.057	0.759	-0.564	0.079	-0.391	0.008	0.056	0.371
Chlorogenic acid	0.042	0.921	0.956	-0.123	0.195	0.209	0.107	0.171	0.841	-0.560	0.057	-0.357	0.064	0.149	0.320
Total phenolic acids	0.128	0.732	0.778	-0.060	0.034	-0.084	-0.076	0.547	0.944	-0.585	0.201	-0.224	0.105	0.145	0.172
Gallocatechin	0.070	0.882	0.924	-0.055	0.102	0.213	0.211	0.234	0.877	-0.642	0.177	-0.366	0.088	0.222	0.195
Epigallocatechin	0.019	0.938	0.980	-0.149	0.207	0.239	0.144	0.067	0.785	-0.545	0.080	-0.396	0.062	0.160	0.342
Procyanidin B1	-0.006	0.876	0.932	-0.137	0.184	0.264	0.217	0.105	0.787	-0.595	0.146	-0.432	0.044	0.156	0.319
Catechin	0.187	0.852	0.908	0.042	0.034	0.040	0.062	0.411	0.950	-0.555	0.278	-0.240	0.213	0.250	0.216
Procyanidin B2	-0.293	0.650	0.635	-0.391	0.391	0.458	0.319	-0.203	0.371	-0.619	-0.299	-0.556	-0.291	-0.007	0.104
Epicatechin	0.124	0.791	0.807	0.043	0.067	0.495	0.438	-0.080	0.676	-0.461	0.086	-0.299	0.150	0.316	0.176
Epigallocatechin gallate	0.158	0.674	0.703	-0.022	-0.010	-0.116	-0.097	0.612	0.918	-0.532	0.200	-0.162	0.124	0.176	0.145
Procyanidin B1 3-O-gallate	0.073	0.718	0.699	0.088	0.146	0.225	0.572	0.018	0.509	-0.343	0.288	-0.249	0.072	0.186	0.312
Gallocatechin gallate	-0.017	0.891	0.929	-0.161	0.224	0.337	0.192	-0.117	0.598	-0.476	0.042	-0.404	0.049	0.136	0.323
Quercetin 3-O-di-galactoside/glucoside	0.044	0.893	0.911	-0.071	0.063	0.412	0.365	-0.081	0.696	-0.540	0.086	-0.384	0.078	0.259	0.230
Myricetin 3-O-galactoside 1	0.026	0.449	0.577	0.032	0.069	-0.093	0.179	0.158	0.440	-0.279	0.379	-0.224	0.157	-0.006	0.308
Epicatechin gallate	0.136	0.804	0.845	0.050	0.116	0.359	0.454	0.004	0.684	-0.482	0.220	-0.296	0.177	0.267	0.204
Rutin	0.262	0.809	0.855	0.212	0.029	0.205	0.338	0.207	0.791	-0.409	0.333	-0.159	0.306	0.267	0.243
Quercetin 3-O-galactoside	0.221	0.481	0.548	0.354	0.008	0.200	0.355	0.105	0.426	-0.197	0.266	-0.060	0.297	0.151	0.069

Kaempferol 3-O-rutinoside 2	-0.019	0.540	0.628	-0.041	0.041	0.240	0.550	-0.089	0.412	-0.439	0.282	-0.345	0.081	0.118	0.171
Kaempferol 3-O-glucoside	0.038	0.199	0.246	0.189	-0.023	0.065	0.730	0.082	0.275	-0.284	0.283	-0.124	0.076	0.121	-0.022
Unknown quercetin conjugate	-0.107	0.774	0.857	-0.225	0.186	0.219	0.286	0.038	0.672	-0.613	0.170	-0.497	-0.020	0.082	0.337
Unknown kaempferol conjugate	0.114	0.791	0.849	0.005	0.067	0.204	0.228	0.198	0.806	-0.502	0.272	-0.305	0.187	0.290	0.293
Total flavonoids	0.138	0.895	0.946	-0.010	0.082	0.116	0.116	0.317	0.922	-0.580	0.228	-0.299	0.167	0.234	0.241
Caffeine	-0.084	0.878	0.878	-0.206	0.282	0.392	0.238	-0.133	0.573	-0.526	-0.062	-0.440	-0.064	0.119	0.270
Ethoxyacetic acid	0.220	0.307	0.380	0.268	0.081	0.535	0.381	-0.396	0.112	0.161	0.159	0.046	0.320	-0.023	0.597
Acetic acid	0.546	0.183	0.296	0.419	-0.221	-0.459	-0.239	0.558	0.388	0.190	0.879	0.388	0.598	0.013	0.562
Pterin-6-carboxylic acid	0.393	0.703	0.752	0.216	-0.119	-0.312	-0.190	0.732	0.942	-0.321	0.456	0.061	0.379	0.229	0.249
Hexanoic acid	0.754	-0.260	-0.261	0.805	-0.198	0.082	-0.208	0.220	-0.154	0.737	0.165	0.843	0.710	0.042	-0.216
Octanoic acid	0.100	-0.532	-0.520	0.194	-0.020	0.191	0.037	-0.500	-0.654	0.705	-0.222	0.391	0.138	-0.234	0.133
Nonanoic acid	0.061	-0.305	-0.312	0.203	-0.016	0.133	0.135	-0.530	-0.516	0.649	-0.179	0.302	0.120	-0.093	0.293
Total acid	0.260	-0.479	-0.472	0.364	-0.064	0.174	0.004	-0.392	-0.580	0.782	-0.145	0.525	0.287	-0.165	0.102
2-Methyl-1-propanol	0.899	0.355	0.456	0.809	-0.196	-0.023	-0.108	0.491	0.509	0.396	0.530	0.652	0.931	0.115	0.298
2-Pentanol	-0.384	0.419	0.395	-0.315	0.335	0.725	0.398	-0.534	0.020	-0.389	-0.335	-0.541	-0.327	-0.042	0.048
1-Butanol	0.234	0.335	0.472	0.123	-0.040	-0.143	-0.098	0.207	0.343	-0.053	0.282	0.022	0.433	0.034	0.211
2-Hexanol	0.293	0.264	0.394	0.222	-0.015	0.022	0.021	0.131	0.287	0.007	0.205	0.075	0.493	-0.014	0.217
3-Methyl-1-butanol	1	0.085	0.153	0.897	-0.296	-0.191	-0.291	0.573	0.295	0.598	0.572	0.870	0.969	0.058	0.139
1-(2-Aminoethylamino)-2-propanol	0.085	1	0.960	-0.043	0.215	0.264	0.114	0.133	0.783	-0.480	0.086	-0.312	0.106	0.418	0.235
2-Octen-4-ol	0.153	0.960	1	-0.006	0.183	0.223	0.099	0.167	0.831	-0.472	0.170	-0.284	0.194	0.206	0.325
3-Methyl-1-pentanol	0.897	-0.043	-0.006	1	-0.173	-0.072	-0.075	0.455	0.110	0.672	0.501	0.853	0.877	0.100	0.065
(E)-3-Hexen-1-ol	-0.296	0.215	0.183	-0.173	1	0.381	0.233	-0.282	-0.062	-0.138	-0.277	-0.356	-0.262	-0.100	0.092
1-Hexanol	-0.191	0.264	0.223	-0.072	0.381	1	0.491	-0.639	-0.074	-0.215	-0.378	-0.364	-0.162	0.082	-0.138
2-Methyl-4-heptanol	-0.291	0.114	0.099	-0.075	0.233	0.491	1	-0.447	-0.088	-0.248	-0.176	-0.354	-0.263	0.014	-0.033
2-Ethyl-1-hexanol	0.573	0.133	0.167	0.455	-0.282	-0.639	-0.447	1	0.545	0.076	0.542	0.484	0.518	0.147	-0.027
Linalool	0.295	0.783	0.831	0.110	-0.062	-0.074	-0.088	0.545	1	-0.466	0.286	-0.097	0.274	0.229	0.203
1-Octanol	0.598	-0.480	-0.472	0.672	-0.138	-0.215	-0.248	0.076	-0.466	1	0.190	0.845	0.601	-0.090	0.172
2,3-Butanediol	0.572	0.086	0.170	0.501	-0.277	-0.378	-0.176	0.542	0.286	0.190	1	0.447	0.547	0.006	0.405
Phenylethyl alcohol	0.870	-0.312	-0.284	0.853	-0.356	-0.364	-0.354	0.484	-0.097	0.845	0.447	1	0.836	0.014	0.063
Total alcohol	0.969	0.106	0.194	0.877	-0.262	-0.162	-0.263	0.518	0.274	0.601	0.547	0.836	1	0.057	0.173
3-Hydroxybutanal	0.058	0.418	0.206	0.100	-0.100	0.082	0.014	0.147	0.229	-0.090	0.006	0.014	0.057	1	-0.217
Octanal	0.139	0.235	0.325	0.065	0.092	-0.138	-0.033	-0.027	0.203	0.172	0.405	0.063	0.173	-0.217	1
Total aldehyde	0.141	0.237	0.327	0.067	0.092	-0.136	-0.032	-0.027	0.204	0.172	0.406	0.064	0.175	-0.216	1.000

Benzaldehyde	0.061	-0.163	-0.317	0.079	-0.260	-0.419	-0.329	0.403	-0.150	0.099	0.091	0.214	0.008	0.589	-0.517
Myristicin	0.462	-0.454	-0.465	0.349	-0.346	-0.599	-0.625	0.502	-0.201	0.597	0.090	0.707	0.382	-0.039	-0.209
Benzeneacetaldehyde	0.216	0.327	0.378	0.160	0.022	-0.107	0.142	0.369	0.436	-0.326	0.172	-0.058	0.281	0.167	-0.184
Naphthalene	0.143	0.815	0.842	0.018	0.061	0.095	0.066	0.457	0.907	-0.642	0.249	-0.278	0.149	0.295	0.039
1,2-Dihydro-1,1,6-trimethylnaphthalene	0.143	0.815	0.842	0.018	0.061	0.095	0.066	0.457	0.907	-0.642	0.249	-0.278	0.149	0.295	0.039
Benzyl alcohol	-0.505	0.123	0.081	-0.440	0.282	0.834	0.489	-0.722	-0.134	-0.468	-0.603	-0.619	-0.506	-0.012	-0.356
Total benzenoid	-0.465	0.102	0.040	-0.413	0.238	0.780	0.439	-0.633	-0.134	-0.477	-0.624	-0.579	-0.475	0.082	-0.518
Ethyl butanoate	0.093	0.353	0.381	0.176	0.185	0.787	0.526	-0.478	0.102	-0.026	-0.073	-0.131	0.173	0.102	0.201
Ethyl 3-methylbutanoate	0.419	0.058	0.108	0.538	0.033	0.466	0.231	-0.251	-0.061	0.416	0.252	0.349	0.441	-0.009	0.389
3-Methylbutyl acetate	-0.274	0.291	0.303	-0.161	0.292	0.854	0.641	-0.711	-0.030	-0.260	-0.257	-0.449	-0.191	0.016	0.138
3-Hydroxymandelic acid ethyl ester	-0.048	0.102	0.181	0.045	0.172	0.174	0.422	-0.373	-0.027	0.046	0.284	-0.120	0.050	-0.183	0.673
Ethyl hexanoate	0.275	-0.033	0.003	0.311	0.040	0.485	0.201	-0.417	-0.203	0.491	-0.115	0.265	0.336	-0.057	0.308
Hexyl acetate	-0.345	0.150	0.166	-0.212	0.311	0.785	0.592	-0.762	-0.157	-0.175	-0.318	-0.426	-0.258	-0.075	0.167
Ethyl (E)-3-hexenoate	0.446	0.242	0.327	0.489	0.095	0.353	0.357	-0.008	0.200	0.067	0.441	0.200	0.498	-0.056	0.133
Ethyl lactate	0.828	-0.151	-0.093	0.709	-0.353	-0.519	-0.527	0.546	0.030	0.726	0.660	0.873	0.817	-0.010	0.373
Propyl hexanoate	0.574	0.087	0.142	0.617	-0.034	0.131	0.036	-0.029	0.043	0.658	0.251	0.537	0.627	0.008	0.583
Ethyl octanoate	0.050	-0.394	-0.352	0.089	-0.013	0.084	0.003	-0.489	-0.526	0.628	-0.064	0.274	0.106	-0.266	0.530
Ethyl 3-hydroxybutanoate	0.332	-0.500	-0.520	0.292	-0.205	-0.248	-0.368	-0.023	-0.513	0.739	-0.118	0.572	0.285	-0.100	-0.009
Ethyl nonanoate	0.060	-0.119	-0.112	0.180	0.027	0.620	0.323	-0.573	-0.288	0.312	-0.117	0.061	0.124	0.030	0.269
Isoamyl lactate	0.669	0.163	0.228	0.630	-0.236	-0.509	-0.268	0.749	0.407	0.250	0.874	0.543	0.676	0.100	0.285
Methyl decanoate	0.210	0.081	0.014	0.402	-0.046	0.236	0.211	-0.361	-0.235	0.520	-0.001	0.249	0.279	0.277	0.270
Ethyl 2-furancarboxylate	0.544	-0.330	-0.346	0.495	-0.115	-0.515	-0.527	0.401	-0.246	0.738	0.153	0.733	0.504	-0.031	-0.057
Ethyl decanoate	-0.205	-0.371	-0.336	-0.138	0.151	0.073	0.093	-0.577	-0.508	0.386	-0.214	0.054	-0.162	-0.301	0.410
Diethyl succinate	0.542	-0.379	-0.376	0.384	-0.522	-0.654	-0.631	0.565	-0.088	0.552	0.191	0.731	0.463	-0.014	-0.183
Ethyl 9-decenoate	0.081	-0.401	-0.365	0.081	-0.097	-0.088	-0.077	-0.345	-0.439	0.543	-0.003	0.365	0.092	-0.209	0.408
Ethyl benzeneacetate	0.321	0.128	0.137	0.242	-0.209	-0.383	-0.292	0.808	0.525	-0.288	0.358	0.196	0.228	0.158	-0.393
2-Phenethyl acetate	0.221	-0.203	-0.154	0.176	-0.896	-0.083	0.026	-0.028	-0.039	0.154	0.217	0.263	0.227	0.050	0.038
Ethyl dodecanoate	-0.015	0.046	0.003	0.080	0.047	0.227	0.303	-0.266	-0.165	0.243	-0.372	0.020	0.058	0.059	-0.116
Total ester	-0.025	0.074	0.154	0.065	0.154	0.171	0.406	-0.385	-0.056	0.096	0.279	-0.081	0.073	-0.192	0.686
3-Hydroxy-2-butanone	-0.229	-0.222	-0.200	-0.196	-0.054	0.024	0.209	-0.354	-0.246	0.098	-0.323	-0.071	-0.196	-0.180	-0.059
Limonene	0.134	0.745	0.750	0.046	0.052	0.114	0.069	0.482	0.829	-0.621	0.268	-0.265	0.143	0.357	-0.036
Terpinolene	-0.009	0.953	0.862	-0.113	0.303	0.380	0.158	-0.021	0.619	-0.458	-0.087	-0.348	-0.004	0.459	0.135

Total terpene	0.131	0.842	0.841	0.031	0.117	0.196	0.104	0.399	0.852	-0.622	0.215	-0.294	0.139	0.350	0.025
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Values in **bold** indicate significant ($p < 0.05$) correlation

Table S6 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated Chardonnay wines

(continued 5/7)

Variables	Total aldehyde	Benzaldehyde	Myristicin	Benzeneacetaldehyde	Naphthalene	1,2-Dihydro-1,1,6-trimethylnaphthalene	Benzyl alcohol	Total benzeneid	Ethylbutanoate	Ethyl 3-methylbutanoate	3-Methylbutyl acetate	3-Hydroxymandellic acid ethyl ester	Ethyl hexanoate	Hexyl acetate	Ethyl (E)-3-hexenoate
TA	0.099	0.253	0.426	-0.330	-0.815	-0.815	-0.298	-0.289	-0.387	-0.195	-0.342	0.016	0.116	-0.209	-0.464
pH	-0.087	-0.216	-0.438	0.416	0.869	0.869	0.273	0.276	0.380	0.139	0.333	-0.006	-0.130	0.186	0.502
TPC	0.233	-0.240	-0.354	0.396	0.928	0.928	0.021	0.002	0.176	-0.105	0.117	0.004	-0.188	-0.013	0.170
TFC	0.323	-0.371	-0.557	0.368	0.887	0.887	0.134	0.079	0.342	0.010	0.320	0.215	-0.087	0.180	0.318
DPPH	0.145	-0.199	-0.400	0.482	0.964	0.964	0.065	0.063	0.226	-0.058	0.169	0.008	-0.204	0.020	0.253
FRAP	0.251	-0.338	-0.502	0.214	0.772	0.772	0.231	0.178	0.299	0.001	0.326	0.104	-0.107	0.204	0.245
ABTS	0.285	-0.324	-0.513	0.374	0.842	0.842	0.160	0.121	0.361	0.038	0.342	0.167	-0.049	0.206	0.289
L*	0.285	-0.276	-0.002	-0.606	-0.636	-0.636	0.187	0.086	0.166	0.252	0.281	0.219	0.442	0.424	-0.202
a*	-0.224	0.229	0.225	0.423	0.547	0.547	-0.290	-0.203	-0.340	-0.316	-0.453	-0.340	-0.469	-0.529	-0.023
b*	-0.287	0.198	-0.088	0.561	0.734	0.734	-0.071	0.016	-0.110	-0.262	-0.196	-0.257	-0.449	-0.351	0.174
Quinic acid	0.017	0.004	-0.001	0.324	0.803	0.803	-0.229	-0.208	-0.168	-0.272	-0.282	-0.255	-0.388	-0.379	-0.009
Gallic acid	0.372	-0.362	-0.455	0.297	0.753	0.753	0.127	0.076	0.236	-0.048	0.275	0.199	-0.109	0.175	0.231
Chlorogenic acid	0.322	-0.366	-0.432	0.277	0.852	0.852	0.108	0.054	0.309	0.029	0.264	0.110	-0.049	0.141	0.211
Total phenolic acids	0.174	-0.161	-0.203	0.358	0.909	0.909	-0.105	-0.113	-0.003	-0.204	-0.074	-0.097	-0.313	-0.189	0.096
Gallocatechin	0.197	-0.262	-0.539	0.454	0.939	0.939	0.134	0.109	0.292	-0.028	0.272	0.149	-0.192	0.109	0.350
Epigallocatechin	0.344	-0.351	-0.513	0.322	0.815	0.815	0.139	0.087	0.349	0.038	0.331	0.194	-0.036	0.203	0.267
Procyanidin B1	0.321	-0.372	-0.595	0.331	0.852	0.852	0.156	0.094	0.376	0.022	0.359	0.247	-0.054	0.206	0.321
Catechin	0.217	-0.179	-0.408	0.520	0.954	0.954	-0.047	-0.055	0.199	-0.059	0.120	0.107	-0.219	-0.032	0.311
Procyanidin B2	0.105	-0.329	-0.500	0.167	0.553	0.553	0.432	0.396	0.249	-0.099	0.391	0.039	-0.123	0.301	0.126
Epicatechin	0.178	-0.385	-0.626	0.361	0.710	0.710	0.382	0.342	0.636	0.269	0.566	0.290	0.203	0.380	0.439
Epigallocatechin gallate	0.147	-0.122	-0.125	0.320	0.865	0.865	-0.140	-0.143	-0.023	-0.203	-0.126	-0.153	-0.279	-0.236	0.034
Procyanidin B1 3-O-gallate	0.314	-0.279	-0.571	0.352	0.575	0.575	0.072	0.019	0.389	0.151	0.351	0.396	0.005	0.207	0.416
Gallocatechin gallate	0.324	-0.344	-0.558	0.294	0.684	0.684	0.217	0.165	0.422	0.115	0.435	0.257	0.036	0.312	0.328
Quercetin 3-O-digalactoside/glucoside	0.232	-0.344	-0.616	0.382	0.764	0.764	0.303	0.265	0.530	0.178	0.503	0.256	0.076	0.336	0.387
Myricetin 3-O-galactoside 1	0.308	-0.155	-0.492	0.441	0.526	0.526	-0.198	-0.242	0.123	-0.055	0.143	0.419	-0.207	0.055	0.378
Epicatechin	0.206	-0.306	-0.635	0.521	0.751	0.751	0.258	0.236	0.532	0.198	0.487	0.333	0.053	0.309	0.498

gallate															
Rutin	0.245	-0.311	-0.559	0.538	0.799	0.799	0.069	0.038	0.459	0.196	0.349	0.353	0.013	0.177	0.508
Quercetin 3-O-galactoside	0.071	-0.131	-0.540	0.532	0.534	0.534	0.033	0.025	0.329	0.181	0.304	0.361	-0.108	0.174	0.565
Kaempferol 3-O-rutinoside 2	0.172	-0.267	-0.655	0.489	0.556	0.556	0.181	0.149	0.446	0.121	0.459	0.428	0.000	0.304	0.474
Kaempferol 3-O-glucoside	-0.021	-0.092	-0.475	0.564	0.390	0.390	0.034	0.043	0.226	0.055	0.239	0.354	-0.167	0.123	0.434
Unknown quercetin conjugate	0.338	-0.349	-0.655	0.396	0.786	0.786	0.132	0.072	0.330	-0.043	0.372	0.329	-0.116	0.233	0.333
Unknown kaempferol conjugate	0.295	-0.296	-0.591	0.467	0.819	0.819	0.094	0.054	0.435	0.088	0.361	0.324	0.014	0.204	0.376
Total flavonoids	0.243	-0.232	-0.466	0.485	0.944	0.944	0.024	0.004	0.256	-0.032	0.197	0.138	-0.179	0.043	0.322
Caffeine	0.271	-0.355	-0.515	0.196	0.677	0.677	0.280	0.226	0.385	0.095	0.417	0.162	0.006	0.303	0.262
Ethoxyacetic acid	0.598	-0.671	-0.537	-0.073	0.069	0.069	0.237	0.093	0.855	0.791	0.758	0.755	0.737	0.724	0.572
Acetic acid	0.562	-0.004	0.041	0.289	0.321	0.321	-0.694	-0.727	-0.056	0.145	-0.256	0.398	-0.048	-0.320	0.402
Pterin-6-carboxylic acid	0.251	-0.013	-0.032	0.443	0.838	0.838	-0.396	-0.384	-0.049	-0.097	-0.235	-0.051	-0.258	-0.344	0.152
Hexanoic acid	-0.214	0.129	0.544	-0.170	-0.261	-0.261	-0.177	-0.135	0.155	0.499	-0.114	-0.220	0.424	-0.089	0.326
Octanoic acid	0.133	-0.240	0.232	-0.584	-0.788	-0.788	0.109	0.036	0.314	0.539	0.272	0.242	0.714	0.431	0.060
Nonanoic acid	0.293	-0.280	0.044	-0.522	-0.683	-0.683	0.058	-0.043	0.375	0.561	0.338	0.445	0.690	0.477	0.091
Total acid	0.102	-0.189	0.292	-0.539	-0.729	-0.729	0.042	-0.019	0.326	0.597	0.226	0.205	0.723	0.367	0.138
2-Methyl-1-propanol	0.301	-0.157	0.148	0.310	0.379	0.379	-0.373	-0.375	0.369	0.535	0.011	0.157	0.393	-0.084	0.552
2-Pentanol	0.049	-0.358	-0.573	-0.038	0.224	0.224	0.555	0.492	0.498	0.201	0.643	0.120	0.184	0.609	0.137
1-Butanol	0.211	-0.115	-0.198	0.546	0.379	0.379	-0.260	-0.253	0.203	0.030	0.086	0.297	0.076	0.023	0.330
2-Hexanol	0.217	-0.255	-0.271	0.575	0.338	0.338	-0.155	-0.161	0.354	0.138	0.220	0.418	0.224	0.147	0.482
3-Methyl-1-butanol	0.141	0.061	0.462	0.216	0.143	0.143	-0.505	-0.465	0.093	0.419	-0.274	-0.048	0.275	-0.345	0.446
1-(2-Aminoethylamino)-2-propanol	0.237	-0.163	-0.454	0.327	0.815	0.815	0.123	0.102	0.353	0.058	0.291	0.102	-0.033	0.150	0.242
2-Octen-4-ol	0.327	-0.317	-0.465	0.378	0.842	0.842	0.081	0.040	0.381	0.108	0.303	0.181	0.003	0.166	0.327
3-Methyl-1-pentanol	0.067	0.079	0.349	0.160	0.018	0.018	-0.440	-0.413	0.176	0.538	-0.161	0.045	0.311	-0.212	0.489
(E)-3-Hexen-1-ol	0.092	-0.260	-0.346	0.022	0.061	0.061	0.282	0.238	0.185	0.033	0.292	0.172	0.040	0.311	0.095
1-Hexanol	-0.136	-0.419	-0.599	-0.107	0.095	0.095	0.834	0.780	0.787	0.466	0.854	0.174	0.485	0.785	0.353
2-Methyl-4-heptanol	-0.032	-0.329	-0.625	0.142	0.066	0.066	0.489	0.439	0.526	0.231	0.641	0.422	0.201	0.592	0.357
2-Ethyl-1-hexanol	-0.027	0.403	0.502	0.369	0.457	0.457	-0.722	-0.633	-0.478	-0.251	-0.711	-0.373	-0.417	-0.762	-0.008
Linalool	0.204	-0.150	-0.201	0.436	0.907	0.907	-0.134	-0.134	0.102	-0.061	-0.030	-0.027	-0.203	-0.157	0.200
1-Octanol	0.172	0.099	0.597	-0.326	-0.642	-0.642	-0.468	-0.477	-0.026	0.416	-0.260	0.046	0.491	-0.175	0.067
2,3-Butanediol	0.406	0.091	0.090	0.172	0.249	0.249	-0.603	-0.624	-0.073	0.252	-0.257	0.284	-0.115	-0.318	0.441

