

Supplementary Materials: An Early Calcium Loading during Cherry Tree Dormancy Improves Fruit Quality Features at Harvest

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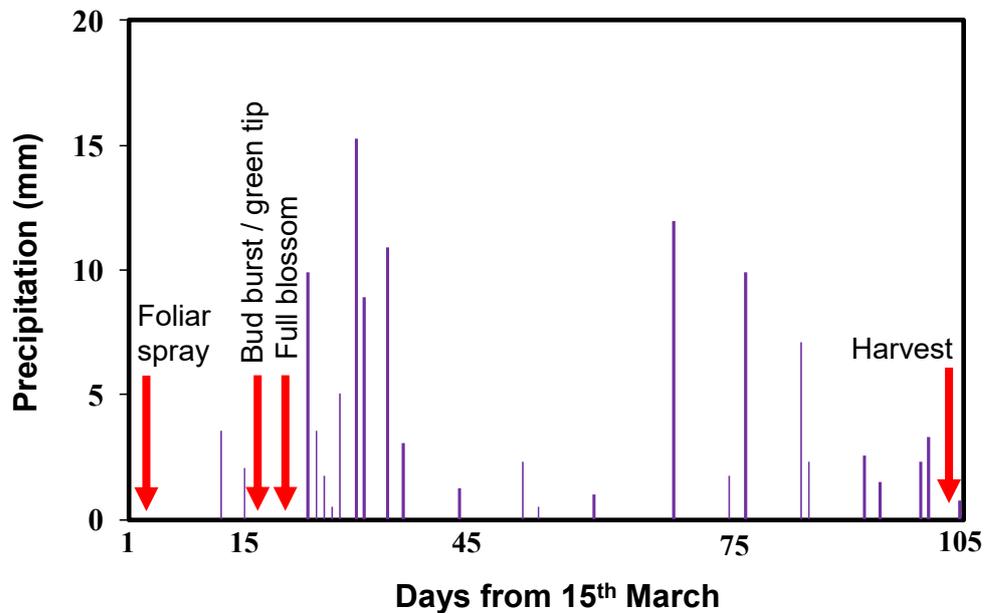


Figure S1. Precipitation graph from 15th March until harvest and timeline of foliar sprays, bud burst, and full blossom.

Table S1. Quantitative results of mineral elements in phloem, flower, and vegetative buds.

Phloem													
	%, DW	P			K			Mg			Na		
		Mean	SD	P-value	Mean	SD	P-value	Mean	SD	P-value	Mean	SD	P-value
	Control	0.33	0.06	NS	0.34	0.03	NS	0.17	0.02	NS	0.03	0.005	NS
Molarity of CaCl ₂ *2H ₂ O	0.25	0.35	0.06		0.4	0.04		0.23	0.02		0.039	0.007	
	0.5	0.31	0.04		0.36	0.03		0.18	0.02		0.03	0.002	
	1	0.34	0.05		0.34	0.04		0.21	0.03		0.034	0.003	
ppm, DW													
	Mean	Zn		P-value	Cu		P-value	Mn		P-value	Fe		P-value
		Mean	SD		Mean	SD		Mean	SD		Mean	SD	
	Control	22.96	3.9	NS	100.12	15.93	NS	99.45	9.95	NS	64.25	12.06	NS
Molarity of CaCl ₂ *2H ₂ O	0.25	28.89	6.55		72.73	20.33		84.94	5.15		64.7	20.46	
	0.5	21.73	5.04		94.97	20.71		86.93	13.35		50.89	8.22	
	1	29.16	7.28		101.56	15.32		87.73	3.42		58.9	16.78	
Flower buds													
	%, DW	P			K			Mg			Na		
		Mean	SD	P-value	Mean	SD	P-value	Mean	SD	P-value	Mean	SD	P-value
	Control	0.81b	0.06	*	1.14	0.1	NS	0.3	0.01	NS	0.027	0.001	NS
Molarity of CaCl ₂ *2H ₂ O	0.25	0.86b	0.05		1.09	0.03		0.29	0.01		0.029	0.002	
	0.5	0.9ab	0.06		1.09	0.06		0.33	0.02		0.033	0.002	
	1	0.99a	0.07		1.19	0.12		0.33	0.02		0.034	0.009	
ppm, DW													
	Mean	Zn		P-value	Cu		P-value	Mn		P-value	Fe		P-value
		Mean	SD		Mean	SD		Mean	SD		Mean	SD	
	Control	20.64	0.67	NS	53.28	15.29	NS	14.4	1.66	NS	44.62	11.31	NS
Molarity of CaCl ₂ *2H ₂ O	0.25	18.78	2.51		32.37	5.92		12.99	0.58		40.62	8.19	
	0.5	19.04	2.48		47.09	18.23		13.1	2.03		37.69	11.45	
	1	20.55	3.38		48.23	3.13		13.59	0.93		37.11	7.3	
Vegetative buds													
	%, DW	P			K			Mg			Na		
		Mean	SD	P-value	Mean	SD	P-value	Mean	SD	P-value	Mean	SD	P-value
	Control	0.84	0.03	NS	0.89	0.06	NS	0.33	0.02	NS	0.048	0.007	NS
Molarity of CaCl ₂ *2H ₂ O	0.25	0.79	0.05		0.88	0.06		0.3	0.01		0.049	0.017	
	0.5	0.8	0.05		0.94	0.08		0.31	0.03		0.055	0.011	
	1	0.83	0.05		0.89	0.12		0.34	0.03		0.037	0.009	
ppm, DW													
	Mean	Zn		P-value	Cu		P-value	Mn		P-value	Fe		P-value
		Mean	SD		Mean	SD		Mean	SD		Mean	SD	
	Control	18.95	4.22	NS	53.79	17.67	NS	15.6	0.6	NS	48.57	5.44	NS
Molarity of CaCl ₂ *2H ₂ O	0.25	22.53	0.13		39.4	8.96		15.96	0.58		45.25	10.62	
	0.5	21.48	1.64		56.42	19.67		16.35	1.55		46.31	2.32	
	1	21.31	2.96		38.75	6.34		14.37	0.08		45.01	9.84	

