

Table S1. Individual phenolic and flavonoid concentration in passion fruit peel powder developed using different carriers (gum arabic, maltodextrin, and waxy starch).

No.	Tentative ID	Retention time	M-H	MS ^E Fragment ions	UV max	Elemental formula
Phenolic acids						
1	Caffeic acid*	1.62	179.0562	266.0874 , 215.0311, 176.0561, 104.0345, 96.9650, 74.0235	229	C ₈ H ₃ O ₅
2	1,3-di-Caffeoylquinic acid	5.29	515.1234	515.1203 , 353.0007, 341.1035, 191.0888, 179.0355, 173.0070, 135.6668, 111.0076, 89.0255	229; 262	C ₂₅ H ₂₃ O ₁₂
3	3,4-di-Caffeoylquinic acid	5.48	515.1295	515.1351 , 473.1112, 353.3334, 341.1131, 335.0388, 191.0004, 179.0988, 173.0093, 135.6668, 111.0077	229; 274	C ₂₅ H ₂₃ O ₁₂
4	3,5-di-Caffeoylquinic acid	5.65	515.1205	515.1243 , 353.0445, 300.0888, 271.8899, 191.0999, 179.0991, 173.0996, 135.0665, 111.0071	229; 274	C ₂₅ H ₂₃ O ₁₂
5	Protocatechuic acid glucoside	13.20	315.1075	243.1686 , 592.1250, 349.0259, 331.1505, 205.0503, 167.0382, 135.0468, 101.0243, 89.0237	229	C ₁₃ H ₁₆ O ₉
6	Galloyl shikimic acid di-HHDP-glucose	17.23	311.1096	311.223 , 179.0888, 149.8980	325	C ₁₄ H ₁₅ O ₉
7	Hydroxy gallic acid	20.77	187.0963	125.0929 , 579.4675, 332.8404, 301.0642, 257.0844, 209.0735, 190.9932, 177.0537, 113.0179, 89.0220, 59.0104	229	C ₇ H ₆ O ₅
8	Valoneic acid	21.87	469.2275	469.4345 , 315.0480, 300.0211, 277.1638, 243.0618, 159.0445, 89.0216	229	-
9	Coumaric acid derivative	26.04	265.1458	265.1475 , 517.3561, 311.1669, 183.0117, 96.9590	229	C ₁₅ H ₂₁ O ₄
10	Vanillic acid glucoside	24.48	329.2294	329.2326 , 253.2153, 187.0959, 171.1025, 152.9954, 125.0953	229	C ₁₈ H ₃₅ O ₅
Flavonoids						
11	(Epi)catechin-(epi)catechin	22.24	575.1392	575.1748 , 285.0396, 301.0350, 257.0447	229	C ₃₀ H ₂₃ O ₁₂
12	Acacetin rhamnoside	8.02	429.0858	151.0260 , 429.0883, 283.0690	229; 271	C ₁₄ H ₂₁ O ₁₅
13	Quercetin*	12.13	301.0920	95.0498 , 139.0387, 152.0120, 301.0875, 161.0395, 71.0147	229; 279	C ₁₃ H ₉ O ₇
14	4'-Methoxyluteolin-8-C-β-D-glucopyranoside	14.66	461.1633	101.0237 , 269.1015, 195.0653, 161.0461, 71.0105	229	C ₁₉ H ₉ O ₁₄
15	Kaempferol 3,7,4'-O-triglucoside	18.22	761.2019	230.0652 , 336.1085, 161.0439, 147.0439, 59.0129	229	C ₃₃ H ₄₀ O ₂₁
16	Luteolin-3-O-(6''-O-malonyl)-glucoside	26.31	533.3900	533.6756 , 327.2171, 287.0389, 179.0996, 161.0904, 111.6158, 101.0230	229	-

Stilbenes						
17	Piceatannol I	19.09	243.0652	243.0662 , 201.0544, 159.0457	229; 282	C ₆ H ₁₅ O ₂₃
18	Piceatannol II	21.98	243.0660	300.0239 , 285.0355, 265.0533, 161.0458, 133.0710, 113.0280	229	C ₅ H ₇ O ₁₁
Unknown compounds						
19	Unknown 1	9.62	219.0495	111.0078 , 173.0123, 67.0175	229; 283	C ₁₅ H ₇ O ₂
20	Unknown 2	10.42	313.0933	101.0222 , 269.0965, 161.0469, 313.0930, 71.0126	229; 262	C ₁₂ H ₁₇ O ₈
21	Unknown 3	12.97	456.1484	119.0369 , 456.1532, 161.0421, 89.0225, 85.0304	229; 281	-
22	Unknown 4	13.35	355.1017	554.2530 , 165.0511, 161.0486, 139.0370, 113.0199	229	C ₂₃ H ₁₅ O ₄
23	Unknown 5	13.97	439.1762	163.0601 , 205.0735, 119.0329, 89.0236, 59.0139	229; 262	C ₁₃ H ₁₁ O ₁₇
24	Unknown 6	14.84	340.1030	161.0436 , 281.1368, 113.0221, 101.0217	229; 261	-
25	Unknown 7	15.22	486.1611	307.1017 , 440.1565, 163.0590, 125.0230, 103.0388, 59.0116	229	-
25	Unknown 8	19.91	455.2090	263.1495 , 161.0415, 101.0242, 89.0229, 59.0122	229	C ₃₁ H ₃ O ₅
27	Unknown 9	22.91	485.1211	485.1206 , 375.0853, 109.0279	229; 326	C ₃₅ H ₁₇ O ₃

*Confirmed using a pure chemical standard;- = unknown or no records in literature sources to our knowledge. Literature sources and *standards were used to corroborate existing observations. MS^E fragments in bold type face refer to the base peak (the highest peak).