Phenylethyl alcohol	0.064	0.214	0.707	-0.058	-0.278	-0.278	-0.619	-0.579	-0.131	0.349	-0.449	-0.120	0.265	-0.426	0.200
Total alcohol	0.175	0.008	0.382	0.281	0.149	0.149	-0.506	-0.475	0.173	0.441	-0.191	0.050	0.336	-0.258	0.498
3-Hydroxybutanal	-0.216	0.589	-0.039	0.167	0.295	0.295	-0.012	0.082	0.102	-0.009	0.016	-0.183	-0.057	-0.075	-0.056
Octanal	1.000	-0.517	-0.209	-0.184	0.039	0.039	-0.356	-0.518	0.201	0.389	0.138	0.673	0.308	0.167	0.133
Total aldehyde	1	-0.518	-0.210	-0.183	0.041	0.041	-0.356	-0.518	0.203	0.391	0.139	0.673	0.309	0.167	0.134
Benzaldehyde	-0.518	1	0.476	0.195	-0.038	-0.038	-0.351	-0.168	-0.564	-0.433	-0.583	-0.569	-0.501	-0.594	-0.282
Myristicin	-0.210	0.476	1	-0.194	-0.375	-0.375	-0.591	-0.492	-0.576	-0.140	-0.789	-0.641	-0.046	-0.716	-0.430
Benzeneacetaldehyde	-0.183	0.195	-0.194	1	0.580	0.580	-0.109	0.021	-0.012	-0.228	-0.059	0.020	-0.348	-0.188	0.350
Naphthalene	0.041	-0.038	-0.375	0.580	1	1.000	0.008	0.037	0.148	-0.121	0.089	-0.040	-0.327	-0.072	0.315
1,2-Dihydro-1,1,6-trimethylnaphthalene	0.041	-0.038	-0.375	0.580	1.000	1	0.008	0.037	0.148	-0.121	0.089	-0.040	-0.327	-0.072	0.315
Benzyl alcohol	-0.356	-0.351	-0.591	-0.109	0.008	0.008	1	0.974	0.568	0.137	0.767	0.052	0.255	0.735	0.110
Total benzenoid	-0.518	-0.168	-0.492	0.021	0.037	0.037	0.974	1	0.479	0.043	0.665	-0.099	0.161	0.619	0.076
Ethyl butanoate	0.203	-0.564	-0.576	-0.012	0.148	0.148	0.568	0.479	1	0.772	0.904	0.486	0.775	0.843	0.544
Ethyl 3-methylbutanoate	0.391	-0.433	-0.140	-0.228	-0.121	-0.121	0.137	0.043	0.772	1	0.560	0.436	0.807	0.560	0.478
3-Methylbutyl acetate	0.139	-0.583	-0.789	-0.059	0.089	0.089	0.767	0.665	0.904	0.560	1	0.556	0.570	0.969	0.473
3-Hydroxymandelic acid ethyl ester	0.673	-0.569	-0.641	0.020	-0.040	-0.040	0.052	-0.099	0.486	0.436	0.556	1	0.328	0.576	0.568
Ethyl hexanoate	0.309	-0.501	-0.046	-0.348	-0.327	-0.327	0.255	0.161	0.775	0.807	0.570	0.328	1	0.594	0.227
Hexyl acetate	0.167	-0.594	-0.716	-0.188	-0.072	-0.072	0.735	0.619	0.843	0.560	0.969	0.576	0.594	1	0.398
Ethyl (E)-3-hexenoate	0.134	-0.282	-0.430	0.350	0.315	0.315	0.110	0.076	0.544	0.478	0.473	0.568	0.227	0.398	1
Ethyl lactate	0.373	0.182	0.606	0.015	-0.144	-0.144	-0.793	-0.779	-0.220	0.256	-0.524	0.040	0.152	-0.529	0.168
Propyl hexanoate	0.584	-0.412	0.062	-0.216	-0.171	-0.171	-0.222	-0.326	0.569	0.787	0.272	0.448	0.828	0.284	0.308
Ethyl octanoate	0.530	-0.403	0.084	-0.568	-0.708	-0.708	-0.018	-0.144	0.344	0.572	0.294	0.492	0.748	0.434	-0.006
Ethyl 3-hydroxybutanoate	-0.010	0.274	0.735	-0.264	-0.628	-0.628	-0.344	-0.287	-0.279	0.083	-0.432	-0.344	0.274	-0.382	-0.355
Ethyl nonanoate	0.270	-0.459	-0.355	-0.306	-0.325	-0.325	0.439	0.337	0.806	0.747	0.742	0.543	0.851	0.763	0.356
Isoamyl lactate	0.286	0.190	0.180	0.338	0.351	0.351	-0.737	-0.725	-0.169	0.097	-0.404	0.201	-0.190	-0.464	0.445
Methyl decanoate	0.271	-0.111	-0.174	-0.149	-0.311	-0.311	-0.004	-0.064	0.480	0.533	0.354	0.452	0.586	0.349	0.223
Ethyl 2-furancarboxylate	-0.057	0.382	0.825	-0.095	-0.395	-0.395	-0.645	-0.573	-0.508	-0.113	-0.732	-0.420	-0.007	-0.713	-0.291
Ethyl decanoate	0.409	-0.403	-0.015	-0.560	-0.665	-0.665	0.132	0.012	0.248	0.441	0.330	0.461	0.529	0.513	-0.072
Diethyl succinate	-0.183	0.475	0.932	-0.089	-0.278	-0.278	-0.640	-0.541	-0.634	-0.242	-0.837	-0.589	-0.163	-0.811	-0.369
Ethyl 9-decenoate	0.408	-0.256	0.264	-0.576	-0.637	-0.637	-0.115	-0.211	0.088	0.455	0.077	0.286	0.459	0.227	-0.102
Ethyl benzenoacetate	-0.393	0.449	0.310	0.463	0.559	0.559	-0.314	-0.184	-0.437	-0.328	-0.527	-0.498	-0.623	-0.589	0.063
2-Phenethyl	0.038	0.002	0.052	-0.093	-0.137	-0.137	-0.034	-0.046	0.186	0.273	0.117	0.162	0.244	0.120	0.152

acetate																
Ethyl dodecanoate	-0.116	-0.140	-0.200	-0.099	-0.209	-0.209	0.256	0.224	0.349	0.071	0.354	0.241	0.389	0.399	0.226	
Total ester	0.686	-0.576	-0.605	-0.015	-0.084	-0.084	0.043	-0.111	0.495	0.467	0.553	0.998	0.371	0.581	0.557	
3-Hydroxy-2-butanone	-0.060	-0.210	-0.077	-0.329	-0.355	-0.355	0.290	0.236	0.222	0.134	0.324	0.263	0.310	0.482	0.109	
Limonene	-0.034	0.080	-0.367	0.618	0.975	0.975	-0.005	0.045	0.132	-0.141	0.074	-0.068	-0.352	-0.089	0.335	
Terpinolene	0.136	-0.113	-0.412	0.240	0.690	0.690	0.249	0.242	0.373	0.077	0.342	0.014	0.017	0.220	0.166	
Total terpene	0.027	-0.016	-0.412	0.573	0.982	0.982	0.061	0.093	0.221	-0.070	0.159	-0.032	-0.267	-0.007	0.344	

Values in **bold** indicate significant ($p < 0.05$) correlation

Table S6 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated Chardonnay wines

(continued 6/7)

Variables	Ethyl lactate	Propyl hexanoate	Ethyl octanoate	Ethyl 3-hydroxybutanoate	Ethyl nonanoate	Isoamyl lactate	Methyl decanoate	Ethyl 2-furancarboxylate	Ethyl decanoate	Diethyl succinate	Ethyl 9-decanoate	Ethyl benzeneacetate	2-Phenethyl acetate	Ethyl dodecanoate	Total ester
TA	0.256	0.105	0.571	0.660	0.112	-0.174	0.253	0.470	0.439	0.391	0.413	-0.470	0.126	0.173	0.049
pH	-0.258	-0.089	-0.607	-0.621	-0.139	0.178	-0.210	-0.453	-0.499	-0.383	-0.481	0.459	-0.129	-0.078	-0.042
TPC	-0.137	-0.036	-0.514	-0.574	-0.264	0.226	-0.278	-0.367	-0.470	-0.253	-0.457	0.373	-0.153	-0.124	-0.029
TFC	-0.214	0.020	-0.420	-0.619	-0.125	0.183	-0.147	-0.461	-0.387	-0.479	-0.444	0.165	-0.245	-0.076	0.179
DPPH	-0.171	-0.064	-0.578	-0.590	-0.244	0.230	-0.236	-0.410	-0.524	-0.291	-0.496	0.412	-0.094	-0.162	-0.028
FRAP	-0.275	-0.078	-0.403	-0.526	-0.185	0.021	-0.157	-0.421	-0.325	-0.449	-0.374	0.105	-0.283	-0.041	0.072
ABTS	-0.202	0.024	-0.373	-0.531	-0.128	0.105	-0.081	-0.431	-0.319	-0.421	-0.349	0.105	-0.136	-0.057	0.138
L*	-0.075	0.253	0.724	0.350	0.335	-0.491	0.341	0.082	0.783	-0.127	0.656	-0.785	-0.083	0.222	0.255
a*	0.102	-0.255	-0.629	-0.274	-0.464	0.405	-0.519	0.012	-0.578	0.308	-0.409	0.843	0.045	-0.286	-0.364
b*	-0.040	-0.305	-0.777	-0.455	-0.338	0.402	-0.434	-0.195	-0.799	0.046	-0.678	0.785	0.051	-0.269	-0.295
Quinic acid	0.006	-0.167	-0.648	-0.487	-0.453	0.419	-0.494	-0.144	-0.630	0.057	-0.549	0.716	-0.157	-0.247	-0.288
Gallic acid	-0.189	-0.035	-0.285	-0.483	-0.162	0.032	-0.248	-0.437	-0.173	-0.346	-0.165	0.092	-0.116	-0.144	0.175
Chlorogenic acid	-0.186	0.051	-0.380	-0.526	-0.167	0.119	-0.142	-0.394	-0.329	-0.368	-0.346	0.168	-0.193	-0.071	0.082
Total phenolic acids	-0.081	-0.124	-0.590	-0.561	-0.390	0.316	-0.446	-0.293	-0.533	-0.119	-0.476	0.549	-0.170	-0.229	-0.132
Galocatechin	-0.211	-0.060	-0.546	-0.649	-0.184	0.227	-0.174	-0.473	-0.491	-0.403	-0.486	0.307	-0.101	-0.106	0.110
Epigallocatechin	-0.190	0.060	-0.336	-0.517	-0.129	0.106	-0.048	-0.406	-0.285	-0.422	-0.324	0.063	-0.173	-0.026	0.166
Procyanidin B1	-0.217	0.041	-0.387	-0.608	-0.073	0.172	-0.088	-0.473	-0.382	-0.482	-0.427	0.090	-0.147	-0.036	0.213
Catechin	-0.068	0.002	-0.552	-0.610	-0.245	0.381	-0.171	-0.356	-0.526	-0.271	-0.500	0.408	-0.083	-0.124	0.070
Procyanidin B2	-0.470	-0.266	-0.334	-0.390	-0.132	-0.328	-0.281	-0.467	-0.202	-0.481	-0.276	-0.024	-0.334	-0.048	0.011
Epicatechin	-0.228	0.200	-0.268	-0.506	0.234	0.048	0.148	-0.490	-0.305	-0.491	-0.319	-0.013	0.062	0.057	0.267
Epigallocatechin gallate	-0.044	-0.091	-0.572	-0.524	-0.368	0.336	-0.443	-0.244	-0.553	-0.055	-0.488	0.576	-0.146	-0.200	-0.184
Procyanidin B1 3-O-gallate	-0.131	0.153	-0.280	-0.464	-0.024	0.271	0.264	-0.336	-0.327	-0.468	-0.393	-0.022	-0.062	0.117	0.365
Galocatechin gallate	-0.196	0.070	-0.233	-0.437	-0.046	0.026	0.078	-0.397	-0.185	-0.470	-0.241	-0.096	-0.135	0.032	0.234
Quercetin 3-O-digalactoside/glucoside	-0.248	0.093	-0.322	-0.505	0.051	0.048	0.114	-0.499	-0.317	-0.493	-0.344	-0.002	0.040	0.012	0.229
Myricetin 3-O-galactoside 1	0.011	0.053	-0.249	-0.478	-0.153	0.432	0.080	-0.238	-0.240	-0.334	-0.285	0.029	-0.032	0.020	0.389
Epicatechin gallate	-0.181	0.120	-0.340	-0.561	0.072	0.172	0.085	-0.466	-0.317	-0.490	-0.337	0.073	-0.006	0.006	0.305
Rutin	-0.043	0.211	-0.379	-0.612	0.015	0.398	0.162	-0.378	-0.400	-0.399	-0.406	0.198	0.043	0.051	0.322
Quercetin 3-O-galactoside	-0.024	0.130	-0.337	-0.465	0.020	0.371	0.306	-0.253	-0.329	-0.350	-0.343	0.133	0.101	0.159	0.332

Kaempferol 3-O-rutinoside 2	-0.197	0.037	-0.262	-0.533	0.054	0.190	0.090	-0.453	-0.248	-0.503	-0.284	-0.059	0.080	0.005	0.399
Kaempferol 3-O-glucoside	-0.176	-0.038	-0.349	-0.452	-0.041	0.263	0.066	-0.330	-0.287	-0.344	-0.283	0.175	0.106	-0.027	0.323
Unknown quercetin conjugate	-0.249	-0.034	-0.356	-0.615	-0.091	0.153	-0.121	-0.512	-0.331	-0.529	-0.393	0.022	-0.130	-0.063	0.293
Unknown kaempferol conjugate	-0.116	0.161	-0.338	-0.653	0.059	0.311	0.016	-0.464	-0.376	-0.447	-0.414	0.139	-0.007	0.018	0.293
Total flavonoids	-0.119	0.004	-0.520	-0.612	-0.213	0.308	-0.146	-0.396	-0.485	-0.334	-0.476	0.327	-0.104	-0.101	0.102
Caffeine	-0.286	-0.010	-0.283	-0.414	-0.091	-0.077	-0.007	-0.420	-0.210	-0.477	-0.266	-0.047	-0.208	-0.015	0.137
Ethoxyacetic acid	0.095	0.744	0.554	-0.191	0.795	0.051	0.576	-0.367	0.408	-0.507	0.340	-0.545	0.285	0.273	0.771
Acetic acid	0.694	0.364	0.016	-0.097	-0.076	0.878	0.059	0.178	-0.198	0.164	-0.037	0.211	0.157	-0.233	0.393
Pterin-6-carboxylic acid	0.230	0.088	-0.505	-0.436	-0.406	0.618	-0.211	-0.061	-0.528	0.069	-0.428	0.592	-0.045	-0.164	-0.077
Hexanoic acid	0.546	0.481	0.233	0.455	0.264	0.246	0.265	0.497	0.082	0.493	0.295	0.150	0.222	0.181	-0.182
Octanoic acid	0.183	0.544	0.877	0.485	0.651	-0.302	0.457	0.184	0.816	0.091	0.780	-0.620	0.266	0.390	0.295
Nonanoic acid	0.164	0.647	0.874	0.348	0.644	-0.200	0.716	0.123	0.834	-0.044	0.747	-0.714	0.292	0.519	0.491
Total acid	0.287	0.615	0.825	0.504	0.633	-0.183	0.520	0.266	0.740	0.168	0.747	-0.528	0.292	0.413	0.259
2-Methyl-1-propanol	0.656	0.687	0.056	0.055	0.198	0.640	0.277	0.256	-0.173	0.234	0.028	0.219	0.221	0.077	0.171
2-Pentanol	-0.547	-0.064	-0.044	-0.273	0.213	-0.458	0.120	-0.519	-0.006	-0.627	-0.170	-0.345	-0.130	0.081	0.105
1-Butanol	0.233	0.233	-0.048	-0.237	-0.006	0.426	0.132	-0.032	-0.148	-0.105	-0.171	-0.008	0.065	0.115	0.288
2-Hexanol	0.230	0.336	0.031	-0.211	0.209	0.391	0.217	-0.086	-0.150	-0.183	-0.202	-0.077	0.111	0.216	0.411
3-Methyl-1-butanol	0.828	0.574	0.050	0.332	0.060	0.669	0.210	0.544	-0.205	0.542	0.081	0.321	0.221	-0.015	-0.025
1-(2-Aminoethylamino)-2-propanol	-0.151	0.087	-0.394	-0.500	-0.119	0.163	0.081	-0.330	-0.371	-0.379	-0.401	0.128	-0.203	0.046	0.074
2-Octen-4-ol	-0.093	0.142	-0.352	-0.520	-0.112	0.228	0.014	-0.346	-0.336	-0.376	-0.365	0.137	-0.154	0.003	0.154
3-Methyl-1-pentanol	0.709	0.617	0.089	0.292	0.180	0.630	0.402	0.495	-0.138	0.384	0.081	0.242	0.176	0.080	0.065
(E)-3-Hexen-1-ol	-0.353	-0.034	-0.013	-0.205	0.027	-0.236	-0.046	-0.115	0.151	-0.522	-0.097	-0.209	-0.896	0.047	0.154
1-Hexanol	-0.519	0.131	0.084	-0.248	0.620	-0.509	0.236	-0.515	0.073	-0.654	-0.088	-0.383	-0.083	0.227	0.171
2-Methyl-4-heptanol	-0.527	0.036	0.003	-0.368	0.323	-0.268	0.211	-0.527	0.093	-0.631	-0.077	-0.292	0.026	0.303	0.406
2-Ethyl-1-hexanol	0.546	-0.029	-0.489	-0.023	-0.573	0.749	-0.361	0.401	-0.577	0.565	-0.345	0.808	-0.028	-0.266	-0.385
Linalool	0.030	0.043	-0.526	-0.513	-0.288	0.407	-0.235	-0.246	-0.508	-0.088	-0.439	0.525	-0.039	-0.165	-0.056
1-Octanol	0.726	0.658	0.628	0.739	0.312	0.250	0.520	0.738	0.386	0.552	0.543	-0.288	0.154	0.243	0.096
2,3-Butanediol	0.660	0.251	-0.064	-0.118	-0.117	0.874	-0.001	0.153	-0.214	0.191	-0.003	0.358	0.217	-0.372	0.279
Phenylethyl alcohol	0.873	0.537	0.274	0.572	0.061	0.543	0.249	0.733	0.054	0.731	0.365	0.196	0.263	0.020	-0.081
Total alcohol	0.817	0.627	0.106	0.285	0.124	0.676	0.279	0.504	-0.162	0.463	0.092	0.228	0.227	0.058	0.073
3-Hydroxybutanal	-0.010	0.008	-0.266	-0.100	0.030	0.100	0.277	-0.031	-0.301	-0.014	-0.209	0.158	0.050	0.059	-0.192
Octanal	0.373	0.583	0.530	-0.009	0.269	0.285	0.270	-0.057	0.410	-0.183	0.408	-0.393	0.038	-0.116	0.686

Total aldehyde	0.373	0.584	0.530	-0.010	0.270	0.286	0.271	-0.057	0.409	-0.183	0.408	-0.393	0.038	-0.116	0.686
Benzaldehyde	0.182	-0.412	-0.403	0.274	-0.459	0.190	-0.111	0.382	-0.403	0.475	-0.256	0.449	0.002	-0.140	-0.576
Myristicin	0.606	0.062	0.084	0.735	-0.355	0.180	-0.174	0.825	-0.015	0.932	0.264	0.310	0.052	-0.200	-0.605
Benzeneacetaldehyde	0.015	-0.216	-0.568	-0.264	-0.306	0.338	-0.149	-0.095	-0.560	-0.089	-0.576	0.463	-0.093	-0.099	-0.015
Naphthalene	-0.144	-0.171	-0.708	-0.628	-0.325	0.351	-0.311	-0.395	-0.665	-0.278	-0.637	0.559	-0.137	-0.209	-0.084
1,2-Dihydro-1,1,6-trimethylnaphthalene	-0.144	-0.171	-0.708	-0.628	-0.325	0.351	-0.311	-0.395	-0.665	-0.278	-0.637	0.559	-0.137	-0.209	-0.084
Benzyl alcohol	-0.793	-0.222	-0.018	-0.344	0.439	-0.737	-0.004	-0.645	0.132	-0.640	-0.115	-0.314	-0.034	0.256	0.043
Total benzenoid	-0.779	-0.326	-0.144	-0.287	0.337	-0.725	-0.064	-0.573	0.012	-0.541	-0.211	-0.184	-0.046	0.224	-0.111
Ethyl butanoate	-0.220	0.569	0.344	-0.279	0.806	-0.169	0.480	-0.508	0.248	-0.634	0.088	-0.437	0.186	0.349	0.495
Ethyl 3-methylbutanoate	0.256	0.787	0.572	0.083	0.747	0.097	0.533	-0.113	0.441	-0.242	0.455	-0.328	0.273	0.071	0.467
3-Methylbutyl acetate	-0.524	0.272	0.294	-0.432	0.742	-0.404	0.354	-0.732	0.330	-0.837	0.077	-0.527	0.117	0.354	0.553
3-Hydroxymandelic acid ethyl ester	0.040	0.448	0.492	-0.344	0.543	0.201	0.452	-0.420	0.461	-0.589	0.286	-0.498	0.162	0.241	0.998
Ethyl hexanoate	0.152	0.828	0.748	0.274	0.851	-0.190	0.586	-0.007	0.529	-0.163	0.459	-0.623	0.244	0.389	0.371
Hexyl acetate	-0.529	0.284	0.434	-0.382	0.763	-0.464	0.349	-0.713	0.513	-0.811	0.227	-0.589	0.120	0.399	0.581
Ethyl (E)-3-hexenoate	0.168	0.308	-0.006	-0.355	0.356	0.445	0.223	-0.291	-0.072	-0.369	-0.102	0.063	0.152	0.226	0.557
Ethyl lactate	1	0.527	0.273	0.523	-0.032	0.709	0.221	0.718	-0.007	0.683	0.301	0.142	0.216	-0.154	0.071
Propyl hexanoate	0.527	1	0.682	0.274	0.654	0.276	0.689	0.224	0.409	0.027	0.444	-0.435	0.230	0.313	0.487
Ethyl octanoate	0.273	0.682	1	0.425	0.703	-0.196	0.526	0.120	0.873	-0.016	0.808	-0.765	0.267	0.286	0.542
Ethyl 3-hydroxybutanoate	0.523	0.274	0.425	1	0.046	-0.131	0.239	0.789	0.204	0.687	0.396	-0.266	0.071	-0.072	-0.299
Ethyl nonanoate	-0.032	0.654	0.703	0.046	1	-0.236	0.593	-0.304	0.522	-0.414	0.385	-0.642	0.348	0.395	0.572
Isoamyl lactate	0.709	0.276	-0.196	-0.131	-0.236	1	0.072	0.292	-0.367	0.284	-0.195	0.485	0.117	-0.125	0.190
Methyl decanoate	0.221	0.689	0.526	0.239	0.593	0.072	1	0.144	0.304	-0.148	0.227	-0.613	0.262	0.532	0.475
Ethyl 2-furancarboxylate	0.718	0.224	0.120	0.789	-0.304	0.292	0.144	1	-0.072	0.841	0.186	0.070	-0.161	-0.070	-0.389
Ethyl decanoate	-0.007	0.409	0.873	0.204	0.522	-0.367	0.304	-0.072	1	-0.156	0.879	-0.651	0.120	0.208	0.502
Diethyl succinate	0.683	0.027	-0.016	0.687	-0.414	0.284	-0.148	0.841	-0.156	1	0.196	0.345	0.193	-0.195	-0.559
Ethyl 9-decenoate	0.301	0.444	0.808	0.396	0.385	-0.195	0.227	0.186	0.879	0.196	1	-0.473	0.255	-0.035	0.337
Ethyl benzeneacetate	0.142	-0.435	-0.765	-0.266	-0.642	0.485	-0.613	0.070	-0.651	0.345	-0.473	1	-0.054	-0.412	-0.527
2-Phenethyl acetate	0.216	0.230	0.267	0.071	0.348	0.117	0.262	-0.161	0.120	0.193	0.255	-0.054	1	0.139	0.184
Ethyl dodecanoate	-0.154	0.313	0.286	-0.072	0.395	-0.125	0.532	-0.070	0.208	-0.195	-0.035	-0.412	0.139	1	0.251
Total ester	0.071	0.487	0.542	-0.299	0.572	0.190	0.475	-0.389	0.502	-0.559	0.337	-0.527	0.184	0.251	1
3-Hydroxy-2-butanone	-0.269	0.123	0.449	-0.148	0.326	-0.273	0.132	-0.302	0.637	-0.191	0.407	-0.283	0.283	0.619	0.282
Limonene	-0.141	-0.209	-0.742	-0.610	-0.301	0.383	-0.297	-0.393	-0.742	-0.285	-0.728	0.590	-0.131	-0.195	-0.115