Different letters indicate significant difference based on Duncan's Multiple Range Test; P ≤ 0.05
NS: No Significance

Table S2. Quantitative results of sweet cherry quality traits.

Ratio flesh fruit ⁻¹ (%)				Penetration force (N)				Polyphenols (µg gallic acid g ⁻¹ FW)				Anthocyanins (µg cyanidin g ⁻¹ FW)				Flavonols (µg rutin g ⁻¹ FW)			
		Mean	SD			Mean	SD			Mean	SD			Mean	SD			Mean	SD
Treatment (T)	Control	94.88	0.4	Treatment (T)	Control	13.78	2.04	Treatment (T)	Control	179.92	34.73	Treatment (T)	Control	26.15	8.44	Treatment (T)	Control	33.68	4.23
	CaCl ₂ * 0.25M	94.86	0.19		CaCl ₂ * 0.25M	13.78	2.48		CaCl ₂ * 0.25M	179.01	26.05		CaCl ₂ * 0.25M	24.79	7.14		CaCl ₂ * 0.25M	34.97	8.99
	0.5M	94.92	0.25		2H ₂ O	13.97	2.21		0.5M	185.84	23.75		0.5M	23.02	4.64		0.5M	37.5	14.33
	1M	94.92	0.26		1M	14.5	1.78		1M	192.69	29.89		1M	27.77	3.93		1M	33.52	9.44
Age of spur (AS) buds				Age of spur (AS) buds				Age of spur (AS) buds				Age of spur (AS) buds				Age of spur (AS) buds			
1st buds		94.91	0.36	1st buds		14.39	2.24	1st buds		185.51	32.69	1st buds		25.3	6.75	1st buds		33.52	10.81
2nd buds		94.89	0.17	2nd buds		13.66	2.01	2nd buds		183.22	24.29	2nd buds		24.01	5.9	2nd buds		36.32	8.51
Ripening (R) Color				Ripening (R) Color				Ripening (R) Color				Ripening (R) Color				Ripening (R) Color			
Red		94.81	0.29	Red		14.43	2.04	Red		176.56	26.32	Red		19.96	2.69	Red		35.63	11.55
Black		94.98	0.25	Black		13.58	2.18	Black		192.17	29.01	Black		29.34	5.31	Black		34.21	7.68
Factors			Factors			Factors			Factors			Factors			Factors				
T	P-value	N	T	P-value	N	T	P-value	N	T	P-value	N	T	P-value	N	T	P-value	N		
T	0.9	12	T	0.205	60	T	0.6	12	T	0.313	12	T	0.767	12	T	0.204	12		
AS	0.805	24	AS	0.003	**	120	AS	0.776	24	AS	0.269	24	AS	0.353	24	AS	0.693	12	
R	0.026	*	24	R	0.008	**	120	R	0.049	*	24	R	<0.001	***	24	R	<0.001	***	12
T*AS	0.28	6	T*AS	0.273	30	T*AS	0.696	6	T*AS	0.352	6	T*AS	0.776	6	T*AS	0.122	6		
T*R	0.151	6	T*R	0.531	30	T*R	0.272	6	T*R	0.079	6	T*R	0.815	6	T*R	0.153	6		
AS*R	0.049	*	12	AS*R	0.881	60	AS*R	0.322	12	AS*R	0.617	12	AS*R	0.266	12	AS*R	0.974	6	
T*AS*R	0.266	3	T*AS*R	0.673	15	T*AS*R	0.134	3	T*AS*R	0.364	3	T*AS*R	0.256	3	T*AS*R	0.164	3		

