

Table S3 Aroma-active compounds in GET characterized by GC-O-MS.

Numbers	Compounds	RT (min) <sup>a</sup>	Aroma descriptions <sup>b</sup>	Aroma intensities <sup>c</sup>
*1	Propanoic acid, 2-methyl-, ethyl ester	8.58	Fruity, banana-like	3.2
*2	2-Methyl-butanal	9.00	Musty, chocolate, nutty	2.5
*3	Pentanal	9.08	Fermented bready	1.7
*4	Butanoic acid, 2-methyl-, methyl ester	9.96	Fruity, apple, green pear	2.0
5	Butanoic acid, ethyl ester	10.87	Green, soapy	3.0
6	Butanoic acid, 2-methyl-, ethyl ester	11.45	Sweat, flower	3.0
*7	Methyl tiglate	17.88	Ethereal	3.2
*8	Hexanoic acid, ethyl ester	19.25	Fruity, sweet	2.2
*9	Acetic acid, hexyl ester	21.11	Fruity, green	3.0
*10	( <i>Z</i> )-3-Hexen-1-ol, acetate	23.22	Green, sweet, fruity	2.0
11	1-Hexanol	24.92	Resin, flower, green	2.2
*12	3-Hexen-1-ol	26.13	Green, leafy, grassy	2.2
*13	Butanoic acid, hexyl ester	27.79	Green, fruity	3.0
14	1-Heptanol	29.20	Fruity, soapy	3.0
15	( <i>Z</i> )-Butanoic acid, 3-hexenyl ester	29.54	Green, apple-like, fruity	2.2
*16	Benzaldehyde	32.33	Bitter, almond-like	2.3
*17	Linalool	33.08	Citrus, floral, sweet	4.0
*18	1-Octanol	34.17	Metallic	2.3
*19	( <i>E, E</i> )-3,5-Octadien-2-one	34.20	Fruity, green grassy	2.0
*20	Benzoic acid, methyl ester	36.51	Fruity	3.0
*21	Safranal	37.36	Woody, spicy, phenolic	2.2
22	( <i>Z</i> )-Hexanoic acid, 3-hexenyl ester	37.53	Green, waxy, winey, grassy	2.3
23	( <i>Z</i> )-Hex-3-enyl ( <i>E</i> )-2-methylbut-2-enoate	38.00	Green	3.0
24	Benzoic acid, ethyl ester	38.15	Green, fruity	2.2
25	Acetic acid, phenylethyl ester	40.48	Sweet, floral, fruity	2.2
26	Acetic acid, 2-phenylethyl ester	40.53	Floral, rose, honey	2.2
27	Alloaromadendrene	41.70	Pungent	2.0
*28	Methyl salicylate	42.33	Mint-like	3.2
29	Benzoic acid, 2-hydroxy-, ethyl ester	43.40	Woody	2.0
30	Geraniol	44.40	Citrus, flower	1.8
31	<i>cis</i> -Calamenene	44.48	Flower	2.7
*32	$\alpha$ -Ionone	44.70	Floral, violet	2.0
*33	Geranylacetone	44.74	Magnolia, green	3.2
*34	Benzyl alcohol	45.31	Fruity, rose-like	3.0
35	Butanoic acid, phenylmethyl ester	45.50	Fruity, sweet	3.0
*36	Phenylethyl alcohol	47.05	Rose, honey	3.2
*37	<i>cis</i> -Jasmone	47.88	Jasmine-like, herbal, floral, woody	4.0
*38	Nerolidol	50.96	Fresh, floral, fruity	2.0
39	2-Phenylethyl tiglate	54.39	Medical	2.8
40	Dihydroactinodiolide	57.60	Musk coumarin	3.0
*41	Indole	59.77	Floral, animal-like	2.0

The symbol of "\*" represented the volatile compounds with OAV greater than 1 in GET;

<sup>a</sup> RT represented the retention time of volatile compounds identified by GC-O-MS;

<sup>b</sup> Aroma descriptions obtained from GC-O-MS;

<sup>c</sup> Aroma intensity perceived by panelists.