Terpinolene	-0.247	0.030	-0.345	-0.396	-0.085	-0.032	0.101	-0.317	-0.290	-0.397	-0.338	0.046	-0.271	0.063	-0.009
Total terpene	-0.168	-0.143	-0.687	-0.608	-0.256	0.319	-0.232	-0.410	-0.678	-0.339	-0.680	0.502	-0.167	-0.152	-0.076

Values in **bold** indicate significant ($p < 0.05$) correlation

Table S6 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated Chardonnay wines

(continued 7/7)

Variables	3-Hydroxy-2-butanone	Limonene	Terpinolene	Total terpene
TA	0.195	-0.768	-0.626	-0.809
pH	-0.185	0.823	0.693	0.868
TPC	-0.247	0.847	0.746	0.895
TFC	-0.239	0.811	0.794	0.878
DPPH	-0.280	0.893	0.742	0.932
FRAP	-0.203	0.692	0.873	0.794
ABTS	-0.183	0.746	0.838	0.830
L*	0.405	-0.700	-0.083	-0.607
a*	-0.227	0.549	0.001	0.465
b*	-0.417	0.777	0.171	0.694
Quinic acid	-0.305	0.780	0.369	0.749
Gallic acid	-0.094	0.629	0.704	0.704
Chlorogenic acid	-0.187	0.751	0.843	0.839
Total phenolic acids	-0.265	0.841	0.589	0.854
Gallocatechin	-0.259	0.865	0.755	0.912
Epigallocatechin	-0.170	0.711	0.853	0.805
Procyanidin B1	-0.224	0.774	0.747	0.835
Catechin	-0.279	0.884	0.681	0.910
Procyanidin B2	-0.155	0.497	0.727	0.594
Epicatechin	-0.173	0.638	0.707	0.712
Epigallocatechin gallate	-0.269	0.813	0.537	0.817
Procyanidin B1 3-O-gallate	-0.218	0.543	0.666	0.618
Gallocatechin gallate	-0.133	0.588	0.849	0.698
Quercetin 3-O-di-galactoside/glucoside	-0.183	0.684	0.836	0.777
Myricetin 3-O-galactoside 1	-0.150	0.486	0.228	0.469
Epicatechin gallate	-0.191	0.678	0.699	0.742
Rutin	-0.214	0.733	0.658	0.782
Quercetin 3-O-galactoside	-0.185	0.514	0.330	0.521

Kaempferol 3-O-rutinoside 2	-0.137	0.507	0.403	0.528
Kaempferol 3-O-glucoside	-0.140	0.390	0.086	0.355
Unknown quercetin conjugate	-0.214	0.720	0.631	0.761
Unknown kaempferol conjugate	-0.206	0.765	0.606	0.792
Total flavonoids	-0.262	0.866	0.746	0.910
Caffeine	-0.147	0.589	0.898	0.711
Ethoxyacetic acid	0.153	0.022	0.259	0.108
Acetic acid	-0.303	0.325	-0.041	0.274
Pterin-6-carboxylic acid	-0.273	0.780	0.519	0.783
Hexanoic acid	0.100	-0.236	-0.218	-0.237
Octanoic acid	0.575	-0.798	-0.420	-0.757
Nonanoic acid	0.613	-0.727	-0.234	-0.658
Total acid	0.536	-0.740	-0.380	-0.698
2-Methyl-1-propanol	-0.167	0.342	0.211	0.362
2-Pentanol	-0.096	0.235	0.542	0.328
1-Butanol	-0.065	0.357	0.158	0.342
2-Hexanol	-0.069	0.348	0.113	0.333
3-Methyl-1-butanol	-0.229	0.134	-0.009	0.131
1-(2-Aminoethylamino)-2-propanol	-0.222	0.745	0.953	0.842
2-Octen-4-ol	-0.200	0.750	0.862	0.841
3-Methyl-1-pentanol	-0.196	0.046	-0.113	0.031
(E)-3-Hexen-1-ol	-0.054	0.052	0.303	0.117
1-Hexanol	0.024	0.114	0.380	0.196
2-Methyl-4-heptanol	0.209	0.069	0.158	0.104
2-Ethyl-1-hexanol	-0.354	0.482	-0.021	0.399
Linalool	-0.246	0.829	0.619	0.852
1-Octanol	0.098	-0.621	-0.458	-0.622
2,3-Butanediol	-0.323	0.268	-0.087	0.215
Phenylethyl alcohol	-0.071	-0.265	-0.348	-0.294
Total alcohol	-0.196	0.143	-0.004	0.139
3-Hydroxybutanal	-0.180	0.357	0.459	0.350
Octanal	-0.059	-0.036	0.135	0.025
Total aldehyde	-0.060	-0.034	0.136	0.027

Benzaldehyde	-0.210	0.080	-0.113	-0.016
Myristicin	-0.077	-0.367	-0.412	-0.412
Benzeneacetaldehyde	-0.329	0.618	0.240	0.573
Naphthalene	-0.355	0.975	0.690	0.982
1,2-Dihydro-1,1,6-trimethyl-naphthalene	-0.355	0.975	0.690	0.982
Benzyl alcohol	0.290	-0.005	0.249	0.061
Total benzenoid	0.236	0.045	0.242	0.093
Ethyl butanoate	0.222	0.132	0.373	0.221
Ethyl 3-methylbutanoate	0.134	-0.141	0.077	-0.070
3-Methylbutyl acetate	0.324	0.074	0.342	0.159
3-Hydroxymandelic acid ethyl ester	0.263	-0.068	0.014	-0.032
Ethyl hexanoate	0.310	-0.352	0.017	-0.267
Hexyl acetate	0.482	-0.089	0.220	-0.007
Ethyl (E)-3-hexenoate	0.109	0.335	0.166	0.344
Ethyl lactate	-0.269	-0.141	-0.247	-0.168
Propyl hexanoate	0.123	-0.209	0.030	-0.143
Ethyl octanoate	0.449	-0.742	-0.345	-0.687
Ethyl 3-hydroxybutanoate	-0.148	-0.610	-0.396	-0.608
Ethyl nonanoate	0.326	-0.301	-0.085	-0.256
Isoamyl lactate	-0.273	0.383	-0.032	0.319
Methyl decanoate	0.132	-0.297	0.101	-0.232
Ethyl 2-furancarboxylate	-0.302	-0.393	-0.317	-0.410
Ethyl decanoate	0.637	-0.742	-0.290	-0.678
Diethyl succinate	-0.191	-0.285	-0.397	-0.339
Ethyl 9-decenoate	0.407	-0.728	-0.338	-0.680
Ethyl benzeneacetate	-0.283	0.590	0.046	0.502
2-Phenethyl acetate	0.283	-0.131	-0.271	-0.167
Ethyl dodecanoate	0.619	-0.195	0.063	-0.152
Total ester	0.282	-0.115	-0.009	-0.076
3-Hydroxy-2-butanone	1	-0.408	-0.187	-0.381
Limonene	-0.408	1	0.633	0.984
Terpinolene	-0.187	0.633	1	0.751
Total terpene	-0.381	0.984	0.751	1

Values in **bold** indicate significant ($p < 0.05$) correlation

Table S7 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated model wines

(continued 1/7)

Variables	TA	pH	TPC	TFC	DPPH	FRAP	ABTS	L*	a*	b*	Quinic acid	Gallic acid	Chlorogenic acid	Total phenolic acids	Gallocatechin
TA	1	-0.986	-0.567	-0.596	-0.595	-0.489	-0.517	0.682	-0.492	-0.847	-0.476	-0.490	-0.499	-0.719	-0.712
pH	-0.986	1	0.558	0.577	0.566	0.466	0.499	-0.714	0.550	0.863	0.500	0.483	0.468	0.724	0.703
TPC	-0.567	0.558	1	0.763	0.893	0.929	0.933	-0.269	0.170	0.410	0.350	0.815	0.854	0.940	0.947
TFC	-0.596	0.577	0.763	1	0.769	0.758	0.785	-0.236	0.136	0.436	0.288	0.712	0.823	0.822	0.801
DPPH	-0.595	0.566	0.893	0.769	1	0.967	0.981	-0.162	0.010	0.343	0.392	0.701	0.968	0.894	0.867
FRAP	-0.489	0.466	0.929	0.758	0.967	1	0.979	-0.100	-0.021	0.257	0.418	0.671	0.948	0.884	0.884
ABTS	-0.517	0.499	0.933	0.785	0.981	0.979	1	-0.123	0.003	0.288	0.337	0.758	0.966	0.905	0.880
L*	0.682	-0.714	-0.269	-0.236	-0.162	-0.100	-0.123	1	-0.959	-0.942	-0.582	-0.205	-0.014	-0.500	-0.492
a*	-0.492	0.550	0.170	0.136	0.010	-0.021	0.003	-0.959	1	0.849	0.530	0.135	-0.127	0.397	0.369
b*	-0.847	0.863	0.410	0.436	0.343	0.257	0.288	-0.942	0.849	1	0.593	0.333	0.201	0.628	0.622
Quinic acid	-0.476	0.500	0.350	0.288	0.392	0.418	0.337	-0.582	0.530	0.593	1	-0.147	0.265	0.511	0.519
Gallic acid	-0.490	0.483	0.815	0.712	0.701	0.671	0.758	-0.205	0.135	0.333	-0.147	1	0.721	0.771	0.736
Chlorogenic acid	-0.499	0.468	0.854	0.823	0.968	0.948	0.966	-0.014	-0.127	0.201	0.265	0.721	1	0.839	0.808
Total phenolic acids	-0.719	0.724	0.940	0.822	0.894	0.884	0.905	-0.500	0.397	0.628	0.511	0.771	0.839	1	0.973
Gallocatechin	-0.712	0.703	0.947	0.801	0.867	0.884	0.880	-0.492	0.369	0.622	0.519	0.736	0.808	0.973	1
Epigallocatechin	-0.500	0.477	0.938	0.820	0.961	0.981	0.977	-0.085	-0.037	0.262	0.384	0.705	0.959	0.891	0.878
Procyanidin B1	-0.719	0.688	0.912	0.816	0.954	0.911	0.933	-0.279	0.108	0.468	0.337	0.789	0.932	0.924	0.911
Catechin	-0.727	0.696	0.908	0.773	0.910	0.883	0.884	-0.443	0.293	0.592	0.453	0.739	0.843	0.942	0.957
Procyanidin B2	-0.472	0.476	0.824	0.784	0.806	0.868	0.853	-0.117	0.020	0.256	0.484	0.543	0.842	0.811	0.807
Epicatechin	-0.516	0.513	0.789	0.831	0.808	0.842	0.824	-0.271	0.180	0.445	0.629	0.467	0.781	0.830	0.847
Epigallocatechin gallate	-0.524	0.513	0.674	0.521	0.608	0.694	0.608	-0.489	0.398	0.530	0.758	0.267	0.532	0.720	0.813
Procyanidin B1 3-O-gallate	-0.632	0.650	0.831	0.767	0.811	0.828	0.842	-0.507	0.427	0.609	0.585	0.626	0.758	0.922	0.923
Gallocatechin gallate	-0.492	0.458	0.887	0.827	0.965	0.970	0.965	-0.031	-0.104	0.224	0.380	0.655	0.976	0.853	0.838
Quercetin 3-O-di-galactoside/glucoside	-0.471	0.490	0.832	0.763	0.720	0.796	0.809	-0.236	0.176	0.355	0.344	0.678	0.739	0.818	0.825
Myricetin 3-O-galactoside 1	-0.247	0.192	0.487	0.314	0.546	0.468	0.497	-0.081	0.002	0.171	-0.105	0.568	0.476	0.434	0.428
Epicatechin gallate	-0.706	0.699	0.850	0.789	0.887	0.869	0.884	-0.396	0.252	0.540	0.422	0.719	0.858	0.911	0.913
Rutin	-0.562	0.555	0.631	0.554	0.726	0.743	0.678	-0.428	0.298	0.518	0.850	0.191	0.636	0.731	0.749
Quercetin 3-O-galactoside	-0.503	0.485	0.503	0.414	0.531	0.497	0.500	-0.476	0.358	0.496	0.365	0.422	0.461	0.598	0.576
Kaempferol 3-O-rutinoside 2	-0.647	0.646	0.774	0.722	0.806	0.783	0.809	-0.400	0.279	0.528	0.370	0.679	0.774	0.838	0.852

Kaempferol 3-O-glucoside	-0.530	0.528	0.512	0.429	0.600	0.625	0.555	-0.487	0.365	0.538	0.894	0.051	0.498	0.631	0.673
Unknown quercetin conjugate	-0.596	0.558	0.816	0.763	0.832	0.769	0.828	-0.176	0.034	0.343	0.014	0.894	0.859	0.804	0.775
Unknown kaempferol conjugate	-0.641	0.601	0.826	0.757	0.920	0.903	0.875	-0.225	0.052	0.399	0.545	0.538	0.902	0.851	0.853
Total flavonoids	-0.694	0.668	0.942	0.813	0.941	0.931	0.930	-0.381	0.237	0.538	0.469	0.743	0.890	0.961	0.972
Caffeine	-0.453	0.451	0.866	0.745	0.869	0.919	0.910	-0.055	-0.053	0.200	0.431	0.592	0.893	0.824	0.813
Ethoxyacetic acid	-0.409	0.394	-0.274	-0.156	-0.159	-0.282	-0.246	-0.078	-0.055	0.185	-0.125	-0.158	-0.213	-0.222	-0.215
Acetic acid	-0.813	0.735	0.392	0.374	0.379	0.269	0.289	-0.513	0.316	0.666	0.199	0.386	0.272	0.447	0.480
Pterin-6-carboxylic acid	-0.655	0.650	0.362	0.476	0.247	0.193	0.243	-0.343	0.242	0.451	0.045	0.464	0.281	0.417	0.412
Hexanoic acid	-0.343	0.415	-0.126	-0.089	-0.157	-0.194	-0.212	-0.467	0.493	0.404	0.197	-0.040	-0.188	0.059	-0.011
Octanoic acid	-0.068	0.062	-0.016	0.126	-0.024	0.024	-0.051	0.193	-0.222	-0.184	-0.176	0.067	0.101	-0.039	-0.068
Nonanoic acid	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
Total acid	-0.282	0.332	-0.108	-0.018	-0.132	-0.140	-0.186	-0.255	0.260	0.214	0.062	-0.007	-0.102	0.018	-0.047
2-Methyl-1-propanol	-0.756	0.748	0.438	0.421	0.521	0.402	0.459	-0.410	0.262	0.545	0.221	0.477	0.484	0.558	0.507
2-Pentanol	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
1-Butanol	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
2-Hexanol	-0.244	0.252	0.153	0.186	0.369	0.238	0.282	0.052	-0.136	0.113	-0.094	0.274	0.330	0.198	0.165
3-Methyl-1-butanol	-0.767	0.723	0.350	0.475	0.387	0.327	0.295	-0.328	0.147	0.477	0.203	0.340	0.391	0.430	0.423
1-(2-Aminoethylamino)-2-propanol	-0.589	0.510	0.161	0.246	0.178	0.109	0.127	-0.427	0.286	0.528	0.189	0.146	0.109	0.236	0.301
2-Octen-4-ol	-0.512	0.523	0.882	0.706	0.784	0.827	0.838	-0.072	-0.021	0.220	0.171	0.771	0.796	0.794	0.772
3-Methyl-1-pentanol	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
(E)-3-Hexen-1-ol	-0.414	0.408	0.062	0.033	0.066	-0.006	-0.002	-0.546	0.468	0.491	0.265	0.071	-0.031	0.205	0.180
1-Hexanol	-0.485	0.528	0.085	0.130	0.100	0.033	0.052	-0.219	0.141	0.397	0.156	0.086	0.059	0.165	0.180
2-Methyl-4-heptanol	-0.204	0.272	-0.105	-0.082	-0.107	-0.153	-0.132	-0.100	0.081	0.206	0.042	-0.071	-0.129	-0.046	-0.030
2-Ethyl-1-hexanol	-0.388	0.410	0.368	0.132	0.308	0.299	0.276	-0.365	0.287	0.373	0.416	0.156	0.212	0.392	0.385
Linalool	-0.490	0.526	0.877	0.617	0.752	0.807	0.816	-0.380	0.352	0.467	0.451	0.675	0.674	0.869	0.844
1-Octanol	-0.486	0.506	0.870	0.668	0.756	0.803	0.809	-0.261	0.213	0.399	0.422	0.640	0.688	0.828	0.794
2,3-Butanediol	-0.875	0.839	0.428	0.509	0.426	0.355	0.364	-0.540	0.359	0.678	0.407	0.331	0.370	0.541	0.560
Phenylethyl alcohol	-0.378	0.440	-0.170	-0.145	-0.202	-0.294	-0.265	-0.805	0.825	0.655	0.348	-0.095	-0.305	0.091	0.032
Total alcohol	-0.534	0.608	0.084	0.062	0.091	-0.020	0.023	-0.742	0.720	0.705	0.357	0.146	-0.023	0.317	0.256
3-Hydroxybutanal	-0.515	0.459	0.042	0.089	-0.002	-0.026	-0.084	-0.620	0.512	0.645	0.563	-0.186	-0.102	0.176	0.257
Octanal	-0.322	0.288	-0.232	-0.006	-0.220	-0.312	-0.267	-0.019	-0.094	0.116	-0.272	-0.048	-0.200	-0.223	-0.181
Total aldehyde	-0.322	0.289	-0.232	-0.006	-0.220	-0.312	-0.267	-0.019	-0.094	0.116	-0.272	-0.048	-0.200	-0.223	-0.181
Benzaldehyde	-0.430	0.487	-0.174	-0.133	-0.195	-0.303	-0.265	-0.853	0.862	0.736	0.391	-0.114	-0.316	0.100	0.051
Myristicin	-0.415	0.354	0.124	0.259	0.119	0.099	0.081	-0.153	0.035	0.284	0.118	0.061	0.093	0.129	0.147
Benzeneacetaldehyde	-0.373	0.399	-0.128	-0.162	-0.118	-0.221	-0.195	-0.461	0.388	0.468	0.141	-0.060	-0.222	0.007	0.004
Naphthalene	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604

1,2-Dihydro-1,1,6-trimethyl-naphthalene	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
Benzyl alcohol	-0.362	0.424	-0.221	-0.171	-0.238	-0.340	-0.299	-0.601	0.585	0.567	0.230	-0.144	-0.332	-0.020	-0.032
Total benzenoid	-0.370	0.432	-0.220	-0.174	-0.236	-0.341	-0.299	-0.628	0.613	0.587	0.246	-0.144	-0.335	-0.011	-0.026
Ethyl butanoate	-0.340	0.346	0.023	0.176	0.248	0.132	0.168	0.211	-0.339	-0.051	-0.259	0.217	0.296	0.053	0.006
Ethyl 3-methylbutanoate	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
3-Methylbutyl acetate	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
3-Hydroxymandelic acid ethyl ester	-0.022	-0.020	-0.363	-0.434	-0.269	-0.371	-0.361	0.009	-0.089	-0.011	-0.224	-0.250	-0.356	-0.370	-0.373
Ethyl hexanoate	-0.314	0.352	0.688	0.464	0.418	0.461	0.528	-0.219	0.237	0.306	0.001	0.714	0.381	0.595	0.557
Hexyl acetate	-0.563	0.571	0.461	0.504	0.317	0.375	0.319	-0.285	0.230	0.418	0.403	0.255	0.311	0.473	0.517
Ethyl (E)-3-hexenoate	-0.218	0.199	0.139	0.067	0.313	0.300	0.226	-0.244	0.168	0.278	0.527	-0.133	0.193	0.235	0.233
Ethyl lactate	-0.635	0.615	0.170	0.191	0.108	0.028	0.034	-0.641	0.542	0.655	0.158	0.250	-0.003	0.283	0.309
Propyl hexanoate	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
Ethyl octanoate	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
Ethyl 3-hydroxybutanoate	-0.635	0.623	0.103	0.060	0.052	0.020	-0.021	-0.527	0.422	0.514	0.195	0.121	-0.017	0.204	0.241
Ethyl nonanoate	-0.873	0.844	0.335	0.425	0.333	0.204	0.246	-0.576	0.399	0.728	0.189	0.411	0.257	0.458	0.459
Isoamyl lactate	-0.878	0.842	0.341	0.420	0.364	0.260	0.286	-0.660	0.462	0.758	0.347	0.322	0.286	0.487	0.513
Methyl decanoate	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
Ethyl 2-furancarboxylate	-0.923	0.876	0.506	0.561	0.580	0.511	0.489	-0.548	0.345	0.732	0.574	0.307	0.492	0.637	0.653
Ethyl decanoate	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
Diethyl succinate	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
Ethyl 9-decenoate	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
Ethyl benzeneacetate	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
2-Phenethyl acetate	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
Ethyl dodecanoate	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
Total ester	-0.031	-0.011	-0.355	-0.427	-0.254	-0.356	-0.349	0.000	-0.083	0.000	-0.203	-0.251	-0.344	-0.357	-0.361
3-Hydroxy-2-butanone	-0.956	0.901	0.470	0.530	0.524	0.411	0.424	-0.571	0.354	0.758	0.394	0.396	0.439	0.592	0.604
Limonene	-0.931	0.893	0.714	0.703	0.821	0.713	0.740	-0.519	0.307	0.711	0.469	0.594	0.740	0.828	0.808
Terpinolene	-0.496	0.422	-0.042	0.027	-0.084	-0.155	-0.137	-0.565	0.458	0.552	0.104	0.039	-0.168	0.068	0.121
Total terpene	-0.922	0.886	0.722	0.716	0.834	0.730	0.756	-0.497	0.286	0.691	0.468	0.600	0.761	0.834	0.811

Values in **bold** indicate significant ($p < 0.05$) correlation

Table S7 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated model wines

(continued 2/7)