Different letters indicate significant difference based on Duncan's Multiple Range Test P ≤ 0.05

Hydroxycinnamic acids (µg caffeic acid g ⁻¹ FW)				Skin cracking (% Christensen method)				Skin cracking (% Waterfall method)				Fruit water absorption (% Christensen method)				Fruit water absorption (% Waterfall method)							
		Mean	SD			Mean	SD			Mean	SD			Mean	SD			Mean	SD				
Treatment (T)	Control	41.28	6.21	Treatment (T)	Control	76.94	7.84	Treatment (T)	Control	35	16.78	Treatment (T)	Control	3.17	0.35	Treatment (T)	Control	1.53	0.34				
	CaCl ₂ * 0.25M	40.17	9.24		CaCl ₂ 1M	70.83	12.84		CaCl ₂ 1M	31.67	13.57		CaCl ₂ 1M	3.25	0.27		CaCl ₂ 1M	1.31	0.57				
	0.5M	43.01	9.36		Age of spur (AS) buds				Age of spur (AS) buds				Age of spur (AS) buds				Age of spur (AS) buds						
	1M	39.01	6.6		1st buds		75.28		11.34	1st buds			32.5	15.52	1st buds		3.3	0.48	1st buds		1.43	0.33	
2nd buds		41.29	6.37	2nd buds		72.5	10.67	2nd buds		34.17	15.14	2nd buds		3.22	0.27	2nd buds		1.48	0.47				
Age of spur (AS) buds				Age of spur (AS) buds				Age of spur (AS) buds				Age of spur (AS) buds				Age of spur (AS) buds							
1st buds		40.45	9.25	1st buds		66.39	10.05	1st buds		22.78	7.13	1st buds		3.37	0.43	1st buds		1.18	0.31				
2nd buds		41.29	6.37	2nd buds		81.39	4.81	2nd buds		43.89	13.34	2nd buds		3.16	0.31	2nd buds		1.73	0.28				
Ripening (R) Color				Ripening (R) Color				Ripening (R) Color				Ripening (R) Color				Ripening (R) Color							
Red		43.41	8.15	Red		66.39	10.05	Red		22.78	7.13	Red		3.37	0.43	Red		1.18	0.31				
Black		38.33	6.83	Black		81.39	4.81	Black		43.89	13.34	Black		3.16	0.31	Black		1.73	0.28				
Factors			Factors			Factors			Factors			Factors			Factors								
T	P-value	N	T	P-value	N	T	P-value	N	T	P-value	N	T	P-value	N	T	P-value	N						
T	0.616	12	T	0.068	12	T	0.463	12	T	0.216	12	T	0.204	12	T	0.204	12						
AS	0.703	24	AS	0.387	12	AS	0.712	12	AS	0.563	12	AS	0.693	12	AS	0.693	12						
R	0.027	*	24	R	<0.001	***	12	R	<0.001	***	12	R	<0.001	***	12	R	<0.001	***	12				
T*AS	0.601	6	T*AS	0.174	6	T*AS	0.07	6	T*AS	0.422	6	T*AS	0.122	6	T*AS	0.122	6						
T*R	0.95	6	T*R	0.487	6	T*R	0.667	6	T*R	0.144	6	T*R	0.153	6	T*R	0.153	6						
AS*R	0.158	12	AS*R	0.861	6	AS*R	0.581	6	AS*R	0.059	6	AS*R	0.974	6	AS*R	0.974	6						
T*AS*R	0.096	3	T*AS*R	0.727	3	T*AS*R	0.581	3	T*AS*R	0.354	3	T*AS*R	0.164	3	T*AS*R	0.164	3						

Different letters indicate significant difference based on Duncan's Multiple Range Test P ≤ 0.05

Table S3. Quantitative results of metabolites in sweet cherry skin.