Variables	Epigallocatechin	Procyanidin B1	Catechin	Procyanidin B2	Epicatechin	Epigallocatechin gallate	Procyanidin B1 3-O-gallate	Gallocatechin gallate	Quercetin 3-O-di-galactoside/gluco side	Myricetin 3-O-galactoside 1	Epicatechin gallate	Rutin	Quercetin 3-O-galactoside	Kaempferol 3-O-rutinoside 2	Kaempferol 3-O-glucoside
TA	-0.500	-0.719	-0.727	-0.472	-0.516	-0.524	-0.632	-0.492	-0.471	-0.247	-0.706	-0.562	-0.503	-0.647	-0.530
pH	0.477	0.688	0.696	0.476	0.513	0.513	0.650	0.458	0.490	0.192	0.699	0.555	0.485	0.646	0.528
TPC	0.938	0.912	0.908	0.824	0.789	0.674	0.831	0.887	0.832	0.487	0.850	0.631	0.503	0.774	0.512
TFC	0.820	0.816	0.773	0.784	0.831	0.521	0.767	0.827	0.763	0.314	0.789	0.554	0.414	0.722	0.429
DPPH	0.961	0.954	0.910	0.806	0.808	0.608	0.811	0.965	0.720	0.546	0.887	0.726	0.531	0.806	0.600
FRAP	0.981	0.911	0.883	0.868	0.842	0.694	0.828	0.970	0.796	0.468	0.869	0.743	0.497	0.783	0.625
ABTS	0.977	0.933	0.884	0.853	0.824	0.608	0.842	0.965	0.809	0.497	0.884	0.678	0.500	0.809	0.555
L*	-0.085	-0.279	-0.443	-0.117	-0.271	-0.489	-0.507	-0.031	-0.236	-0.081	-0.396	-0.428	-0.476	-0.400	-0.487
a*	-0.037	0.108	0.293	0.020	0.180	0.398	0.427	-0.104	0.176	0.002	0.252	0.298	0.358	0.279	0.365
b*	0.262	0.468	0.592	0.256	0.445	0.530	0.609	0.224	0.355	0.171	0.540	0.518	0.496	0.528	0.538
Quinic acid	0.384	0.337	0.453	0.484	0.629	0.758	0.585	0.380	0.344	-0.105	0.422	0.850	0.365	0.370	0.894
Gallic acid	0.705	0.789	0.739	0.543	0.467	0.267	0.626	0.655	0.678	0.568	0.719	0.191	0.422	0.679	0.051
Chlorogenic acid	0.959	0.932	0.843	0.842	0.781	0.532	0.758	0.976	0.739	0.476	0.858	0.636	0.461	0.774	0.498
Total phenolic acids	0.891	0.924	0.942	0.811	0.830	0.720	0.922	0.853	0.818	0.434	0.911	0.731	0.598	0.838	0.631
Gallocatechin	0.878	0.911	0.957	0.807	0.847	0.813	0.923	0.838	0.825	0.428	0.913	0.749	0.576	0.852	0.673
Epigallocatechin	1	0.926	0.869	0.890	0.861	0.626	0.801	0.989	0.803	0.448	0.847	0.713	0.486	0.749	0.575
Procyanidin B1	0.926	1	0.951	0.775	0.765	0.602	0.797	0.920	0.720	0.555	0.920	0.670	0.555	0.833	0.541
Catechin	0.869	0.951	1	0.694	0.780	0.741	0.849	0.855	0.672	0.619	0.906	0.721	0.580	0.845	0.618
Procyanidin B2	0.890	0.775	0.694	1	0.840	0.685	0.801	0.872	0.900	0.017	0.764	0.713	0.428	0.661	0.640
Epicatechin	0.861	0.765	0.780	0.840	1	0.745	0.885	0.854	0.803	0.242	0.814	0.835	0.427	0.756	0.767
Epigallocatechin gallate	0.626	0.602	0.741	0.685	0.745	1	0.803	0.607	0.634	0.128	0.686	0.816	0.434	0.649	0.841
Procyanidin B1 3-O-gallate	0.801	0.797	0.849	0.801	0.885	0.803	1	0.767	0.869	0.270	0.914	0.784	0.557	0.877	0.751

Gallocatechin gallate	0.989	0.920	0.855	0.872	0.854	0.607	0.767	1	0.746	0.457	0.828	0.71	0.449	0.735	0.572
Quercetin 3-O-digallocatechin/gallate	0.803	0.720	0.672	0.900	0.803	0.634	0.869	0.746	1	0.050	0.809	0.59	0.470	0.739	0.550
Myricetin 3-O-galactoside 1	0.448	0.555	0.619	0.017	0.242	0.128	0.270	0.457	0.050	1	0.447	0.12	0.160	0.496	0.002
Epicatechin gallate	0.847	0.920	0.906	0.764	0.814	0.686	0.914	0.828	0.809	0.447	1	0.72	0.548	0.955	0.657
Rutin	0.713	0.670	0.721	0.713	0.835	0.816	0.784	0.711	0.599	0.129	0.725	1	0.657	0.599	0.970
Quercetin 3-O-galactoside	0.486	0.555	0.580	0.428	0.427	0.434	0.557	0.449	0.470	0.160	0.548	0.65	1	0.329	0.577
Kaempferol 3-O-rutinoside 2	0.749	0.833	0.845	0.661	0.756	0.649	0.877	0.735	0.739	0.496	0.955	0.59	0.329	1	0.570
Kaempferol 3-O-glucoside	0.575	0.541	0.618	0.640	0.767	0.841	0.751	0.572	0.550	0.002	0.657	0.97	0.577	0.570	1
Unknown quercetin conjugate	0.797	0.915	0.843	0.612	0.543	0.368	0.627	0.795	0.602	0.653	0.809	0.34	0.374	0.780	0.208
Unknown kaempferol conjugate	0.908	0.928	0.892	0.816	0.803	0.711	0.746	0.929	0.640	0.420	0.840	0.79	0.486	0.747	0.692
Total flavonoids	0.925	0.965	0.990	0.782	0.837	0.754	0.878	0.909	0.747	0.558	0.921	0.75	0.574	0.850	0.644
Caffeine	0.934	0.823	0.731	0.983	0.814	0.648	0.782	0.916	0.872	0.133	0.779	0.70	0.441	0.674	0.610
Ethoxyacetic acid	-0.273	-0.090	-0.181	-0.237	-0.236	-0.287	-0.223	-0.252	-0.214	-0.128	-0.064	-	-0.197	-0.022	-0.129
Acetic acid	0.295	0.532	0.551	0.156	0.233	0.281	0.279	0.294	0.160	0.406	0.425	0.18	0.289	0.401	0.220
Pterin-6-carboxylic acid	0.267	0.453	0.395	0.332	0.185	0.228	0.273	0.261	0.291	0.011	0.345	0.07	0.138	0.288	0.025
Hexanoic acid	-0.223	-0.012	0.005	-0.141	-0.201	0.014	0.007	-0.247	-0.122	-0.182	0.064	0.04	0.187	-0.002	0.049
Octanoic acid	0.034	0.077	-0.055	0.114	-0.184	-0.130	-0.202	0.043	0.016	-0.120	0.032	-	-0.020	-0.053	-0.197
Nonanoic acid	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
Total acid	-0.158	0.018	-0.029	-0.062	-0.234	-0.055	-0.091	-0.170	-0.089	-0.189	0.055	-	0.120	-0.029	-0.055
2-Methyl-1-propanol	0.398	0.628	0.555	0.388	0.272	0.287	0.446	0.409	0.333	0.240	0.593	0.28	0.191	0.599	0.252
2-Pentanol	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
1-Butanol	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
2-Hexanol	0.212	0.357	0.339	-0.053	0.170	-0.067	0.164	0.256	-0.069	0.568	0.336	0.10	0.019	0.382	0.017
3-Methyl-1-butanol	0.346	0.547	0.470	0.348	0.211	0.267	0.244	0.362	0.260	0.138	0.468	0.29	0.311	0.374	0.226
1-(2-Aminoethylamino)-2-propanol	0.094	0.258	0.323	0.071	0.174	0.315	0.249	0.119	0.147	0.199	0.286	0.12	0.018	0.374	0.207

2-Octen-4-ol	0.853	0.798	0.692	0.839	0.643	0.446	0.673	0.798	0.835	0.252	0.740	0.45	0.381	0.646	0.331
3-Methyl-1-pentanol	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
(E)-3-Hexen-1-ol	-0.005	0.159	0.217	-0.028	-0.005	0.156	0.144	-0.033	-0.021	-0.043	0.124	0.34	0.777	-0.083	0.308
1-Hexanol	0.039	0.231	0.190	0.053	0.188	0.099	0.194	0.041	0.103	-0.030	0.293	0.15	-0.040	0.302	0.168
2-Methyl-4-heptanol	-0.148	0.002	-0.039	-0.113	0.033	-0.068	0.028	-0.155	-0.039	-0.120	0.110	0.00	-0.189	0.157	0.041
2-Ethyl-1-hexanol	0.309	0.343	0.340	0.331	0.253	0.312	0.281	0.259	0.247	0.002	0.281	0.45	0.447	0.168	0.412
Linalool	0.793	0.734	0.749	0.727	0.774	0.610	0.867	0.712	0.854	0.322	0.810	0.65	0.525	0.740	0.579
1-Octanol	0.832	0.726	0.694	0.767	0.789	0.504	0.761	0.758	0.830	0.289	0.723	0.62	0.460	0.636	0.515
2,3-Butanediol	0.380	0.573	0.547	0.440	0.403	0.468	0.471	0.381	0.386	0.047	0.552	0.43	0.315	0.492	0.433
Phenylethyl alcohol	-0.311	-0.073	0.041	-0.264	-0.169	0.079	0.085	-0.343	-0.199	-0.078	0.045	0.11	0.274	0.049	0.178
Total alcohol	-0.040	0.213	0.290	-0.080	0.081	0.162	0.308	-0.074	0.002	0.140	0.333	0.26	0.314	0.338	0.278
3-Hydroxybutanal	-0.029	0.128	0.257	0.024	0.170	0.481	0.165	-0.015	-0.022	-0.067	0.124	0.38	0.233	0.100	0.463
Octanal	-0.257	-0.084	-0.173	-0.191	-0.232	-0.273	-0.265	-0.230	-0.190	-0.100	-0.101	-	-0.413	-0.003	-0.308
Total aldehyde	-0.257	-0.084	-0.173	-0.191	-0.232	-0.273	-0.265	-0.230	-0.190	-0.100	-0.101	-	-0.413	-0.003	-0.308
Benzaldehyde	-0.319	-0.075	0.056	-0.274	-0.125	0.102	0.113	-0.345	-0.186	-0.065	0.049	0.12	0.236	0.086	0.212
Myristicin	0.151	0.193	0.119	0.129	0.202	0.012	0.069	0.144	0.207	0.041	0.212	0.15	0.099	0.182	0.149
Benzeneacetaldehyde	-0.217	0.023	0.041	-0.240	-0.111	-0.049	0.001	-0.244	-0.166	-0.081	0.062	0.15	0.414	-0.040	0.171
Naphthalene	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
1,2-Dihydro-1,1,6-trimethylnaphthalene	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
Benzyl alcohol	-0.337	-0.086	-0.035	-0.290	-0.107	-0.028	0.034	-0.362	-0.188	-0.126	0.058	0.05	0.059	0.100	0.135
Total benzenoid	-0.340	-0.086	-0.028	-0.295	-0.110	-0.020	0.040	-0.365	-0.192	-0.123	0.057	0.06	0.089	0.094	0.147
Ethyl butanoate	0.122	0.275	0.122	0.060	-0.036	-0.214	-0.005	0.173	-0.013	0.165	0.249	-	-0.057	0.242	-0.136
Ethyl 3-methylbutanoate	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
3-Methylbutyl acetate	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
3-Hydroxymandelic acid ethyl ester	-0.373	-0.241	-0.262	-0.503	-0.463	-0.421	-0.464	-0.364	-0.488	0.064	-0.336	-	0.109	-0.390	-0.220
Ethyl hexanoate	0.528	0.471	0.419	0.484	0.434	0.152	0.490	0.418	0.690	0.200	0.450	0.14	0.272	0.406	0.057

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Hexyl acetate	0.400	0.443	0.462	0.447	0.482	0.528	0.421	0.388	0.403	0.055	0.443	0.34	-0.043	0.432	0.319
Ethyl (E)-3-hexenoate	0.228	0.197	0.286	0.138	0.326	0.355	0.298	0.238	0.115	0.101	0.241	0.65	0.661	0.098	0.627
Ethyl lactate	0.012	0.263	0.353	-0.089	0.050	0.193	0.203	-0.015	0.009	0.359	0.327	0.08	0.088	0.399	0.099
Propyl hexanoate	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
Ethyl octanoate	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
Ethyl 3-hydroxybutanoate	-0.046	0.195	0.270	0.007	-0.107	0.325	0.136	-0.055	0.010	0.041	0.222	0.08	0.135	0.223	0.130
Ethyl nonanoate	0.233	0.487	0.489	0.222	0.193	0.248	0.309	0.231	0.210	0.179	0.398	0.21	0.317	0.369	0.181
Isoamyl lactate	0.270	0.510	0.506	0.298	0.313	0.380	0.438	0.262	0.332	0.133	0.558	0.38	0.359	0.552	0.415
Methyl decanoate	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
Ethyl 2-furancarboxylate	0.507	0.669	0.673	0.505	0.551	0.596	0.583	0.520	0.446	0.137	0.646	0.66	0.529	0.538	0.642
Ethyl decanoate	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
Diethyl succinate	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
Ethyl 9-decenoate	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
Ethyl benzeneacetate	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
2-Phenethyl acetate	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
Ethyl dodecanoate	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
Total ester	-0.361	-0.230	-0.249	-0.493	-0.447	-0.405	-0.449	-0.351	-0.479	0.069	-0.323	-	0.133	-0.382	-0.196
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3-Hydroxy-2-butanone	0.428	0.661	0.656	0.373	0.411	0.445	0.468	0.443	0.329	0.269	0.589	0.47	0.422	0.528	0.440
Limonene	0.717	0.885	0.864	0.611	0.648	0.566	0.720	0.727	0.553	0.435	0.831	0.67	0.567	0.759	0.598
Terpinolene	-0.168	0.042	0.144	-0.170	-0.107	0.177	0.060	-0.168	-0.051	0.052	0.061	0.02	0.254	0.064	0.101
Total terpene	0.735	0.893	0.864	0.636	0.659	0.567	0.727	0.745	0.572	0.422	0.840	0.68	0.569	0.764	0.602
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Values in **bold** indicate significant ($p < 0.05$) correlation

Table S7 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated model wines

(continued 3/7)

Variables	Unknown quercetin conjugate	Unknown kaempferol conjugates	Total flavonoids	Caffeine	Ethoxyacetic acid	Acetic acid	Pterin-6-carboxylic acid	Hexanoic acid	Octanoic acid	Nonanoic acid	Total acid	2-Methyl-1-propanol	2-Pentanol	1-Butanol	2-Hexanol
TA	-0.596	-0.641	-0.694	-0.453	-0.409	-0.813	-0.655	-0.343	-0.068	-0.956	-0.282	-0.756	-0.956	-0.956	-0.244
pH	0.558	0.601	0.668	0.451	0.394	0.735	0.650	0.415	0.062	0.901	0.332	0.748	0.901	0.901	0.252
TPC	0.816	0.826	0.942	0.866	-0.274	0.392	0.362	-0.126	-0.016	0.470	-0.108	0.438	0.470	0.470	0.153
TFC	0.763	0.757	0.813	0.745	-0.156	0.374	0.476	-0.089	0.126	0.530	-0.018	0.421	0.530	0.530	0.186
DPPH	0.832	0.920	0.941	0.869	-0.159	0.379	0.247	-0.157	-0.024	0.524	-0.132	0.521	0.524	0.524	0.369
FRAP	0.769	0.903	0.931	0.919	-0.282	0.269	0.193	-0.194	0.024	0.411	-0.140	0.402	0.411	0.411	0.238
ABTS	0.828	0.875	0.930	0.910	-0.246	0.289	0.243	-0.212	-0.051	0.424	-0.186	0.459	0.424	0.424	0.282
L*	-0.176	-0.225	-0.381	-0.055	-0.078	-0.513	-0.343	-0.467	0.193	-0.571	-0.255	-0.410	-0.571	-0.571	0.052
a*	0.034	0.052	0.237	-0.053	-0.055	0.316	0.242	0.493	-0.222	0.354	0.260	0.262	0.354	0.354	-0.136
b*	0.343	0.399	0.538	0.200	0.185	0.666	0.451	0.404	-0.184	0.758	0.214	0.545	0.758	0.758	0.113
Quinic acid	0.014	0.545	0.469	0.431	-0.125	0.199	0.045	0.197	-0.176	0.394	0.062	0.221	0.394	0.394	-0.094
Gallic acid	0.894	0.538	0.743	0.592	-0.158	0.386	0.464	-0.040	0.067	0.396	-0.007	0.477	0.396	0.396	0.274
Chlorogenic acid	0.859	0.902	0.890	0.893	-0.213	0.272	0.281	-0.188	0.101	0.439	-0.102	0.484	0.439	0.439	0.330
Total phenolic acids	0.804	0.851	0.961	0.824	-0.222	0.447	0.417	0.059	-0.039	0.592	0.018	0.558	0.592	0.592	0.198
Gallocatechin	0.775	0.853	0.972	0.813	-0.215	0.480	0.412	-0.011	-0.068	0.604	-0.047	0.507	0.604	0.604	0.165
Epigallocatechin	0.797	0.908	0.925	0.934	-0.273	0.295	0.267	-0.223	0.034	0.428	-0.158	0.398	0.428	0.428	0.212
Procyanidin B1	0.915	0.928	0.965	0.823	-0.090	0.532	0.453	-0.012	0.077	0.661	0.018	0.628	0.661	0.661	0.357
Catechin	0.843	0.892	0.990	0.731	-0.181	0.551	0.395	0.005	-0.055	0.656	-0.029	0.555	0.656	0.656	0.339
Procyanidin B2	0.612	0.816	0.782	0.983	-0.237	0.156	0.332	-0.141	0.114	0.373	-0.062	0.388	0.373	0.373	-0.053
Epicatechin	0.543	0.803	0.837	0.814	-0.236	0.233	0.185	-0.201	-0.184	0.411	-0.234	0.272	0.411	0.411	0.170
Epigallocatechin gallate	0.368	0.711	0.754	0.648	-0.287	0.281	0.228	0.014	-0.130	0.445	-0.055	0.287	0.445	0.445	-0.067
Procyanidin B1 3-O-gallate	0.627	0.746	0.878	0.782	-0.223	0.279	0.273	0.007	-0.202	0.468	-0.091	0.446	0.468	0.468	0.164
Gallocatechin gallate	0.795	0.929	0.909	0.916	-0.252	0.294	0.261	-0.247	0.043	0.443	-0.170	0.409	0.443	0.443	0.256
Quercetin 3-O-di-galactoside/glucoside	0.602	0.640	0.747	0.872	-0.214	0.160	0.291	-0.122	0.016	0.329	-0.089	0.333	0.329	0.329	-0.069
Myricetin 3-O-galactoside 1	0.653	0.420	0.558	0.133	-0.128	0.406	0.011	-0.182	-0.120	0.269	-0.189	0.240	0.269	0.269	0.568
Epicatechin gallate	0.809	0.840	0.921	0.779	-0.064	0.425	0.345	0.064	0.032	0.589	0.055	0.593	0.589	0.589	0.336
Rutin	0.345	0.796	0.751	0.701	-0.187	0.256	0.071	0.046	-0.125	0.474	-0.028	0.282	0.474	0.474	0.106
Quercetin 3-O-galactoside	0.374	0.486	0.574	0.441	-0.197	0.289	0.138	0.187	-0.020	0.422	0.120	0.191	0.422	0.422	0.019

Kaempferol 3-O-rutinoside 2	0.780	0.747	0.850	0.674	-0.022	0.401	0.288	-0.002	-0.053	0.528	-0.029	0.599	0.528	0.528	0.382
Kaempferol 3-O-glucoside	0.208	0.692	0.644	0.610	-0.129	0.220	0.025	0.049	-0.197	0.440	-0.055	0.252	0.440	0.440	0.017
Unknown quercetin conjugate	1	0.779	0.841	0.672	-0.101	0.516	0.475	-0.045	0.177	0.566	0.038	0.655	0.566	0.566	0.377
Unknown kaempferol conjugate	0.779	1	0.918	0.849	-0.177	0.451	0.344	-0.031	0.118	0.618	0.020	0.577	0.618	0.618	0.264
Total flavonoids	0.841	0.918	1	0.814	-0.212	0.493	0.378	-0.050	-0.039	0.616	-0.062	0.528	0.616	0.616	0.290
Caffeine	0.672	0.849	0.814	1	-0.245	0.167	0.290	-0.184	0.092	0.361	-0.103	0.407	0.361	0.361	0.025
Ethoxyacetic acid	-0.101	-0.177	-0.212	-0.245	1	0.489	0.310	0.189	0.115	0.468	0.207	0.352	0.468	0.468	0.199
Acetic acid	0.516	0.451	0.493	0.167	0.489	1	0.556	0.123	0.053	0.910	0.118	0.661	0.910	0.910	0.147
Pterin-6-carboxylic acid	0.475	0.344	0.378	0.290	0.310	0.556	1	0.345	0.233	0.648	0.354	0.535	0.648	0.648	-0.066
Hexanoic acid	-0.045	-0.031	-0.050	-0.184	0.189	0.123	0.345	1	0.439	0.256	0.921	0.408	0.256	0.256	0.161
Octanoic acid	0.177	0.118	-0.039	0.092	0.115	0.053	0.233	0.439	1	0.126	0.754	0.201	0.126	0.126	-0.083
Nonanoic acid	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1	0.245	0.739	1.000	1.000	0.195
Total acid	0.038	0.020	-0.062	-0.103	0.207	0.118	0.354	0.921	0.754	0.245	1	0.387	0.245	0.245	0.084
2-Methyl-1-propanol	0.655	0.577	0.528	0.407	0.352	0.661	0.535	0.408	0.201	0.739	0.387	1	0.739	0.739	0.386
2-Pentanol	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1	1.000	0.195
1-Butanol	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1	0.195
2-Hexanol	0.377	0.264	0.290	0.025	0.199	0.147	-0.066	0.161	-0.083	0.195	0.084	0.386	0.195	0.195	1
3-Methyl-1-butanol	0.531	0.537	0.448	0.327	0.422	0.725	0.625	0.440	0.653	0.833	0.607	0.676	0.833	0.833	0.102
1-(2-Aminoethylamino)-2-propanol	0.274	0.224	0.279	0.041	0.434	0.733	0.380	-0.161	-0.131	0.699	-0.168	0.380	0.699	0.699	-0.143
2-Octen-4-ol	0.729	0.681	0.757	0.871	-0.032	0.288	0.433	-0.060	0.248	0.402	0.059	0.428	0.402	0.402	0.089
3-Methyl-1-pentanol	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
(E)-3-Hexen-1-ol	-0.001	0.106	0.164	-0.041	0.061	0.329	0.268	0.512	-0.014	0.376	0.366	0.159	0.376	0.376	-0.028
1-Hexanol	0.139	0.178	0.161	0.034	0.454	0.296	0.364	0.509	0.008	0.430	0.381	0.503	0.430	0.430	0.562
2-Methyl-4-heptanol	-0.056	-0.041	-0.062	-0.130	0.361	0.024	0.148	0.469	-0.050	0.125	0.327	0.295	0.125	0.125	0.543
2-Ethyl-1-hexanol	0.197	0.386	0.347	0.348	-0.035	0.305	0.085	0.250	0.053	0.319	0.202	0.369	0.319	0.319	0.051
Linalool	0.572	0.623	0.792	0.749	-0.239	0.221	0.176	0.018	-0.134	0.316	-0.053	0.322	0.316	0.316	0.105
1-Octanol	0.564	0.639	0.757	0.787	-0.145	0.293	0.173	-0.127	-0.055	0.358	-0.122	0.288	0.358	0.358	0.057
2,3-Butanediol	0.466	0.554	0.529	0.395	0.477	0.824	0.745	0.322	0.155	0.892	0.307	0.763	0.892	0.892	0.027
Phenylethyl alcohol	-0.105	-0.102	-0.040	-0.314	0.140	0.225	0.238	0.765	-0.010	0.274	0.555	0.311	0.274	0.274	0.031
Total alcohol	0.149	0.129	0.216	-0.097	0.196	0.286	0.263	0.772	-0.044	0.383	0.546	0.504	0.383	0.383	0.431
3-Hydroxybutanal	-0.005	0.268	0.204	-0.044	0.116	0.617	0.303	0.211	-0.085	0.624	0.116	0.282	0.624	0.624	-0.261
Octanal	0.030	-0.159	-0.198	-0.230	0.849	0.524	0.487	0.038	0.191	0.426	0.127	0.353	0.426	0.426	0.044
Total aldehyde	0.030	-0.159	-0.198	-0.230	0.849	0.524	0.487	0.039	0.191	0.426	0.127	0.353	0.426	0.426	0.044
Benzaldehyde	-0.109	-0.101	-0.029	-0.328	0.180	0.279	0.196	0.666	-0.123	0.339	0.434	0.313	0.339	0.339	0.048
Myristicin	0.148	0.168	0.133	0.098	0.480	0.572	0.185	-0.117	0.263	0.526	0.036	0.136	0.526	0.526	-0.232

Benzeneacetaldehyde	-0.096	-0.048	-0.025	-0.247	0.332	0.306	0.173	0.636	-0.085	0.325	0.432	0.304	0.325	0.325	0.335
Naphthalene	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
1,2-Dihydro-1,1,6-trimethyl-naphthalene	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
Benzyl alcohol	-0.135	-0.129	-0.104	-0.337	0.372	0.252	0.189	0.679	-0.121	0.277	0.449	0.388	0.277	0.277	0.312
Total benzenoid	-0.138	-0.129	-0.099	-0.341	0.358	0.256	0.185	0.688	-0.127	0.284	0.453	0.381	0.284	0.284	0.298
Ethyl butanoate	0.320	0.167	0.108	0.101	0.606	0.176	0.307	0.294	0.404	0.320	0.400	0.453	0.320	0.320	0.654
Ethyl 3-methylbutanoate	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
3-Methylbutyl acetate	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
3-Hydroxymandelic acid ethyl ester	-0.221	-0.274	-0.318	-0.449	0.448	0.328	-0.183	0.072	0.003	0.150	0.063	0.030	0.150	0.150	0.087
Ethyl hexanoate	0.470	0.249	0.467	0.501	-0.080	0.248	0.282	-0.115	-0.082	0.192	-0.122	0.186	0.192	0.192	-0.088
Hexyl acetate	0.324	0.458	0.472	0.387	0.178	0.416	0.713	0.258	0.235	0.540	0.289	0.320	0.540	0.540	-0.065
Ethyl (E)-3-hexenoate	-0.069	0.292	0.277	0.157	-0.152	0.050	-0.413	0.055	-0.132	0.192	-0.022	-0.100	0.192	0.192	0.157
Ethyl lactate	0.304	0.157	0.273	-0.112	0.496	0.766	0.391	0.396	0.152	0.639	0.362	0.572	0.639	0.639	0.215
Propyl hexanoate	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
Ethyl octanoate	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
Ethyl 3-hydroxybutanoate	0.184	0.149	0.200	-0.024	0.478	0.585	0.601	0.565	0.372	0.645	0.579	0.509	0.645	0.645	-0.032
Ethyl nonanoate	0.482	0.384	0.437	0.199	0.530	0.846	0.639	0.324	0.156	0.891	0.310	0.735	0.891	0.891	0.246
Isoamyl lactate	0.458	0.450	0.469	0.261	0.543	0.842	0.541	0.262	0.132	0.883	0.254	0.718	0.883	0.883	0.079
Methyl decanoate	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
Ethyl 2-furancarboxylate	0.471	0.682	0.656	0.477	0.365	0.785	0.566	0.274	0.109	0.943	0.248	0.653	0.943	0.943	0.113
Ethyl decanoate	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
Diethyl succinate	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
Ethyl 9-decenoate	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
Ethyl benzeneacetate	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
2-Phenethyl acetate	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
Ethyl dodecanoate	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
Total ester	-0.220	-0.261	-0.304	-0.439	0.441	0.328	-0.195	0.075	-0.001	0.157	0.063	0.028	0.157	0.157	0.094
3-Hydroxy-2-butanone	0.566	0.618	0.616	0.361	0.468	0.910	0.648	0.256	0.126	1.000	0.245	0.739	1.000	1.000	0.195
Limonene	0.765	0.830	0.847	0.632	0.243	0.760	0.511	0.152	0.050	0.901	0.131	0.769	0.901	0.901	0.355
Terpinolene	0.054	-0.026	0.073	-0.214	0.372	0.681	0.393	0.070	-0.097	0.603	0.014	0.213	0.603	0.603	-0.343
Total terpene	0.773	0.843	0.852	0.658	0.232	0.740	0.507	0.148	0.070	0.890	0.137	0.770	0.890	0.890	0.351

Values in **bold** indicate significant ($p < 0.05$) correlation

Table S7 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated model wines

(continued 4/7)

Variables	3-Methyl-1-butanol	1-(2-Aminoethylamino)-2-propanol	2-Octen-4-ol	3-Methyl-1-pentanol	(E)-3-Hexen-1-ol	1-Hexanol	2-Methyl-4-heptanol	2-Ethyl-1-hexanol	Linalool	1-Octanol	2,3-Butanediole	Phenylethyl alcohol	Total alcohol	3-Hydroxybutanol	Octanol
TA	-0.767	-0.589	-0.512	-0.956	-0.414	-0.485	-0.204	-0.388	-0.490	-0.486	-0.875	-0.378	-0.534	-0.515	-0.322
pH	0.723	0.510	0.523	0.901	0.408	0.528	0.272	0.410	0.526	0.506	0.839	0.440	0.608	0.459	0.288
TPC	0.350	0.161	0.882	0.470	0.062	0.085	-0.105	0.368	0.877	0.870	0.428	-0.170	0.084	0.042	-0.232
TFC	0.475	0.246	0.706	0.530	0.033	0.130	-0.082	0.132	0.617	0.668	0.509	-0.145	0.062	0.089	-0.006
DPPH	0.387	0.178	0.784	0.524	0.066	0.100	-0.107	0.308	0.752	0.756	0.426	-0.202	0.091	-0.002	-0.220
FRAP	0.327	0.109	0.827	0.411	-0.006	0.033	-0.153	0.299	0.807	0.803	0.355	-0.294	-0.020	-0.026	-0.312
ABTS	0.295	0.127	0.838	0.424	-0.002	0.052	-0.132	0.276	0.816	0.809	0.364	-0.265	0.023	-0.084	-0.267
L*	-0.328	-0.427	-0.072	-0.571	-0.546	-0.219	-0.100	-0.365	-0.380	-0.261	-0.540	-0.805	-0.742	-0.620	-0.019
a*	0.147	0.286	-0.021	0.354	0.468	0.141	0.081	0.287	0.352	0.213	0.359	0.825	0.720	0.512	-0.094
b*	0.477	0.528	0.220	0.758	0.491	0.397	0.206	0.373	0.467	0.399	0.678	0.655	0.705	0.645	0.116
Quinic acid	0.203	0.189	0.171	0.394	0.265	0.156	0.042	0.416	0.451	0.422	0.407	0.348	0.357	0.563	-0.272
Gallic acid	0.340	0.146	0.771	0.396	0.071	0.086	-0.071	0.156	0.675	0.640	0.331	-0.095	0.146	-0.186	-0.048
Chlorogenic acid	0.391	0.109	0.796	0.439	-0.031	0.059	-0.129	0.212	0.674	0.688	0.370	-0.305	-0.023	-0.102	-0.200
Total phenolic acids	0.430	0.236	0.794	0.592	0.205	0.165	-0.046	0.392	0.869	0.828	0.541	0.091	0.317	0.176	-0.223
Galocatechin	0.423	0.301	0.772	0.604	0.180	0.180	-0.030	0.385	0.844	0.794	0.560	0.032	0.256	0.257	-0.181
Epigallocatechin	0.346	0.094	0.853	0.428	-0.005	0.039	-0.148	0.309	0.793	0.832	0.380	-0.311	-0.040	-0.029	-0.257
Procyanidin B1	0.547	0.258	0.798	0.661	0.159	0.231	0.002	0.343	0.734	0.726	0.573	-0.073	0.213	0.128	-0.084
Catechin	0.470	0.323	0.692	0.656	0.217	0.190	-0.039	0.340	0.749	0.694	0.547	0.041	0.290	0.257	-0.173
Procyanidin B2	0.348	0.071	0.839	0.373	-0.028	0.053	-0.113	0.331	0.727	0.767	0.440	-0.264	-0.080	0.024	-0.191
Epicatechin	0.211	0.174	0.643	0.411	-0.005	0.188	0.033	0.253	0.774	0.789	0.403	-0.169	0.081	0.170	-0.232
Epigallocatechin gallate	0.267	0.315	0.446	0.445	0.156	0.099	-0.068	0.312	0.610	0.504	0.468	0.079	0.162	0.481	-0.273
Procyanidin B1 3-O-gallate	0.244	0.249	0.673	0.468	0.144	0.194	0.028	0.281	0.867	0.761	0.471	0.085	0.308	0.165	-0.265
Galocatechin gallate	0.362	0.119	0.798	0.443	-0.033	0.041	-0.155	0.259	0.712	0.758	0.381	-0.343	-0.074	-0.015	-0.230
Quercetin 3-O-di-galactoside/glucoside	0.260	0.147	0.835	0.329	-0.021	0.103	-0.039	0.247	0.854	0.830	0.386	-0.199	0.002	-0.022	-0.190
Myricetin 3-O-galactoside 1	0.138	0.199	0.252	0.269	-0.043	-0.030	-0.120	0.002	0.322	0.289	0.047	-0.078	0.140	-0.067	-0.100
Epicatechin gallate	0.468	0.286	0.740	0.589	0.124	0.293	0.110	0.281	0.810	0.723	0.552	0.045	0.333	0.124	-0.101
Rutin	0.293	0.129	0.456	0.474	0.346	0.158	0.002	0.450	0.657	0.624	0.438	0.110	0.261	0.381	-0.363
Quercetin 3-O-galactoside	0.311	0.018	0.381	0.422	0.777	-0.040	-0.189	0.447	0.525	0.460	0.315	0.274	0.314	0.233	-0.413

Kaempferol 3-O-rutinoside 2	0.374	0.374	0.646	0.528	-0.083	0.302	0.157	0.168	0.740	0.636	0.492	0.049	0.338	0.100	-0.003
Kaempferol 3-O-glucoside	0.226	0.207	0.331	0.440	0.308	0.168	0.041	0.412	0.579	0.515	0.433	0.178	0.278	0.463	-0.308
Unknown quercetin conjugate	0.531	0.274	0.729	0.566	-0.001	0.139	-0.056	0.197	0.572	0.564	0.466	-0.105	0.149	-0.005	0.030
Unknown kaempferol conjugate	0.537	0.224	0.681	0.618	0.106	0.178	-0.041	0.386	0.623	0.639	0.554	-0.102	0.129	0.268	-0.159
Total flavonoids	0.448	0.279	0.757	0.616	0.164	0.161	-0.062	0.347	0.792	0.757	0.529	-0.040	0.216	0.204	-0.198
Caffeine	0.327	0.041	0.871	0.361	-0.041	0.034	-0.130	0.348	0.749	0.787	0.395	-0.314	-0.097	-0.044	-0.230
Ethoxyacetic acid	0.422	0.434	-0.032	0.468	0.061	0.454	0.361	-0.035	-0.239	-0.145	0.477	0.140	0.196	0.116	0.849
Acetic acid	0.725	0.733	0.288	0.910	0.329	0.296	0.024	0.305	0.221	0.293	0.824	0.225	0.286	0.617	0.524
Pterin-6-carboxylic acid	0.625	0.380	0.433	0.648	0.268	0.364	0.148	0.085	0.176	0.173	0.745	0.238	0.263	0.303	0.487
Hexanoic acid	0.440	-0.161	-0.060	0.256	0.512	0.509	0.469	0.250	0.018	-0.127	0.322	0.765	0.772	0.211	0.038
Octanoic acid	0.653	-0.131	0.248	0.126	-0.014	0.008	-0.050	0.053	-0.134	-0.055	0.155	-0.010	-0.044	-0.085	0.191
Nonanoic acid	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
Total acid	0.607	-0.168	0.059	0.245	0.366	0.381	0.327	0.202	-0.053	-0.122	0.307	0.555	0.546	0.116	0.127
2-Methyl-1-propanol	0.676	0.380	0.428	0.739	0.159	0.503	0.295	0.369	0.322	0.288	0.763	0.311	0.504	0.282	0.353
2-Pentanol	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
1-Butanol	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
2-Hexanol	0.102	-0.143	0.089	0.195	-0.028	0.562	0.543	0.051	0.105	0.057	0.027	0.031	0.431	-0.261	0.044
3-Methyl-1-butanol	1	0.461	0.445	0.833	0.279	0.333	0.067	0.273	0.166	0.243	0.767	0.204	0.268	0.429	0.431
1-(2-Aminoethylamino)-2-propanol	0.461	1	0.045	0.699	-0.047	0.073	-0.150	-0.122	0.052	0.072	0.595	0.069	-0.006	0.639	0.513
2-Octen-4-ol	0.445	0.045	1	0.402	-0.024	0.101	-0.072	0.320	0.787	0.848	0.422	-0.269	-0.017	-0.212	-0.025
3-Methyl-1-pentanol	0.833	0.699	0.402	1	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
(E)-3-Hexen-1-ol	0.279	-0.047	-0.024	0.376	1	0.096	-0.014	0.426	0.121	0.041	0.375	0.606	0.517	0.371	-0.150
1-Hexanol	0.333	0.073	0.101	0.430	0.096	1	0.936	0.160	0.160	0.091	0.428	0.266	0.585	0.204	0.323
2-Methyl-4-heptanol	0.067	-0.150	-0.072	0.125	-0.014	0.936	1	0.096	0.036	-0.056	0.154	0.271	0.566	0.034	0.242
2-Ethyl-1-hexanol	0.273	-0.122	0.320	0.319	0.426	0.160	0.096	1	0.360	0.397	0.409	0.273	0.346	0.237	-0.172
Linalool	0.166	0.052	0.787	0.316	0.121	0.160	0.036	0.360	1	0.936	0.322	0.001	0.253	-0.026	-0.347
1-Octanol	0.243	0.072	0.848	0.358	0.041	0.091	-0.056	0.397	0.936	1	0.343	-0.161	0.081	-0.044	-0.222
2,3-Butanediol	0.767	0.595	0.422	0.892	0.375	0.428	0.154	0.409	0.322	0.343	1	0.306	0.371	0.584	0.507
Phenylethyl alcohol	0.204	0.069	-0.269	0.274	0.606	0.266	0.271	0.273	0.001	-0.161	0.306	1	0.878	0.415	0.035
Total alcohol	0.268	-0.006	-0.017	0.383	0.517	0.585	0.566	0.346	0.253	0.081	0.371	0.878	1	0.259	0.024
3-Hydroxybutanal	0.429	0.639	-0.212	0.624	0.371	0.204	0.034	0.237	-0.026	-0.044	0.584	0.415	0.259	1	0.169
Octanal	0.431	0.513	-0.025	0.426	-0.150	0.323	0.242	-0.172	-0.347	-0.222	0.507	0.035	0.024	0.169	1
Total aldehyde	0.432	0.512	-0.024	0.426	-0.150	0.323	0.242	-0.172	-0.347	-0.222	0.507	0.035	0.024	0.169	1.000
Benzaldehyde	0.190	0.218	-0.309	0.339	0.513	0.306	0.299	0.232	-0.004	-0.147	0.298	0.965	0.853	0.503	0.065
Myristicin	0.548	0.620	0.211	0.526	-0.030	0.061	-0.092	0.012	0.154	0.326	0.442	-0.097	-0.127	0.390	0.460

Benzeneacetaldehyde	0.201	-0.109	-0.188	0.325	0.734	0.629	0.618	0.408	0.014	-0.088	0.334	0.647	0.742	0.322	0.099
Naphthalene	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
1,2-Dihydro-1,1,6-trimethyl-naphthalene	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
Benzyl alcohol	0.144	0.035	-0.289	0.277	0.403	0.696	0.742	0.269	-0.025	-0.158	0.338	0.800	0.871	0.365	0.259
Total benzenoid	0.146	0.040	-0.296	0.284	0.432	0.673	0.716	0.277	-0.022	-0.157	0.335	0.825	0.883	0.380	0.236
Ethyl butanoate	0.470	-0.020	0.287	0.320	-0.032	0.490	0.405	-0.066	-0.069	-0.044	0.261	-0.047	0.227	-0.364	0.456
Ethyl 3-methylbutanoate	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
3-Methylbutyl acetate	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
3-Hydroxymandelic acid ethyl ester	0.116	0.114	-0.351	0.150	0.354	0.016	-0.003	0.010	-0.352	-0.274	0.021	0.132	0.054	0.181	0.282
Ethyl hexanoate	0.101	0.044	0.732	0.192	-0.009	0.071	-0.005	0.264	0.772	0.829	0.214	-0.118	0.050	-0.171	-0.047
Hexyl acetate	0.543	0.379	0.471	0.540	-0.047	0.428	0.234	0.032	0.392	0.395	0.621	0.059	0.166	0.371	0.297
Ethyl (E)-3-hexenoate	0.073	-0.009	-0.020	0.192	0.460	-0.022	-0.104	0.365	0.260	0.238	-0.017	0.102	0.153	0.256	-0.542
Ethyl lactate	0.573	0.526	0.114	0.639	0.270	0.292	0.152	0.236	0.153	0.131	0.660	0.536	0.566	0.425	0.545
Propyl hexanoate	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
Ethyl octanoate	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
Ethyl 3-hydroxybutanoate	0.700	0.520	0.158	0.645	0.340	0.252	0.050	0.121	0.016	-0.064	0.651	0.480	0.411	0.442	0.445
Ethyl nonanoate	0.767	0.564	0.346	0.891	0.376	0.414	0.155	0.345	0.163	0.232	0.797	0.349	0.437	0.458	0.531
Isoamyl lactate	0.747	0.698	0.323	0.883	0.336	0.350	0.131	0.415	0.272	0.289	0.883	0.410	0.453	0.577	0.532
Methyl decanoate	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
Ethyl 2-furancarboxylate	0.781	0.616	0.433	0.943	0.435	0.421	0.108	0.344	0.427	0.444	0.877	0.230	0.340	0.642	0.246
Ethyl decanoate	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
Diethyl succinate	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
Ethyl 9-decenoate	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
Ethyl benzeneacetate	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
2-Phenethyl acetate	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
Ethyl dodecanoate	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
Total ester	0.120	0.113	-0.347	0.157	0.368	0.017	-0.005	0.023	-0.339	-0.262	0.021	0.134	0.060	0.187	0.261
3-Hydroxy-2-butanone	0.833	0.699	0.402	1.000	0.376	0.430	0.125	0.319	0.316	0.358	0.892	0.274	0.383	0.624	0.426
Limonene	0.716	0.513	0.616	0.901	0.321	0.350	0.067	0.387	0.578	0.589	0.781	0.179	0.396	0.386	0.157
Terpinolene	0.406	0.838	-0.161	0.603	0.390	-0.071	-0.263	-0.043	-0.091	-0.121	0.537	0.353	0.120	0.704	0.401
Total terpene	0.718	0.493	0.636	0.890	0.309	0.341	0.059	0.388	0.585	0.600	0.775	0.161	0.381	0.363	0.146

Values in **bold** indicate significant ($p < 0.05$) correlation

Table S7 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated model wines

(continued 5/7)

Variables	Total aldehyde	Benzaldehyde	Myristicin	Benzeneacetaldehyde	Naphthalene	1,2-Dihydro-1,1,6-trimethylnaphthalene	Benzyl alcohol	Total benzeneid	Ethylbutanoate	Ethyl 3-methylbutanoate	3-Methylbutyl acetate	3-Hydroxymandelic acid ethyl ester	Ethyl hexanoate	Hexyl acetate	Ethyl (E)-3-hexenoate
TA	-0.322	-0.430	-0.415	-0.373	-0.956	-0.956	-0.362	-0.370	-0.340	-0.956	-0.956	-0.022	-0.314	-0.563	-0.218
pH	0.289	0.487	0.354	0.399	0.901	0.901	0.424	0.432	0.346	0.901	0.901	-0.020	0.352	0.571	0.199
TPC	-0.232	-0.174	0.124	-0.128	0.470	0.470	-0.221	-0.220	0.023	0.470	0.470	-0.363	0.688	0.461	0.139
TFC	-0.006	-0.133	0.259	-0.162	0.530	0.530	-0.171	-0.174	0.176	0.530	0.530	-0.434	0.464	0.504	0.067
DPPH	-0.220	-0.195	0.119	-0.118	0.524	0.524	-0.238	-0.236	0.248	0.524	0.524	-0.269	0.418	0.317	0.313
FRAP	-0.312	-0.303	0.099	-0.221	0.411	0.411	-0.340	-0.341	0.132	0.411	0.411	-0.371	0.461	0.375	0.300
ABTS	-0.267	-0.265	0.081	-0.195	0.424	0.424	-0.299	-0.299	0.168	0.424	0.424	-0.361	0.528	0.319	0.226
L*	-0.019	-0.853	-0.153	-0.461	-0.571	-0.571	-0.601	-0.628	0.211	-0.571	-0.571	0.009	-0.219	-0.285	-0.244
a*	-0.094	0.862	0.035	0.388	0.354	0.354	0.585	0.613	-0.339	0.354	0.354	-0.089	0.237	0.230	0.168
b*	0.116	0.736	0.284	0.468	0.758	0.758	0.567	0.587	-0.051	0.758	0.758	-0.011	0.306	0.418	0.278
Quinic acid	-0.272	0.391	0.118	0.141	0.394	0.394	0.230	0.246	-0.259	0.394	0.394	-0.224	0.001	0.403	0.527
Gallic acid	-0.048	-0.114	0.061	-0.060	0.396	0.396	-0.144	-0.144	0.217	0.396	0.396	-0.250	0.714	0.255	-0.133
Chlorogenic acid	-0.200	-0.316	0.093	-0.222	0.439	0.439	-0.332	-0.335	0.296	0.439	0.439	-0.356	0.381	0.311	0.193
Total phenolic acids	-0.223	0.100	0.129	0.007	0.592	0.592	-0.020	-0.011	0.053	0.592	0.592	-0.370	0.595	0.473	0.235
Gallocatechin	-0.181	0.051	0.147	0.004	0.604	0.604	-0.032	-0.026	0.006	0.604	0.604	-0.373	0.557	0.517	0.233
Epigallocatechin	-0.257	-0.319	0.151	-0.217	0.428	0.428	-0.337	-0.340	0.122	0.428	0.428	-0.373	0.528	0.400	0.228
Procyanidin B1	-0.084	-0.075	0.193	0.023	0.661	0.661	-0.086	-0.086	0.275	0.661	0.661	-0.241	0.471	0.443	0.197
Catechin	-0.173	0.056	0.119	0.041	0.656	0.656	-0.035	-0.028	0.122	0.656	0.656	-0.262	0.419	0.462	0.286
Procyanidin B2	-0.191	-0.274	0.129	-0.240	0.373	0.373	-0.290	-0.295	0.060	0.373	0.373	-0.503	0.484	0.447	0.138
Epicatechin	-0.232	-0.125	0.202	-0.111	0.411	0.411	-0.107	-0.110	-0.036	0.411	0.411	-0.463	0.434	0.482	0.326
Epigallocatechin gallate	-0.273	0.102	0.012	-0.049	0.445	0.445	-0.028	-0.020	-0.214	0.445	0.445	-0.421	0.152	0.528	0.355
Procyanidin B1 3-O-gallate	-0.265	0.113	0.069	0.001	0.468	0.468	0.034	0.040	-0.005	0.468	0.468	-0.464	0.490	0.421	0.298
Gallocatechin gallate	-0.230	-0.345	0.144	-0.244	0.443	0.443	-0.362	-0.365	0.173	0.443	0.443	-0.364	0.418	0.388	0.238
Quercetin 3-O-digalactoside/glucoside	-0.190	-0.186	0.207	-0.166	0.329	0.329	-0.188	-0.192	-0.013	0.329	0.329	-0.488	0.690	0.403	0.115
Myricetin 3-O-galactoside 1	-0.100	-0.065	0.041	-0.081	0.269	0.269	-0.126	-0.123	0.165	0.269	0.269	0.064	0.200	0.055	0.101
Epicatechin	-0.101	0.049	0.212	0.062	0.589	0.589	0.058	0.057	0.249	0.589	0.589	-0.336	0.450	0.443	0.241

gallate															
Rutin	-0.363	0.123	0.152	0.159	0.474	0.474	0.052	0.064	-0.067	0.474	0.474	-0.220	0.149	0.342	0.651
Quercetin 3-O-galactoside	-0.413	0.236	0.099	0.414	0.422	0.422	0.059	0.089	-0.057	0.422	0.422	0.109	0.272	-0.043	0.661
Kaempferol 3-O-rutinoside 2	-0.003	0.086	0.182	-0.040	0.528	0.528	0.100	0.094	0.242	0.528	0.528	-0.390	0.406	0.432	0.098
Kaempferol 3-O-glucoside	-0.308	0.212	0.149	0.171	0.440	0.440	0.135	0.147	-0.136	0.440	0.440	-0.220	0.057	0.319	0.627
Unknown quercetin conjugate	0.030	-0.109	0.148	-0.096	0.566	0.566	-0.135	-0.138	0.320	0.566	0.566	-0.221	0.470	0.324	-0.069
Unknown kaempferol conjugate	-0.159	-0.101	0.168	-0.048	0.618	0.618	-0.129	-0.129	0.167	0.618	0.618	-0.274	0.249	0.458	0.292
Total flavonoids	-0.198	-0.029	0.133	-0.025	0.616	0.616	-0.104	-0.099	0.108	0.616	0.616	-0.318	0.467	0.472	0.277
Caffeine	-0.230	-0.328	0.098	-0.247	0.361	0.361	-0.337	-0.341	0.101	0.361	0.361	-0.449	0.501	0.387	0.157
Ethoxyacetic acid	0.849	0.180	0.480	0.332	0.468	0.468	0.372	0.358	0.606	0.468	0.468	0.448	-0.080	0.178	-0.152
Acetic acid	0.524	0.279	0.572	0.306	0.910	0.910	0.252	0.256	0.176	0.910	0.910	0.328	0.248	0.416	0.050
Pterin-6-carboxylic acid	0.487	0.196	0.185	0.173	0.648	0.648	0.189	0.185	0.307	0.648	0.648	-0.183	0.282	0.713	-0.413
Hexanoic acid	0.039	0.666	-0.117	0.636	0.256	0.256	0.679	0.688	0.294	0.256	0.256	0.072	-0.115	0.258	0.055
Octanoic acid	0.191	-0.123	0.263	-0.085	0.126	0.126	-0.121	-0.127	0.404	0.126	0.126	0.003	-0.082	0.235	-0.132
Nonanoic acid	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
Total acid	0.127	0.434	0.036	0.432	0.245	0.245	0.449	0.453	0.400	0.245	0.245	0.063	-0.122	0.289	-0.022
2-Methyl-1-propanol	0.353	0.313	0.136	0.304	0.739	0.739	0.388	0.381	0.453	0.739	0.739	0.030	0.186	0.320	-0.100
2-Pentanol	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
1-Butanol	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
2-Hexanol	0.044	0.048	-0.232	0.335	0.195	0.195	0.312	0.298	0.654	0.195	0.195	0.087	-0.088	-0.065	0.157
3-Methyl-1-butanol	0.432	0.190	0.548	0.201	0.833	0.833	0.144	0.146	0.470	0.833	0.833	0.116	0.101	0.543	0.073
1-(2-Aminoethylamino)-2-propanol	0.512	0.218	0.620	-0.109	0.699	0.699	0.035	0.040	-0.020	0.699	0.699	0.114	0.044	0.379	-0.009
2-Octen-4-ol	-0.024	-0.309	0.211	-0.188	0.402	0.402	-0.289	-0.296	0.287	0.402	0.402	-0.351	0.732	0.471	-0.020
3-Methyl-1-pentanol	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
(E)-3-Hexen-1-ol	-0.150	0.513	-0.030	0.734	0.376	0.376	0.403	0.432	-0.032	0.376	0.376	0.354	-0.009	-0.047	0.460
1-Hexanol	0.323	0.306	0.061	0.629	0.430	0.430	0.696	0.673	0.490	0.430	0.430	0.016	0.071	0.428	-0.022
2-Methyl-4-heptanol	0.242	0.299	-0.092	0.618	0.125	0.125	0.742	0.716	0.405	0.125	0.125	-0.003	-0.005	0.234	-0.104
2-Ethyl-1-hexanol	-0.172	0.232	0.012	0.408	0.319	0.319	0.269	0.277	-0.066	0.319	0.319	0.010	0.264	0.032	0.365
Linalool	-0.347	-0.004	0.154	0.014	0.316	0.316	-0.025	-0.022	-0.069	0.316	0.316	-0.352	0.772	0.392	0.260
1-Octanol	-0.222	-0.147	0.326	-0.088	0.358	0.358	-0.158	-0.157	-0.044	0.358	0.358	-0.274	0.829	0.395	0.238
2,3-Butanediol	0.507	0.298	0.442	0.334	0.892	0.892	0.338	0.335	0.261	0.892	0.892	0.021	0.214	0.621	-0.017