Metabolite reporting checklist			
Level	Aspect	Information	Fill in
general aspect	Type of metabolome analysis	targeted metabolite analysis	FALSE
		non-targeted metabolite class scale profiling	TRUE
		non-targeted metabolite scale profiling	FALSE
	Type of quantification	non-targeted finger printing of mass features	FALSE
		absolute or quantification	Quantification
	Type of reference samples	chemically defined	-
	Type of replication	biologically defined	Frozen grinding skin tissue of <i>Pyrus avium</i> (0.5gr)
		analytical (same analytical sample preparation)	0
	Type of technology	technological (same biological preparation)	0
		biological (same experimental condition)	3
	Sample preparation	full experiment	12
		reference publication	Lipec, J., Schauer, N., Kopka, J., Willmitzer, L., and Fernie, A.R. (2006). Gas chromatography mass spectrometry-based metabolite profiling in plants. <i>Nat. Protoc.</i> 1: 387-396.
metabolite/mass feature	Metabolite	method of sample preparation	Derivatization with methoxyamine hydrochloride and then with N-methyl-N-(trimethylsilyl) tri-fluoroacetamide reagent (MSTFA)
		method of chromatography/separation	Samples (1 µL) was injected and split ratio was 70:1. A TR-5MS capillary column 20m*0.25 mm x 0.25 µm was used. Injector temperature was 220 °C. Carrier gas flow rate was 1ml min ⁻¹ . Temperature program was held at 70 °C for 2 min, then increased to 260 °C (rate 5 °C min ⁻¹), where it remained for 16 min.
	Identification	method of ionization	Ion source 230 °C and the interface 250 °C
		method of detection	Detection type (SQ MS), m/z 50-600 was recorded
		metabolite name	NIST, GOLM, Pubchem
		metabolite sum formula	C ₁₁ H ₁₆ O ₄ , n=number
		metabolite structure and public source of metabolite identifier	NIST, GOLM, Pubchem, KEGG
		identification process	Manually
		by authentic mass isotopomer added to one or all biological sample(s)	TRUE
		by authentic reference compound within a co-processed reference mixture	FALSE
		by authentic reference compound previously mapped to the analytical system	TRUE
		reference library	NIST, GOLM
Quantification	Validity testing	type of mass spectrum	m/z 50-600
		by match of molecular mass (single mass fragment)	-
	by match of fragments	-	
	by match of fragmentation pattern	-	
	by match of mass spectrum to reference library	Yes	
	type of retention index	time elution	
	by match of retention time (index) to reference library	3%	
	type of quantification	Relative abundance	
	Recovery testing (chemical analog)	-	
	Recovery testing (internally added mass isotopomer)	Adonitol	
	Recovery testing (mixture of most divergent samples from the experiment)	-	
	Test for linear range	-	
Limit of quantification (LOQ)	0.001		
Limit of detection (LOD)	0.0005		