Phenylethyl alcohol	0.035	0.965	-0.097	0.647	0.274	0.274	0.800	0.825	-0.047	0.274	0.274	0.132	-0.118	0.059	0.102
Total alcohol	0.024	0.853	-0.127	0.742	0.383	0.383	0.871	0.883	0.227	0.383	0.383	0.054	0.050	0.166	0.153
3-Hydroxybutanal	0.169	0.503	0.390	0.322	0.624	0.624	0.365	0.380	-0.364	0.624	0.624	0.181	-0.171	0.371	0.256
Octanal	1.000	0.065	0.460	0.099	0.426	0.426	0.259	0.236	0.456	0.426	0.426	0.282	-0.047	0.297	-0.542
Total aldehyde	1	0.065	0.460	0.099	0.426	0.426	0.259	0.236	0.457	0.426	0.426	0.282	-0.047	0.297	-0.541
Benzaldehyde	0.065	1	-0.010	0.605	0.339	0.339	0.805	0.831	-0.089	0.339	0.339	0.141	-0.106	0.056	0.153
Myristicin	0.460	-0.010	1	-0.044	0.526	0.526	-0.034	-0.036	0.037	0.526	0.526	0.251	0.287	0.345	0.118
Benzeneacetaldehyde	0.099	0.605	-0.044	1	0.325	0.325	0.832	0.838	0.181	0.325	0.325	0.426	-0.047	-0.053	0.268
Naphthalene	0.426	0.339	0.526	0.325	1	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
1,2-Dihydro-1,1,6-trimethylnaphthalene	0.426	0.339	0.526	0.325	1.000	1	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
Benzyl alcohol	0.259	0.805	-0.034	0.832	0.277	0.277	1	0.999	0.134	0.277	0.277	0.199	-0.080	0.087	-0.005
Total benzenoid	0.236	0.831	-0.036	0.838	0.284	0.284	0.999	1	0.115	0.284	0.284	0.207	-0.083	0.075	0.023
Ethyl butanoate	0.457	-0.089	0.037	0.181	0.320	0.320	0.134	0.115	1	0.320	0.320	0.074	-0.122	0.155	-0.097
Ethyl 3-methylbutanoate	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1	1.000	0.150	0.192	0.540	0.192
3-Methylbutyl acetate	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1	0.150	0.192	0.540	0.192
3-Hydroxymandelic acid ethyl ester	0.282	0.141	0.251	0.426	0.150	0.150	0.199	0.207	0.074	0.150	0.150	1	-0.190	-0.399	0.192
Ethyl hexanoate	-0.047	-0.106	0.287	-0.047	0.192	0.192	-0.080	-0.083	-0.122	0.192	0.192	-0.190	1	0.283	-0.139
Hexyl acetate	0.297	0.056	0.345	-0.053	0.540	0.540	0.087	0.075	0.155	0.540	0.540	-0.399	0.283	1	-0.196
Ethyl (E)-3-hexenoate	-0.541	0.153	0.118	0.268	0.192	0.192	-0.005	0.023	-0.097	0.192	0.192	0.192	-0.139	-0.196	1
Ethyl lactate	0.545	0.527	0.378	0.361	0.639	0.639	0.514	0.515	0.216	0.639	0.639	0.219	0.141	0.352	-0.097
Propyl hexanoate	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
Ethyl octanoate	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
Ethyl 3-hydroxybutanoate	0.445	0.452	0.206	0.233	0.645	0.645	0.288	0.298	0.343	0.645	0.645	0.084	-0.089	0.515	-0.047
Ethyl nonanoate	0.531	0.405	0.341	0.375	0.891	0.891	0.352	0.358	0.395	0.891	0.891	0.198	0.213	0.369	0.029
Isoamyl lactate	0.532	0.459	0.561	0.353	0.883	0.883	0.421	0.425	0.258	0.883	0.883	0.128	0.194	0.396	0.093
Methyl decanoate	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
Ethyl 2-furancarboxylate	0.246	0.277	0.507	0.319	0.943	0.943	0.220	0.228	0.244	0.943	0.943	0.067	0.166	0.584	0.372
Ethyl decanoate	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
Diethyl succinate	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
Ethyl 9-decenoate	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
Ethyl benzenoacetate	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192
2-Phenethyl	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192

acetate																
Ethyl dodecanoate	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192	
Total ester	0.261	0.146	0.254	0.433	0.157	0.157	0.197	0.206	0.074	0.157	0.157	0.999	-0.193	-0.403	0.227	
3-Hydroxy-2-butanone	0.426	0.339	0.526	0.325	1.000	1.000	0.277	0.284	0.320	1.000	1.000	0.150	0.192	0.540	0.192	
Limonene	0.157	0.223	0.362	0.221	0.901	0.901	0.153	0.161	0.354	0.901	0.901	-0.021	0.314	0.445	0.305	
Terpinolene	0.400	0.430	0.507	0.170	0.603	0.603	0.163	0.184	-0.177	0.603	0.603	0.321	-0.035	0.193	0.079	
Total terpene	0.146	0.201	0.358	0.202	0.890	0.890	0.133	0.140	0.365	0.890	0.890	-0.037	0.318	0.442	0.304	

Values in **bold** indicate significant ($p < 0.05$) correlation.

Benzaldehyde	0.527	0.339	0.339	0.452	0.405	0.459	0.339	0.277	0.339	0.339	0.339	0.339	0.339	0.339	0.146
Myristicin	0.378	0.526	0.526	0.206	0.341	0.561	0.526	0.507	0.526	0.526	0.526	0.526	0.526	0.526	0.254
Benzeneacetaldehyde	0.361	0.325	0.325	0.233	0.375	0.353	0.325	0.319	0.325	0.325	0.325	0.325	0.325	0.325	0.433
Naphthalene	0.639	1.000	1.000	0.645	0.891	0.883	1.000	0.943	1.000	1.000	1.000	1.000	1.000	1.000	0.157
1,2-Dihydro-1,1,6-trimethyl-naphthalene	0.639	1.000	1.000	0.645	0.891	0.883	1.000	0.943	1.000	1.000	1.000	1.000	1.000	1.000	0.157
Benzyl alcohol	0.514	0.277	0.277	0.288	0.352	0.421	0.277	0.220	0.277	0.277	0.277	0.277	0.277	0.277	0.197
Total benzenoid	0.515	0.284	0.284	0.298	0.358	0.425	0.284	0.228	0.284	0.284	0.284	0.284	0.284	0.284	0.206
Ethyl butanoate	0.216	0.320	0.320	0.343	0.395	0.258	0.320	0.244	0.320	0.320	0.320	0.320	0.320	0.320	0.074
Ethyl 3-methylbutanoate	0.639	1.000	1.000	0.645	0.891	0.883	1.000	0.943	1.000	1.000	1.000	1.000	1.000	1.000	0.157
3-Methylbutyl acetate	0.639	1.000	1.000	0.645	0.891	0.883	1.000	0.943	1.000	1.000	1.000	1.000	1.000	1.000	0.157
3-Hydroxymandelic acid ethyl ester	0.219	0.150	0.150	0.084	0.198	0.128	0.150	0.067	0.150	0.150	0.150	0.150	0.150	0.150	0.999
Ethyl hexanoate	0.141	0.192	0.192	-0.089	0.213	0.194	0.192	0.166	0.192	0.192	0.192	0.192	0.192	0.192	-
Hexyl acetate	0.352	0.540	0.540	0.515	0.369	0.396	0.540	0.584	0.540	0.540	0.540	0.540	0.540	0.540	0.193
Ethyl (E)-3-hexenoate	-0.097	0.192	0.192	-0.047	0.029	0.093	0.192	0.372	0.192	0.192	0.192	0.192	0.192	0.192	0.403
Ethyl lactate	1	0.639	0.639	0.688	0.701	0.770	0.639	0.497	0.639	0.639	0.639	0.639	0.639	0.639	0.215
Propyl hexanoate	0.639	1	1.000	0.645	0.891	0.883	1.000	0.943	1.000	1.000	1.000	1.000	1.000	1.000	0.157
Ethyl octanoate	0.639	1.000	1	0.645	0.891	0.883	1.000	0.943	1.000	1.000	1.000	1.000	1.000	1.000	0.157
Ethyl 3-hydroxybutanoate	0.688	0.645	0.645	1	0.676	0.642	0.645	0.587	0.645	0.645	0.645	0.645	0.645	0.645	0.082
Ethyl nonanoate	0.701	0.891	0.891	0.676	1	0.835	0.891	0.753	0.891	0.891	0.891	0.891	0.891	0.891	0.199
Isoamyl lactate	0.770	0.883	0.883	0.642	0.835	1	0.883	0.803	0.883	0.883	0.883	0.883	0.883	0.883	0.132
Methyl decanoate	0.639	1.000	1.000	0.645	0.891	0.883	1	0.943	1.000	1.000	1.000	1.000	1.000	1.000	0.157
Ethyl 2-furancarboxylate	0.497	0.943	0.943	0.587	0.753	0.803	0.943	1	0.943	0.943	0.943	0.943	0.943	0.943	0.081
Ethyl decanoate	0.639	1.000	1.000	0.645	0.891	0.883	1.000	0.943	1	1.000	1.000	1.000	1.000	1.000	0.157
Diethyl succinate	0.639	1.000	1.000	0.645	0.891	0.883	1.000	0.943	1.000	1	1.000	1.000	1.000	1.000	0.157
Ethyl 9-decenoate	0.639	1.000	1.000	0.645	0.891	0.883	1.000	0.943	1.000	1.000	1	1.000	1.000	1.000	0.157
Ethyl benzeneacetate	0.639	1.000	1.000	0.645	0.891	0.883	1.000	0.943	1.000	1.000	1.000	1	1.000	1.000	0.157
2-Phenethyl acetate	0.639	1.000	1.000	0.645	0.891	0.883	1.000	0.943	1.000	1.000	1.000	1.000	1	1.000	0.157

Ethyl dodecanoate	0.639	1.000	1.000	0.645	0.891	0.883	1.000	0.943	1.000	1.000	1.000	1.000	1.000	1	0.157
Total ester	0.215	0.157	0.157	0.082	0.199	0.132	0.157	0.081	0.157	0.157	0.157	0.157	0.157	0.157	1
3-Hydroxy-2-butanone	0.639	1.000	1.000	0.645	0.891	0.883	1.000	0.943	1.000	1.000	1.000	1.000	1.000	1.000	0.157
Limonene	0.505	0.901	0.901	0.449	0.761	0.781	0.901	0.885	0.901	0.901	0.901	0.901	0.901	0.901	-0.008
Terpinolene	0.539	0.603	0.603	0.602	0.532	0.637	0.603	0.533	0.603	0.603	0.603	0.603	0.603	0.603	0.320
Total terpene	0.483	0.890	0.890	0.436	0.748	0.771	0.890	0.880	0.890	0.890	0.890	0.890	0.890	0.890	-0.024

Values in **bold** indicate significant ($p < 0.05$) correlation

Table S7 Correlation matrix of quantified phenolic and volatile compounds, physicochemical properties and tested responses of tea-macerated model wines

(continued 7/7)

Variables	3-Hydroxy-2-butanone	Limonene	Terpinolene	Total terpene
TA	-0.956	-0.931	-0.496	-0.922
pH	0.901	0.893	0.422	0.886
TPC	0.470	0.714	-0.042	0.722
TFC	0.530	0.703	0.027	0.716
DPPH	0.524	0.821	-0.084	0.834
FRAP	0.411	0.713	-0.155	0.730
ABTS	0.424	0.740	-0.137	0.756
L*	-0.571	-0.519	-0.565	-0.497
a*	0.354	0.307	0.458	0.286
b*	0.758	0.711	0.552	0.691
Quinic acid	0.394	0.469	0.104	0.468
Gallic acid	0.396	0.594	0.039	0.600
Chlorogenic acid	0.439	0.740	-0.168	0.761
Total phenolic acids	0.592	0.828	0.068	0.834
Gallocatechin	0.604	0.808	0.121	0.811
Epigallocatechin	0.428	0.717	-0.168	0.735
Procyanidin B1	0.661	0.885	0.042	0.893
Catechin	0.656	0.864	0.144	0.864
Procyanidin B2	0.373	0.611	-0.170	0.636
Epicatechin	0.411	0.648	-0.107	0.659
Epigallocatechin gallate	0.445	0.566	0.177	0.567
Procyanidin B1 3-O-gallate	0.468	0.720	0.060	0.727
Gallocatechin gallate	0.443	0.727	-0.168	0.745
Quercetin 3-O-di-galactoside/glucoside	0.329	0.553	-0.051	0.572
Myricetin 3-O-galactoside 1	0.269	0.435	0.052	0.422
Epicatechin gallate	0.589	0.831	0.061	0.840
Rutin	0.474	0.674	0.020	0.681
Quercetin 3-O-galactoside	0.422	0.567	0.254	0.569
Kaempferol 3-O-rutinoside 2	0.528	0.759	0.064	0.764

Kaempferol 3-O-glucoside	0.440	0.598	0.101	0.602
Unknown quercetin conjugate	0.566	0.765	0.054	0.773
Unknown kaempferol conjugate	0.618	0.830	-0.026	0.843
Total flavonoids	0.616	0.847	0.073	0.852
Caffeine	0.361	0.632	-0.214	0.658
Ethoxyacetic acid	0.468	0.243	0.372	0.232
Acetic acid	0.910	0.760	0.681	0.740
Pterin-6-carboxylic acid	0.648	0.511	0.393	0.507
Hexanoic acid	0.256	0.152	0.070	0.148
Octanoic acid	0.126	0.050	-0.097	0.070
Nonanoic acid	1.000	0.901	0.603	0.890
Total acid	0.245	0.131	0.014	0.137
2-Methyl-1-propanol	0.739	0.769	0.213	0.770
2-Pentanol	1.000	0.901	0.603	0.890
1-Butanol	1.000	0.901	0.603	0.890
2-Hexanol	0.195	0.355	-0.343	0.351
3-Methyl-1-butanol	0.833	0.716	0.406	0.718
1-(2-Aminoethylamino)-2-propanol	0.699	0.513	0.838	0.493
2-Octen-4-ol	0.402	0.616	-0.161	0.636
3-Methyl-1-pentanol	1.000	0.901	0.603	0.890
(E)-3-Hexen-1-ol	0.376	0.321	0.390	0.309
1-Hexanol	0.430	0.350	-0.071	0.341
2-Methyl-4-heptanol	0.125	0.067	-0.263	0.059
2-Ethyl-1-hexanol	0.319	0.387	-0.043	0.388
Linalool	0.316	0.578	-0.091	0.585
1-Octanol	0.358	0.589	-0.121	0.600
2,3-Butanediol	0.892	0.781	0.537	0.775
Phenylethyl alcohol	0.274	0.179	0.353	0.161
Total alcohol	0.383	0.396	0.120	0.381
3-Hydroxybutanal	0.624	0.386	0.704	0.363
Octanal	0.426	0.157	0.401	0.146
Total aldehyde	0.426	0.157	0.400	0.146
Benzaldehyde	0.339	0.223	0.430	0.201
Myristicin	0.526	0.362	0.507	0.358
Benzeneacetaldehyde	0.325	0.221	0.170	0.202
Naphthalene	1.000	0.901	0.603	0.890

1,2-Dihydro-1,1,6-trimethyl-naphthalene	1.000	0.901	0.603	0.890
Benzyl alcohol	0.277	0.153	0.163	0.133
Total benzenoid	0.284	0.161	0.184	0.140
Ethyl butanoate	0.320	0.354	-0.177	0.365
Ethyl 3-methylbutanoate	1.000	0.901	0.603	0.890
3-Methylbutyl acetate	1.000	0.901	0.603	0.890
3-Hydroxymandelic acid ethyl ester	0.150	-0.021	0.321	-0.037
Ethyl hexanoate	0.192	0.314	-0.035	0.318
Hexyl acetate	0.540	0.445	0.193	0.442
Ethyl (E)-3-hexenoate	0.192	0.305	0.079	0.304
Ethyl lactate	0.639	0.505	0.539	0.483
Propyl hexanoate	1.000	0.901	0.603	0.890
Ethyl octanoate	1.000	0.901	0.603	0.890
Ethyl 3-hydroxybutanoate	0.645	0.449	0.602	0.436
Ethyl nonanoate	0.891	0.761	0.532	0.748
Isoamyl lactate	0.883	0.781	0.637	0.771
Methyl decanoate	1.000	0.901	0.603	0.890
Ethyl 2-furancarboxylate	0.943	0.885	0.533	0.880
Ethyl decanoate	1.000	0.901	0.603	0.890
Diethyl succinate	1.000	0.901	0.603	0.890
Ethyl 9-decenoate	1.000	0.901	0.603	0.890
Ethyl benzeneacetate	1.000	0.901	0.603	0.890
2-Phenethyl acetate	1.000	0.901	0.603	0.890
Ethyl dodecanoate	1.000	0.901	0.603	0.890
Total ester	0.157	-0.008	0.320	-0.024
3-Hydroxy-2-butanone	1	0.901	0.603	0.890
Limonene	0.901	1	0.340	0.999
Terpinolene	0.603	0.340	1	0.314
Total terpene	0.890	0.999	0.314	1

Values in **bold** indicate significant ($p < 0.05$) correlation

Table S8 PLS-DA results for bottle-aged tea-macerated Chardonnay wines: variables importance in projection (VIP > 1) and standardized coefficient (β)

PLS-DA (Variables)	VIP	β
Total benzenoid	1.435	0.172
Acetic acid	0.568	-0.125
Ethoxyacetic acid	1.918	0.068
Ethyl lactate	1.186	-0.156
Total ester	1.574	0.034
Total aldehyde	0.528	-0.042
Total alcohol	0.480	-0.100
2-Pentanol	1.571	0.134
Ethyl hexanoate	1.429	0.065
Benzaldehyde	1.503	-0.079
2-Methyl-4-heptanol	1.566	0.104
Ethyl benzeneacetate	1.482	-0.083
a*	1.308	-0.084
Ethyl dodecanoate	1.041	0.067

Table S9 PLS-DA results for bottle-aged tea-macerated model wines: variables importance in projection (VIP > 1) and standardized coefficient (β)

PLS-DA (Variables)	VIP	β
Ethyl butanoate	2.298	0.143
Total aldehyde	1.783	0.111
a*	1.613	-0.100
Total ester	1.409	0.088
L*	1.350	0.084
3-Hydroxybutanal	1.225	-0.076
2-Hexanol	1.172	0.073
b*	1.170	-0.073
Ethyl (<i>E</i>)-3-hexenoate	1.143	-0.071
2-Ethyl-1-hexanol	1.044	-0.065
Total phenolic acids	1.043	-0.065

Table S10 Odor active values (OAVs) of volatile compounds in tea-macerated Chardonnay wine at different aging times

Peak No.	Compound Name	Odor descriptor*	Odor threshold (µg/L)	Reference*	C																							
					2GC				4GC				1BC				3BC											
					Aging time (month)																							
1	3	6	9	1	3	6	9	1	3	6	9	1	3	6	9	1	3	6	9	1	3	6	9					
Fruity																												
1	Ethyl butanoate***	Sweet, fruity, juicy	20	(4)	13.804 ± 1.441 ^a	11.816 ± 3.224 ^b	10.766 ± 1.100 ^c	8.410 ± 0.572 ^d	13.419 ± 0.480 ^e	10.182 ± 1.178 ^f	12.325 ± 1.068 ^g	8.345 ± 0.741 ^h	14.995 ± 1.054 ⁱ	13.199 ± 0.841 ^j	12.696 ± 0.837 ^k	8.723 ± 0.754 ^l	14.352 ± 0.585 ^m	11.707 ± 0.412 ⁿ	11.810 ± 1.699 ^o	8.109 ± 0.493 ^p	16.237 ± 0.610 ^q	11.729 ± 2.259 ^r	8.951 ± 0.866 ^s	9.234 ± 1.024 ^t				
2	Ethyl 3-methylbutanoate***	Sweet, estery, fruity	3	(4)	33.138 ± 0.435 ^a	34.038 ± 0.000 ^b	33.041 ± 0.463 ^c	31.593 ± 0.151 ^d	32.833 ± 0.096 ^e	32.508 ± 0.488 ^f	33.291 ± 0.224 ^g	31.429 ± 0.309 ^h	33.123 ± 0.242 ⁱ	32.769 ± 0.180 ^j	33.158 ± 0.180 ^k	31.774 ± 0.288 ^l	33.170 ± 0.314 ^m	32.712 ± 0.170 ⁿ	33.434 ± 0.562 ^o	31.482 ± 0.097 ^p	33.530 ± 0.773 ^q	32.698 ± 4.105 ^r	32.499 ± 1.024 ^s	32.420 ± 0.574 ^t				
4	3-Methylbutyl acetate***	Sweet, banana, fruity	30	(4)	19.782 ± 3.585 ^a	10.203 ± 6.195 ^b	3.621 ± 0.720 ^c	1.847 ± 0.261 ^d	18.738 ± 1.218 ^e	7.078 ± 1.279 ^f	4.971 ± 0.720 ^g	2.100 ± 0.459 ^h	22.675 ± 2.749 ⁱ	11.846 ± 1.281 ^j	11.846 ± 0.551 ^k	3.014 ± 0.604 ^l	20.497 ± 1.860 ^m	9.318 ± 1.028 ⁿ	5.058 ± 1.679 ^o	1.720 ± 0.272 ^p	23.335 ± 1.607 ^q	10.200 ± 4.105 ^r	3.295 ± 1.024 ^s	3.279 ± 0.912 ^t				
11	Ethyl hexanoate***	Sweet, fruity, pineapple	5	(4)	51.223 ± 7.535 ^a	47.904 ± 15.527 ^b	49.494 ± 9.280 ^c	36.460 ± 0.000 ^d	41.251 ± 2.266 ^e	34.399 ± 4.762 ^f	47.997 ± 5.228 ^g	30.129 ± 5.390 ^h	41.597 ± 4.483 ⁱ	41.151 ± 1.281 ^j	43.769 ± 3.868 ^k	27.422 ± 3.200 ^l	47.560 ± 3.879 ^m	42.546 ± 2.294 ⁿ	47.309 ± 8.501 ^o	27.410 ± 3.735 ^p	50.073 ± 2.914 ^q	35.886 ± 8.651 ^r	25.359 ± 1.024 ^s	27.501 ± 4.456 ^t				
14	Hexyl acetate	Fruity, green, apple	63	(7)	0.404 ± 0.069 ^a	0.207 ± 0.113 ^b	0.064 ± 0.013 ^c	0.029 ± 0.004 ^d	0.314 ± 0.022 ^e	0.108 ± 0.017 ^f	0.072 ± 0.003 ^g	0.037 ± 0.008 ^h	0.321 ± 0.037 ⁱ	0.158 ± 0.028 ^j	0.085 ± 0.019 ^k	0.046 ± 0.004 ^l	0.353 ± 0.031 ^m	0.153 ± 0.015 ⁿ	0.076 ± 0.016 ^o	0.027 ± 0.006 ^p	0.318 ± 0.011 ^q	0.052 ± 0.011 ^r	0.058 ± 0.016 ^s	0.048 ± 0.011 ^t				
18	Ethyl (E)-3-hexenoate	Green, fruity, passion fruit	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--				
20	2-Octen-4-ol	Fruity, berry	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--				
22	Ethyl lactate	Sweet, fruity, acidic	14000	(4)	0.010 ± 0.001 ^a	0.015 ± 0.001 ^b	0.016 ± 0.003 ^c	0.011 ± 0.001 ^d	0.010 ± 0.001 ^e	0.014 ± 0.000 ^f	0.014 ± 0.000 ^g	0.012 ± 0.002 ^h	0.011 ± 0.001 ⁱ	0.014 ± 0.000 ^j	0.014 ± 0.000 ^k	0.011 ± 0.000 ^l	0.010 ± 0.001 ^m	0.015 ± 0.001 ⁿ	0.015 ± 0.001 ^o	0.012 ± 0.000 ^p	0.011 ± 0.000 ^q	0.012 ± 0.000 ^r	0.013 ± 0.001 ^s	0.013 ± 0.001 ^t				
23	Propyl hexanoate	Sweet, fruity, juicy	15	(3)	0.428 ± 0.091 ^a	0.553 ± 0.154 ^b	0.545 ± 0.121 ^c	0.199 ± 0.078 ^d	0.315 ± 0.037 ^e	0.401 ± 0.052 ^f	0.534 ± 0.076 ^g	0.122 ± 0.051 ^h	0.279 ± 0.050 ⁱ	0.477 ± 0.068 ^j	0.540 ± 0.081 ^k	0.102 ± 0.031 ^l	0.354 ± 0.047 ^m	0.412 ± 0.035 ⁿ	0.491 ± 0.112 ^o	0.087 ± 0.047 ^p	0.376 ± 0.028 ^q	0.357 ± 0.133 ^r	0.222 ± 0.043 ^s	0.009 ± 0.000 ^t				
31	Benzaldehyde	Strong, sharp, sweet	2000	(4)	0.004 ± 0.000 ^a	0.007 ± 0.001 ^b	0.010 ± 0.002 ^c	0.011 ± 0.001 ^d	0.002 ± 0.002 ^e	0.010 ± 0.001 ^f	0.014 ± 0.001 ^g	0.004 ± 0.004 ^h	0.014 ± 0.002 ⁱ	0.014 ± 0.008 ^j	0.013 ± 0.006 ^k	0.007 ± 0.000 ^l	0.010 ± 0.000 ^m	0.028 ± 0.002 ⁿ	0.047 ± 0.006 ^o	0.009 ± 0.000 ^p	0.013 ± 0.001 ^q	0.020 ± 0.003 ^r	0.022 ± 0.003 ^s	0.022 ± 0.002 ^t				
33	Ethyl 3-hydroxybutanoate	Fruity, green, grape	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--				
37	Isoamyl lactate	Fruity, creamy, nutty	200	(2)	--	0.084 ± 0.000 ^a	<0.001 ^b	0.046 ± 0.009 ^c	0.063 ± 0.003 ^d	0.137 ± 0.009 ^e	0.126 ± 0.013 ^f	0.077 ± 0.006 ^g	0.065 ± 0.015 ^h	0.137 ± 0.023 ⁱ	0.121 ± 0.007 ^j	0.077 ± 0.010 ^k	0.053 ± 0.008 ^l	0.127 ± 0.030 ^m	0.104 ± 0.034 ⁿ	0.081 ± 0.005 ^o	0.066 ± 0.012 ^p	0.117 ± 0.024 ^q	0.123 ± 0.017 ^r	0.115 ± 0.027 ^s				
42	Ethyl decanoate***	Sweet, waxy, fruity	200	(4)	1.603 ± 0.158 ^a	1.576 ± 0.267 ^b	0.916 ± 0.064 ^c	0.676 ± 0.051 ^d	0.883 ± 0.544 ^e	0.569 ± 0.030 ^f	0.458 ± 0.003 ^g	0.454 ± 0.023 ^h	0.503 ± 0.018 ⁱ	0.522 ± 0.006 ^j	0.443 ± 0.000 ^k	0.424 ± 0.000 ^l	0.687 ± 0.005 ^m	0.657 ± 0.010 ⁿ	0.517 ± 0.018 ^o	0.459 ± 0.002 ^p	0.503 ± 0.002 ^q	0.497 ± 0.026 ^r	0.435 ± 0.008 ^s	0.409 ± 0.005 ^t				
43	Diethyl succinate	Fruity, apple, cooked apple	200000	(4)	0.007 ± 0.000 ^a	0.010 ± 0.001 ^b	0.014 ± 0.003 ^c	0.011 ± 0.000 ^d	0.004 ± 0.003 ^e	0.009 ± 0.001 ^f	0.011 ± 0.000 ^g	0.011 ± 0.002 ^h	0.006 ± 0.000 ⁱ	0.008 ± 0.001 ^j	0.011 ± 0.001 ^k	0.010 ± 0.001 ^l	0.007 ± 0.000 ^m	0.009 ± 0.001 ⁿ	0.012 ± 0.001 ^o	0.011 ± 0.001 ^p	0.007 ± 0.000 ^q	0.008 ± 0.000 ^r	0.010 ± 0.001 ^s	0.011 ± 0.001 ^t				
44	Ethyl 9-decenoate***	Fruity, fatty	76.20	(4)	5.149 ± 0.530 ^a	8.398 ± 1.678 ^b	5.642 ± 0.000 ^c	<0.001 ^d	<0.001 ^e	<0.001 ^f	<0.001 ^g	<0.001 ^h	<0.001 ⁱ	<0.001 ^j	<0.001 ^k	<0.001 ^l	<0.001 ^m	0.368 ± 0.000 ⁿ	<0.001 ^o	<0.001 ^p	<0.001 ^q	<0.001 ^r	<0.001 ^s	<0.001 ^t				
Floral																												
35	Linalool***	Citrus, floral, sweet	25	(4)	0.340 ± 0.244 ^a	0.402 ± 0.00 ^b	0.424 ± 0.119 ^c	<0.001 ^d	5.560 ± 3.855 ^e	11.784 ± 0.970 ^f	10.476 ± 0.698 ^g	8.693 ± 1.156 ^h	12.722 ± 1.050 ⁱ	17.137 ± 0.307 ^j	23.430 ± 3.358 ^k	17.971 ± 1.631 ^l	2.785 ± 0.213 ^m	2.975 ± 0.223 ⁿ	2.598 ± 0.169 ^o	1.432 ± 0.068 ^p	6.846 ± 0.459 ^q	6.577 ± 0.068 ^r	7.114 ± 0.406 ^s	5.264 ± 0.441 ^t				
47	Ethyl benzeneacetate	Sweet, floral, honey	650	(4)	0.003 ± 0.000 ^a	0.003 ± 0.000 ^b	0.005 ± 0.001 ^c	0.003 ± 0.000 ^d	0.007 ± 0.000 ^e	0.014 ± 0.001 ^f	0.024 ± 0.001 ^g	0.031 ± 0.004 ^h	0.009 ± 0.001 ⁱ	0.017 ± 0.001 ^j	0.038 ± 0.005 ^k	0.037 ± 0.004 ^l	0.007 ± 0.001 ^m	0.013 ± 0.001 ⁿ	0.023 ± 0.002 ^o	0.020 ± 0.002 ^p	0.015 ± 0.001 ^q	0.020 ± 0.001 ^r	0.038 ± 0.004 ^s	0.046 ± 0.005 ^t				
48	2-Phenethyl acetate	Floral, rose, sweet	650	(4)	0.060 ± 0.000 ^a	0.059 ± 0.000 ^b	0.057 ± 0.000 ^c	0.057 ± 0.000 ^d	0.049 ± 0.017 ^e	0.058 ± 0.000 ^f	0.057 ± 0.000 ^g	0.057 ± 0.000 ^h	0.059 ± 0.000 ⁱ	0.058 ± 0.000 ^j	0.057 ± 0.000 ^k	0.057 ± 0.000 ^l	0.057 ± 0.000 ^m	0.058 ± 0.000 ⁿ	0.058 ± 0.000 ^o	0.057 ± 0.000 ^p	0.059 ± 0.000 ^q	0.058 ± 0.000 ^r	0.058 ± 0.000 ^s	0.057 ± 0.000 ^t				
51	Benzyl alcohol	Floral, rose, phenolic	200000	(4)	0.003 ± 0.000 ^a	<0.001 ^b	0.001 ± 0.000 ^c	0.003 ± 0.001 ^d	0.003 ± 0.000 ^e	0.002 ± 0.000 ^f	0.001 ± 0.000 ^g	0.002 ± 0.000 ^h	0.003 ± 0.000 ⁱ	0.002 ± 0.000 ^j	0.002 ± 0.000 ^k	0.002 ± 0.000 ^l	0.003 ± 0.000 ^m	0.002 ± 0.000 ⁿ	0.002 ± 0.000 ^o	0.002 ± 0.000 ^p	0.004 ± 0.000 ^q	0.002 ± 0.000 ^r	0.002 ± 0.000 ^s	0.002 ± 0.000 ^t				
52	Phenylethyl alcohol***	Sweet, floral, fresh	28922.73	(8)	0.951 ± 0.070 ^a	1.061 ± 0.090 ^b	1.297 ± 0.270 ^c	0.869 ± 0.034 ^d	0.792 ± 0.054 ^e	0.954 ± 0.035 ^f	1.091 ± 0.055 ^g	0.948 ± 0.132 ^h	0.808 ± 0.044 ⁱ	0.960 ± 0.056 ^j	1.141 ± 0.149 ^k	0.852 ± 0.057 ^l	0.856 ± 0.049 ^m	1.070 ± 0.008 ⁿ	1.168 ± 0.075 ^o	0.927 ± 0.080 ^p	0.944 ± 0.041 ^q	1.003 ± 0.044 ^r	1.104 ± 0.119 ^s	0.996 ± 0.092 ^t				
Fermented																												
5	2-Pentanol***	Mild, green, fusel	11	(7)	--	<0.001 ^a	<0.001 ^b	--	64.683 ± 0.000 ^c	<0.001 ^d	<0.001 ^e	<0.001 ^f	115.980 ± 1.884 ^g	<0.001 ^h	<0.001 ⁱ	<0.001 ^j	100.495 ± 14.096 ^k	<0.001 ^l	<0.001 ^m	<0.001 ⁿ	<0.001 ^o	<0.001 ^p	<0.001 ^q	<0.001 ^r				
6	1-Butanol	Fusel, oil, sweet	150000	(4)	0.005 ± 0.001 ^a	<0.001 ^b	0.006 ± 0.001 ^c	0.003 ± 0.001 ^d	0.006 ± 0.001 ^e	<0.001 ^f	0.007 ± 0.001 ^g	0.004 ± 0.001 ^h	0.010 ± 0.001 ⁱ	0.059 ± 0.000 ^j	0.009 ± 0.000 ^k	0.005 ± 0.001 ^l	0.007 ± 0.001 ^m	0.023 ± 0.004 ⁿ	0.007 ± 0.000 ^o	0.004 ± 0.001 ^p	<0.001 ^q	0.009 ± 0.001 ^r	0.005 ± 0.001 ^s	0.005 ± 0.000 ^t				
10	3-Methyl-1-butanol***	Fusel, oily, alcoholic	30000	(4)	2.061 ± 0.145 ^a	2.246 ± 0.129 ^b	2.693 ± 0.516 ^c	2.002 ± 0.000 ^d	1.993 ± 0.105 ^e	2.299 ± 0.073 ^f	2.424 ± 0.126 ^g	2.177 ± 0.264 ^h	2.173 ± 0.121 ⁱ	2.424 ± 0.130 ^j	2.746 ± 0.323 ^k	2.080 ± 0.128 ^l	2.008 ± 0.085 ^m	2.434 ± 0.128 ⁿ	2.618 ± 0.083 ^o	1.994 ± 0.087 ^p	2.326 ± 0.091 ^q	2.277 ± 0.038 ^r	2.423 ± 0.180 ^s	2.319 ± 0.231 ^t				
21	3-Methyl-1-pentanol	Fusel, cognac, wine	1000	(4)	0.020 ± 0.003 ^a	0.025 ± 0.002 ^b	0.035 ± 0.008 ^c	0.016 ± 0.001 ^d	0.020 ± 0.001 ^e	0.024 ± 0.001 ^f	0.031 ± 0.002 ^g	0.018 ± 0.002 ^h	0.019 ± 0.002 ⁱ	0.035 ± 0.004 ^j	0.017 ± 0.001 ^k	0.021 ± 0.001 ^l	0.027 ± 0.001 ^m	0.036 ± 0.003 ⁿ	0.018 ± 0.001 ^o	0.028 ± 0.001 ^p	0.030 ± 0.001 ^q	0.030 ± 0.003 ^r	0.035 ± 0.004 ^s	0.023 ± 0.003 ^t				
39	Methyl decanoate	Oily, wine, fruity	1200	(4)	0.063 ± 0.000 ^a	0.063 ± 0.000 ^b	0.061 ± 0.000 ^c	0.060 ± 0.000 ^d	0.047 ± 0.023 ^e	0.060 ± 0.000 ^f	<0.001 ^g	0.060 ± 0.000 ^h	0.060 ± 0.000 ⁱ	0.060 ± 0.000 ^j	0.060 ± 0.000 ^k	<0.001 ^l	0.061 ± 0.000 ^m	0.061 ± 0.000 ⁿ	0.060 ± 0.000 ^o	0.060 ± 0.000 ^p	0.060 ± 0.000 ^q	0.060 ± 0.000 ^r	0.060 ± 0.000 ^s	0.060 ± 0.000 ^t				
Waxy																												

3	2-Methyl-1-propanol	Ethereal, winey, cortex	65000	(4)	0.003 ± 0.000 ^a	0.003 ± 0.000 ^b	0.003 ± 0.001 ^a	0.002 ± 0.000 ^c	0.002 ± 0.000 ^c	0.003 ± 0.000 ^{ab}	0.003 ± 0.000 ^a	0.003 ± 0.000 ^b	0.003 ± 0.000 ^a	0.003 ± 0.000 ^a	0.003 ± 0.000 ^b	0.003 ± 0.000 ^a	0.003 ± 0.000 ^a	0.003 ± 0.000 ^b	0.003 ± 0.000 ^a	0.003 ± 0.000 ^b	0.003 ± 0.000 ^a	0.003 ± 0.000 ^b	0.003 ± 0.000 ^a		
27	Ethyl octanoate***	Fruity, wine, waxy	2	(4)	982.564 ± 105.266 ^b	1073.47 ± 170.165 ^a	851.083 ± 118.071 ^c	630.095 ± 82.749 ^d	590.123 ± 36.747 ^b	654.168 ± 11.604 ^d	609.223 ± 34.642 ^c	467.220 ± 58.960 ^e	491.568 ± 49.553 ^b	558.17 ± 29.769 ^d	487.806 ± 56.644 ^c	347.335 ± 24.824 ^d	686.660 ± 26.995 ^b	769.825 ± 24.161 ^a	668.132 ± 54.353 ^b	443.175 ± 45.289 ^c	575.955 ± 13.579 ^d	505.53 ± 7 ± 0.000 ^a	369.835 ± 25.619 ^d	324.798 ± 32.603 ^b	
34	Ethyl nonanoate	Waxy, cognac, estery	1300	(4)	0.008 ± 0.000 ^a	0.007 ± 0.001 ^b	<0.001	0.004 ± 0.001 ^c	0.006 ± 0.001 ^a	0.006 ± 0.000 ^a	0.006 ± 0.000 ^a	0.004 ± 0.001 ^b	0.006 ± 0.001 ^a	0.006 ± 0.001 ^a	0.005 ± 0.001 ^b	0.004 ± 0.001 ^a	0.008 ± 0.000 ^a	0.007 ± 0.001 ^b	0.006 ± 0.001 ^a	0.004 ± 0.001 ^d	0.009 ± 0.000 ^a	0.005 ± 0.001 ^b	0.004 ± 0.000 ^a	0.003 ± 0.001 ^d	
36	1-Octanol	Waxy, green, orange	900	(4)	0.014 ± 0.001 ^{bc}	0.015 ± 0.001 ^{bc}	0.020 ± 0.005 ^a	0.013 ± 0.000 ^c	0.011 ± 0.003 ^b	0.013 ± 0.000 ^a	0.014 ± 0.001 ^a	0.010 ± 0.001 ^b	0.008 ± 0.000 ^c	0.011 ± 0.001 ^b	0.013 ± 0.002 ^a	0.008 ± 0.001 ^c	0.012 ± 0.001 ^c	0.014 ± 0.001 ^b	0.017 ± 0.000 ^a	0.012 ± 0.001 ^c	0.011 ± 0.000 ^a	0.013 ± 0.001 ^b	0.014 ± 0.002 ^a	0.010 ± 0.001 ^c	
50	Ethyl dodecanoate	Sweet, waxy, floral	5900	(5)	0.007 ± 0.001 ^a	<0.001	0.004 ± 0.000 ^b	0.004 ± 0.001 ^b	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	<0.001	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	0.004 ± 0.000 ^a	<0.001
54	Nonanoic acid	Waxy, dirty, cheese	3000	(4)	0.722 ± 0.019 ^a	0.542 ± 0.048 ^b	0.507 ± 0.107 ^a	0.301 ± 0.002 ^d	0.262 ± 0.009 ^a	0.254 ± 0.006 ^a	0.278 ± 0.016 ^a	<0.001	0.233 ± 0.003 ^a	0.229 ± 0.012 ^a	0.227 ± 0.014 ^a	<0.001	0.314 ± 0.009 ^a	0.299 ± 0.008 ^b	0.307 ± 0.025 ^a	0.238 ± 0.007 ^c	0.247 ± 0.003 ^b	0.241 ± 0.010 ^b	0.274 ± 0.006 ^a	<0.001	
Citrus																									
8	Limonene***	Citrus, orange, fresh	10	(5)	--	--	--	--	2.048 ± 0.036 ^a	2.100 ± 0.031 ^{ab}	2.150 ± 0.047 ^a	2.150 ± 0.047 ^a	3.138 ± 0.031 ^a	3.135 ± 0.017 ^a	3.154 ± 0.026 ^a	3.154 ± 0.026 ^a	1.773 ± 0.031 ^a	1.786 ± 0.011 ^a	1.751 ± 0.029 ^a	1.785 ± 0.033 ^a	2.059 ± 0.008 ^a	2.073 ± 0.021 ^a	2.038 ± 0.036 ^a	2.038 ± 0.036 ^a	
30	2-Ethyl-1-hexanol	Citrus, fresh, floral	8000	(4)	--	--	--	--	0.001 ± 0.000 ^a	0.001 ± 0.000 ^b	0.002 ± 0.000 ^a	0.002 ± 0.000 ^a	0.001 ± 0.000 ^c	0.001 ± 0.000 ^b	0.002 ± 0.000 ^a	0.001 ± 0.000 ^a	0.001 ± 0.000 ^b	0.001 ± 0.000 ^a	0.001 ± 0.000 ^a	0.001 ± 0.000 ^b	0.001 ± 0.000 ^a	0.001 ± 0.000 ^{ab}	0.001 ± 0.000 ^a	0.001 ± 0.000 ^a	
Green																									
24	(E)-3-Hexen-1-ol***	Green, leafy	400	(4)	<0.001	<0.001	<0.001	<0.001	1.106 ± 0.000	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	0.030 ± 0.000	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	
41	Benzeneacetaldehyde**	Green, sweet, floral	4	(4)	<0.001	<0.001	<0.001	7.081 ± 1.090	8.413 ± 0.609 ^a	<0.001	<0.001	8.908 ± 3.731 ^a	9.310 ± 0.658 ^c	16.267 ± 3.669 ^a	14.864 ± 2.549 ^b	9.369 ± 0.825 ^a	5.072 ± 1.947 ^d	14.944 ± 2.715 ^a	9.983 ± 1.490 ^b	7.902 ± 1.630 ^b	8.088 ± 0.937 ^c	14.721 ± 3.454 ^a	11.045 ± 3.623 ^b	10.232 ± 0.000 ^a	
Herbal																									
15	Terpinolene	Sweet, fresh pine, citrus	200	(4)	--	--	--	--	0.030 ± 0.003 ^a	0.011 ± 0.001 ^c	0.017 ± 0.001 ^b	0.005 ± 0.001 ^d	0.053 ± 0.009 ^a	0.019 ± 0.000 ^c	0.038 ± 0.006 ^b	0.012 ± 0.001 ^d	0.004 ± 0.001 ^a	0.001 ± 0.000 ^a	0.002 ± 0.000 ^b	<0.001	0.010 ± 0.001 ^a	0.003 ± 0.000 ^{bc}	0.003 ± 0.002 ^{bc}	0.001 ± 0.000 ^c	
25	1-Hexanol	Pungen, ethereal, fusel	8000	(4)	0.256 ± 0.146 ^a	0.196 ± 0.012 ^b	0.248 ± 0.047 ^a	0.168 ± 0.010 ^b	0.403 ± 0.023 ^a	0.190 ± 0.004 ^c	0.209 ± 0.010 ^b	0.187 ± 0.021 ^c	0.412 ± 0.023 ^a	0.188 ± 0.014 ^c	0.222 ± 0.030 ^b	0.165 ± 0.010 ^d	0.423 ± 0.026 ^a	0.203 ± 0.009 ^{bc}	0.231 ± 0.008 ^b	0.174 ± 0.006 ^c	0.472 ± 0.017 ^a	0.186 ± 0.004 ^c	0.208 ± 0.017 ^b	0.194 ± 0.023 ^c	
Acidic																									
28	Acetic acid	Pungent, acidic, cheesy	200000	(4)	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	
Aldehydic																									
19	Octanal***	Aldehydic, waxy, citrus	50	(4)	41.641 ± 4.502 ^b	51.365 ± 5.887 ^a	42.490 ± 6.356 ^b	40.117 ± 2.231 ^b	43.980 ± 2.570 ^b	49.580 ± 0.790 ^c	43.140 ± 1.346 ^b	41.450 ± 3.474 ^b	44.400 ± 1.749 ^b	45.647 ± 7.081 ^a	43.595 ± 6.034 ^b	42.406 ± 1.478 ^b	40.779 ± 0.944 ^b	45.716 ± 0.905 ^a	41.138 ± 2.102 ^b	37.159 ± 2.037 ^c	38.062 ± 3.087 ^b	43.399 ± 1.692 ^d	38.651 ± 2.487 ^b	36.929 ± 3.301 ^b	
Balsamic																									
40	Ethyl 2-furancarboxylate	Balsamic, fruity, floral	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	
Buttery																									
16	3-Hydroxy-2-butanone	Sweet, buttery, creamy	150000	(4)	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	
Creamy																									
38	2,3-Butanediol	Fruity, creamy, buttery	120000	(4)	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	
Fatty																									
49	Hexanoic acid***	Sour, fatty, sweaty	3000	(4)	1.456 ± 0.066 ^b	1.329 ± 0.057 ^b	1.826 ± 0.378 ^a	1.121 ± 0.032 ^c	1.213 ± 0.041 ^b	1.230 ± 0.048 ^b	1.500 ± 0.076 ^c	1.228 ± 0.146 ^b	1.212 ± 0.082 ^b	1.205 ± 0.049 ^b	1.596 ± 0.189 ^a	1.163 ± 0.068 ^b	1.366 ± 0.073 ^b	1.372 ± 0.067 ^b	1.681 ± 0.106 ^c	1.250 ± 0.064 ^c	1.480 ± 0.055 ^b	1.240 ± 0.063 ^c	1.593 ± 0.163 ^a	1.423 ± 0.103 ^b	
53	Octanoic acid	Fatty, waxy, rancid	26000	(4)	0.322 ± 0.016 ^a	0.257 ± 0.031 ^c	0.294 ± 0.082 ^b	0.158 ± 0.005 ^d	0.164 ± 0.008 ^c	0.133 ± 0.004 ^b	0.163 ± 0.007 ^c	0.116 ± 0.014 ^c	0.126 ± 0.009 ^a	0.111 ± 0.003 ^{ab}	0.135 ± 0.019 ^a	0.090 ± 0.004 ^b	0.217 ± 0.007 ^c	0.198 ± 0.005 ^b	0.209 ± 0.011 ^a	0.140 ± 0.014 ^c	0.173 ± 0.008 ^c	0.129 ± 0.002 ^c	0.142 ± 0.016 ^b	0.122 ± 0.006 ^c	
Pungent																									
45	Naphthalene	Pungent, dry, tarry	21	(8)	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	
Spicy																									
32	Myristicin	Spicy, warm, balsamic	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	
Winey																									
7	2-Hexanol***	Chemical, winey, fruity	6	(6)	130.603 ± 24.253 ^b	<0.001	165.496 ± 9.811 ^a	47.295 ± 11.319 ^c	169.340 ± 18.097 ^b	<0.001	202.126 ± 27.845 ^a	137.360 ± 57.070 ^c	206.82 ± 13.514 ^b	837.89 ± 0.000 ^a	198.048 ± 39.916 ^b	117.602 ± 42.974 ^c	207.317 ± 34.794 ^b	647.531 ± 175.732 ^a	159.018 ± 30.799 ^c	<0.001	195.939 ± 5.330 ^b	230.12 ± 8 ± 10.898 ^a	160.324 ± 10.443 ^c	<0.001	

Others																							
9	3-Hydroxymandelic acid ethyl ester	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
12	1-(2-Aminoethylamino)-2-propanol	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
13	3-Hydroxybutanal	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
17	Ethoxyacetic acid	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
26	2-Methyl-4-heptanol	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
29	Pterin-6-carboxylic acid	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
46	1,2-dihydro-1,1,6-trimethyl-naphthalene	Licorice	20	(1)	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001

* = Odor descriptors were obtained from TGSC²⁵

** = Reference: (1) = Crandles et al.²⁶, (2) = Jiang et al.²⁷, (3) = Li et al.²⁸, (4) = Liang et al.²⁹, (5) = Pino & Oscar³⁰, (6) = Vázquez-Araújo et al.³¹, (7) = Wang et al.³², (8) = Wang et al.³³.

*** = Odor-active compound (OAV >1 under at least one of the treatments)

Samples of C, 2GC, 4GC etc. refer to Table S2.

-- = Not available.

a-d = Values with different superscripts in the same wine group are significantly different ($p \leq 0.05$)

12	1-(2-Aminoethylamino)- 2-propanol	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
13	3-Hydroxybutanal	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
17	Ethoxyacetic acid	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
26	2-Methyl-4-heptanol	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
29	Pterin-6-carboxylic acid	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--	--
46	1,2-dihydro-1,1,6- trimethyl-naphthalene	Licorice	20	(1)	--	--	--	--	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001	<0.001

* = Odor descriptors were obtained from TGSC²⁵

** = Reference: (1) = Crandles et al.²⁶, (2) = Jiang et al.²⁷, (3) = Li et al.²⁸, (4) = Liang et al.²⁹, (5) = Pino & Oscar³⁰, (6) = Vázquez-Araújo et al.³¹, (7) = Wang et al.³², (8) = Wang et al.³³.

*** = Odor-active compound (OAV >1 under at least one of the treatments)

Samples of M, 2GM, 4GM etc. refer to Table S3.

-- = Not available.

a-d = Values with different superscripts in the same row are significantly different ($p \leq 0.05$)

Table S12 The distribution of 62 consumers by different categories in sensory JAR analysis of the wines

Category		Control Chardonnay	2% (w/v) Green tea Chardonnay	4% (w/v) Green tea Chardonnay	1% (w/v) Black tea Chardonnay	3% (w/v) Black tea Chardonnay
<i>Age</i>	18 to 24	32	32	32	32	32
	25 to 30	26	26	26	26	26
	31 to 40	3	3	3	3	3
	41 to 55	1	1	1	1	1
<i>Gender</i>	Male	25	25	25	25	25
	Female	37	37	37	37	37
<i>Occupation</i>	Student	54	54	54	54	54
	Wine related work	1	1	1	1	1
	Non-wine related work	7	7	7	7	7
<i>Visual</i>						
Colour intensity	Too pale	12	1	0	0	0
	A little pale	24	9	3	0	0
	Just about right	24	38	32	12	3
	A little dark	2	14	26	43	14
	Too dark	0	0	1	7	45
<i>Olfactory (aroma)</i>						
Fruit smell	Too weak	8	8	7	2	9
	Weak	20	24	23	18	18
	Just about right	22	20	24	31	9
	Strong	12	8	7	11	22
	Too strong	0	1	1	0	4
Vegetal smell	Too weak	12	4	2	4	1
	Weak	27	15	8	21	8
	Just about right	19	23	22	26	19
	Strong	4	19	27	9	21
	Too strong	0	1	3	2	13
Floral smell	Too weak	14	10	9	5	8
	Weak	22	24	23	20	19
	Just about right	20	22	22	25	19
	Strong	6	6	8	11	16
	Too strong	0	0	0	1	0

Odour intensity	Too weak	4	4	3	3	2
	Weak	17	13	10	8	6
	Just about right	35	28	33	39	23
	Strong	6	17	15	11	24
	Too strong	0	0	1	1	7
<i>Olfactive (taste/ flavor)</i>						
Fruit flavor	Too weak	5	16	26	3	15
	Weak	18	28	23	16	18
	Just about right	28	16	9	31	13
	Strong	11	2	3	12	15
	Too strong	0	0	1	0	1
Vegetal taste and flavor	Too weak	11	5	5	4	5
	Weak	29	12	16	18	15
	Just about right	14	19	4	31	11
	Strong	6	21	22	8	17
	Too strong	2	5	15	1	14
Floral flavor	Too weak	16	15	17	4	10
	Weak	19	24	27	19	28
	Just about right	24	19	9	31	12
	Strong	3	4	9	8	10
	Too strong	0	0	0	0	2
<i>Mouthfeel/ texture</i>						
Sweetness	Too weak	5	25	29	5	16
	Weak	15	27	24	18	24
	Just about right	24	10	9	28	19
	Strong	18	0	0	11	2
	Too strong	0	0	0	0	1
Bitterness	Too weak	15	2	0	1	2
	Weak	17	5	2	22	8
	Just about right	22	15	9	28	22
	Strong	7	27	18	10	18
	Too strong	1	13	33	1	12
Sourness	Too weak	4	1	4	3	3
	Weak	13	9	4	7	5
	Just about right	29	9	15	32	20

	Strong	14	32	20	16	21
	Too strong	2	11	19	4	13
Astringency	Too weak	2	1	0	0	1
	Weak	14	7	4	11	9
	Just about right	35	21	15	38	19
	Strong	10	24	29	11	26
	Too strong	1	8	14	2	7
Viscosity	Too weak	3	3	4	0	2
	Weak	22	24	18	12	14
	Just about right	32	19	26	38	31
	Strong	5	16	12	11	12
	Too strong	0	0	2	1	3
Length	Too weak	3	3	1	2	2
	Weak	15	15	11	15	9
	Just about right	33	21	19	30	22
	Strong	11	19	19	14	25
	Too strong	0	4	12	1	4
<i>Overall perception</i>						
Overall liking	Dislike very much	0	13	20	3	12
	Dislike moderately	7	10	10	4	11
	Dislike slightly	8	12	19	7	11
	Neither like nor dislike	10	17	6	13	7
	Like slightly	12	7	6	14	7
	Like moderately	14	1	0	7	10
	Like very much	11	2	1	14	4
Purchase intention	Extremely unlikely	1	11	17	3	12
	Unlikely	14	30	32	7	22
	Neutral	23	17	10	24	15
	Likely	21	3	3	22	10
	Extremely likely	3	1	0	6	3

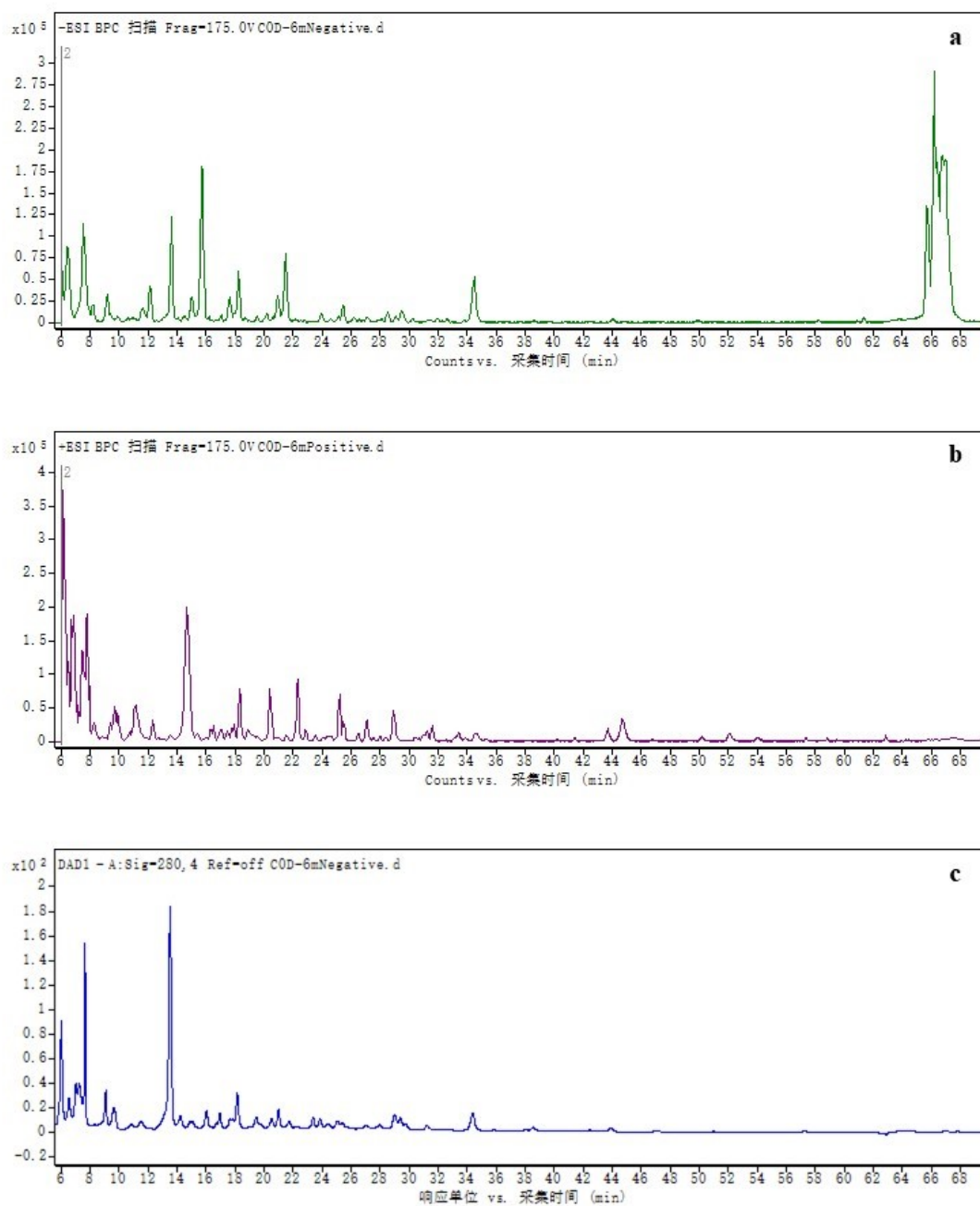


Figure S1 HPLC-DAD-MS chromatograms of phenolic compounds from 6 month-aged control Chardonnay wine detected at negative (a) and positive (b) model and 280 nm (c)

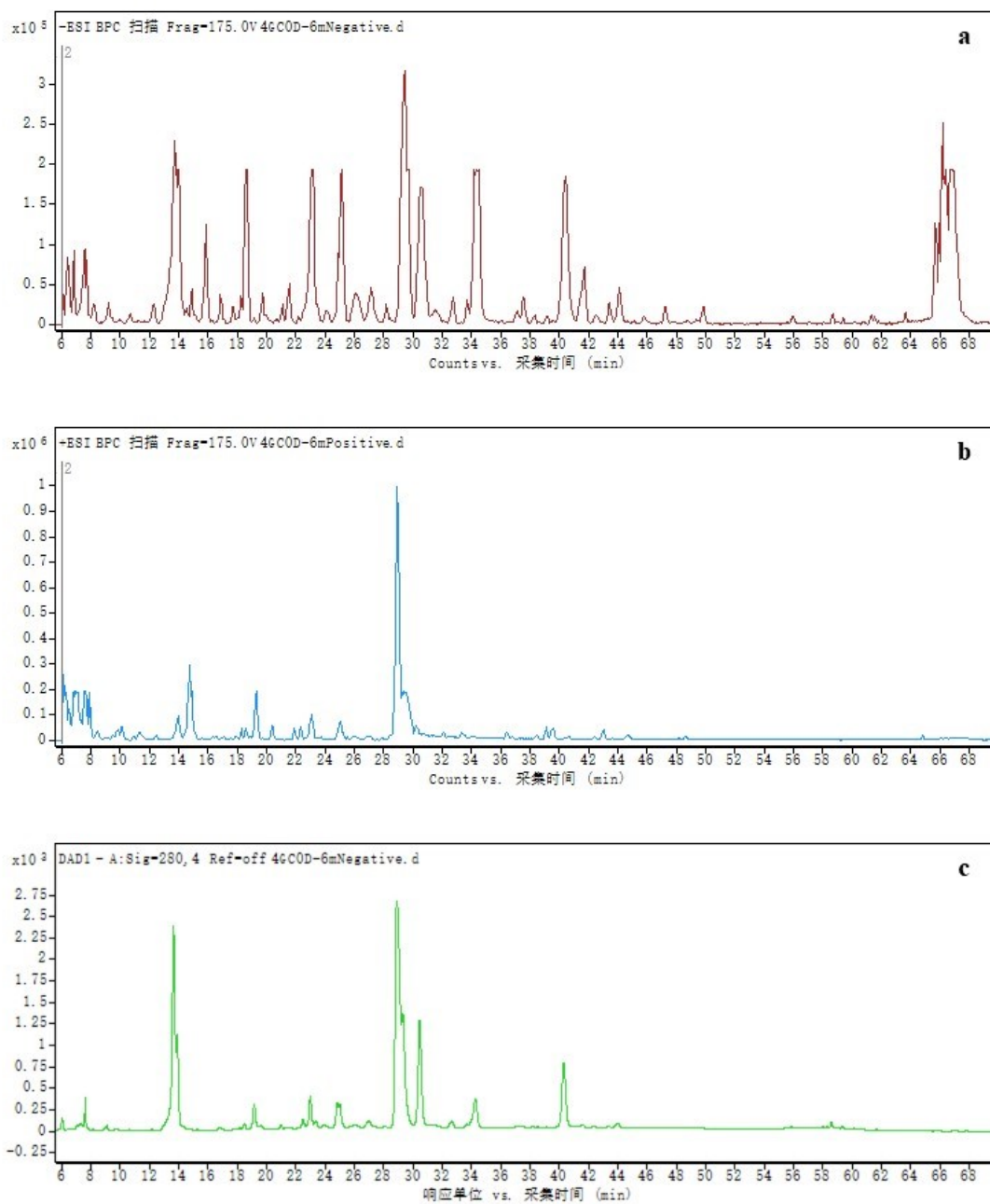


Figure S2 HPLC-DAD-MS chromatograms of phenolic compounds from 6 month-aged green tea-macerated Chardonnay wine detected at negative (a) and positive (b) model and 280 nm (c)

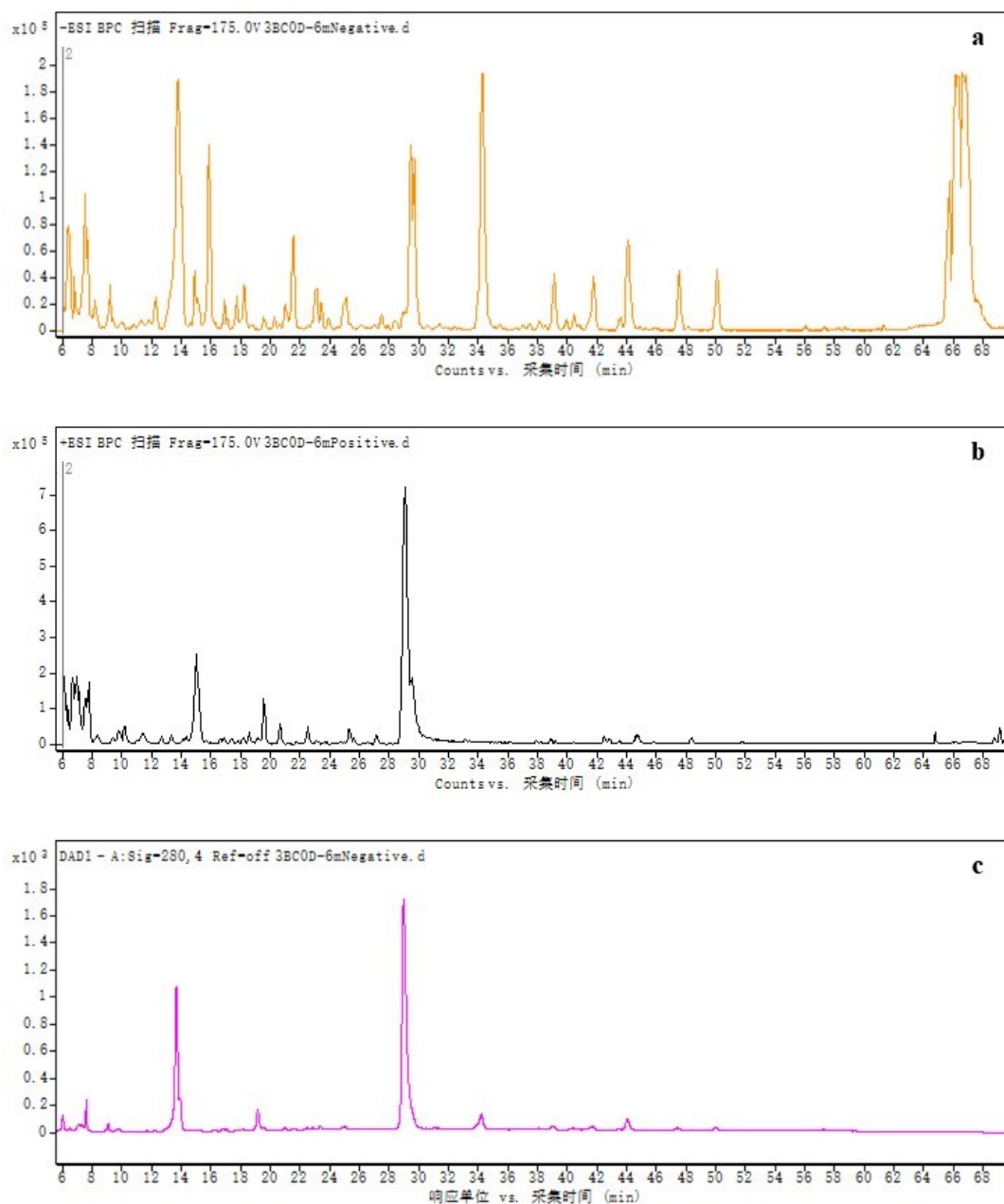
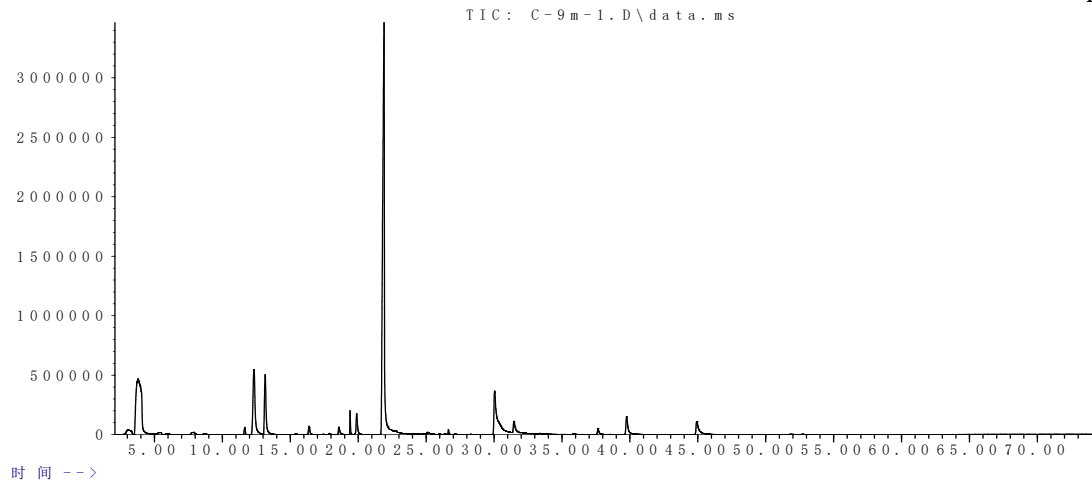


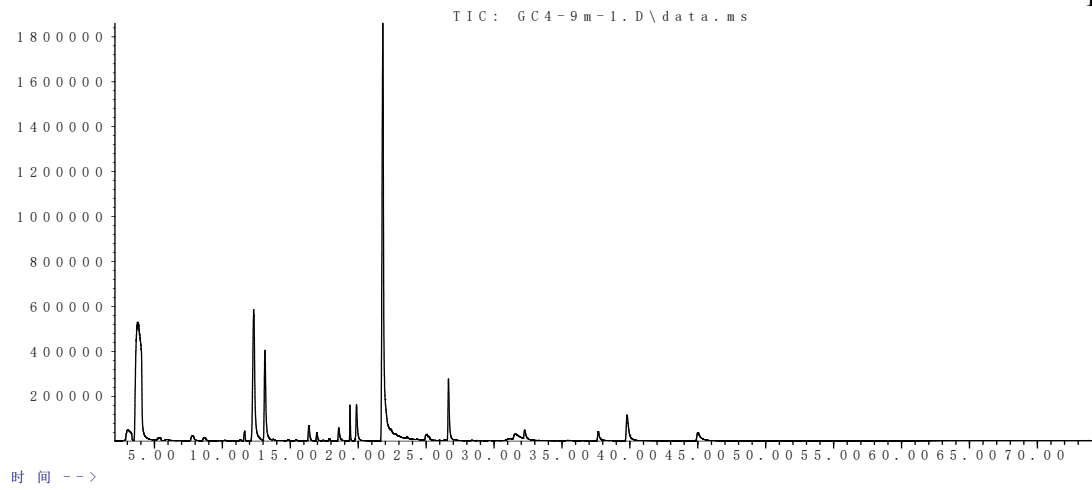
Figure S3 HPLC-DAD-MS chromatograms of phenolic compounds from 6 month-aged black tea-macerated Chardonnay wine detected at negative (a) and positive (b) model and 280 nm (c)

丰度



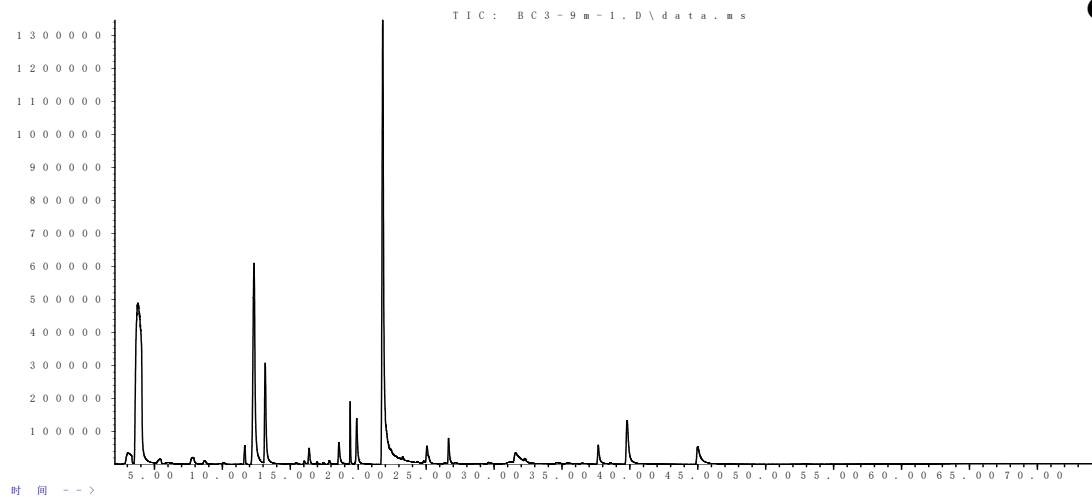
A

丰度



B

丰度



C

Figure S4 Gas chromatography-mass spectrometry (GC-MS) chromatograms showing the volatile compound profile of control Chardonnay wine (A), green tea-macerated

Chardonnay wine (B), and black tea-macerated Chardonnay wine (C) after 9 months of aging

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