<p>Experiment title: Organism/Plant species: <i>Prunus avium L.</i> Organ/tissue: fruit skin Analytical tool: GC-MS</p> <p>Peak no.- number referenced back to the main text Ret. Time- Retention time, in minutes (difference in Ret.Time between ES(+) and ES(-) modes was less than XX minutes) Putative Name- putative identification of the metabolite Mol. Formula- molecular formula of the metabolite or its FA adduct; MSMS fragments- fragments (10 largest) (S)- identification confirmed by a standard compound I, II, III- different isomers Identification level (A; B; C; D)- (A) standard or NMR; (B) MS/MS; (C) MS²; (D) MS only</p>										
Peak no.	Ret. Time	Putative metabolite name	Corresponding metabolite in literature	Metabolite Class	Mol. formula	MS/MS ES(+) fragments (10 largest)	Species detected before	References	Identification level (A-D)	
1	8.185	Oxalic acid (2TMS)	Oxalic acid	Acid	C ₈ H ₁₆ O ₄ Si ₂	147/999, 73/118, 75/67, 52/3, 79/3, 66/2	<i>Solanum lycopersicum L.</i>	nist, goim, pubchem	D	
2	11.39	Silanol, trimethyl-, phosphate (3:1)	Phosphoric acid	Other Compound	C ₉ H ₂₇ O ₄ PSi ₃	299/999, 73/287, 300/222, 301/127, 283/66, 133/62, 225/53, 314/53, 207/52, 147/50	<i>Actinidia chinensis</i>	nist, goim, pubchem	D	
3	11.86	L-Proline, 2tms derivative	Proline	Amino Acid	C ₁₁ H ₂₃ NO ₂ Si ₂	142/999, 73/281, 75/135	<i>Arabidopsis thaliana</i>	nist, goim, pubchem	A	
4	12.15	Butanedioic acid, 2tms derivative	Succinic acid	Acid	C ₁₀ H ₂₀ O ₄ Si ₂	147/999, 73/522, 75/184, 148/108, 149/26, 55/11, 74/7	<i>Arabidopsis thaliana, Crataegus scabrifolia</i>	nist, goim, pubchem	D	
5	15.05	Malic acid, 3tms derivative	Malic acid	Acid	C ₁₃ H ₂₀ O ₅ Si ₃	73/999, 147/785, 233/126, 133/126, 245/126, 148/125, 189/121, 75/113, 74/91, 190/84	<i>Arabidopsis thaliana</i>	nist, goim, pubchem	A	
6	15.5	5-Oxoproline, 2tms derivative	Oxoproline	Amino Acid	C ₁₁ H ₂₃ NO ₂ Si ₂	73/999, 156/231, 75/179, 147/11	<i>Arabidopsis thaliana</i>	nist, goim, pubchem	D	
7	15.67	4-aminobutanoic acid (3TMS)	GABA	Amino Acid	C ₁₃ H ₂₃ O ₂ Si ₃	174/999, 73/959, 147/382, 75/131	<i>Malus domestica</i>	nist, goim, pubchem	D	
8	17.39	D(-)-ribose, tetrakis(trimethylsilyl) ether, trimethylsilyloxime (isomer 2)	Ribose	Sugar	C ₁₈ H ₄₈ NO ₅ Si ₄	73/999, 103/773, 217/388, 147/341, 75/75, 189/62, 160/45, 74/42, 117/42, 133/31	<i>Prunus persica L.</i>	nist, goim, pubchem	D	
9	17.52	D-arabinose, tetrakis(trimethylsilyl) ether, ethyloxime (isomer 2)	Arabinose	Sugar	C ₁₉ H ₄₇ O ₅ NSi ₄	73/999, 103/715, 217/518, 147/362, 160/138, 189/119, 218/90, 117/83, 74/81, 133/75	<i>Actinidia chinensis</i>	nist, goim, pubchem	D	
10	17.61	D-(+)-xylose, tetrakis(trimethylsilyl) ether, methyloxime (syn)	Xylose	Sugar	C ₁₈ H ₄₈ NO ₅ Si ₄	73/999, 204/760, 217/719, 109/542, 147/398, 205/256, 218/82, 129/80, 75/58, 117/43	<i>Prunus avium</i>	nist, goim, pubchem	D	
11	18.22	L(-)-arabitol, 5tms derivative	Arabitol	Alcohol	C ₂₀ H ₃₂ O ₅ Si ₅	73/999, 147/661, 217/631, 103/613, 205/204, 117/64, 129/54, 75/24, 148/5, 189/3	<i>Zizyphus jujube Mill.</i>	nist, goim, pubchem	D	
12	18.32	L-Rhamnose, 4tms derivative	Rhamnose	Sugar	C ₁₈ H ₃₄ O ₅ Si ₄	73/999, 204/987, 147/412, 217/151, 191/75, 75/65, 205/36, 189/1	<i>Arabidopsis thaliana</i>	nist, goim, pubchem	D	
13	20.09	D-(+)-Talofuranose, pentakis(trimethylsilyl) ether (isomer 2)	Talose	Sugar	C ₂₁ H ₃₂ O ₅ Si ₅	73/999, 204/987, 147/412, 217/151, 191/75, 75/65, 205/36, 189/1	<i>Amaranthus tricolor</i>	nist, goim, pubchem	D	
14	20.168	Gulonic acid, -gamma-, lactone, 4tms derivavative	Gulonic acid	Sugar	C ₁₈ H ₃₂ O ₅ Si ₄	73/999, 217/467, 147/427, 204/409, 205/173, 103/150, 129/116, 189/107, 75/96, 117/84	<i>Arabidopsis thaliana</i>	nist, goim, pubchem	D	
14	20.27	Quinic acid (5tms)	Quinic acid	Acid	C ₂₂ H ₃₂ O ₅ Si ₅	73/999, 147/435, 345/359, 255/252, 217/228, 103/214, 191/87, 75/86, 204/58, 205/51	<i>Malus domestica</i>	nist, goim, pubchem	D	
15	20.54	D-Fructose, 1,3,4,5,6-pentakis-O-(trimethylsilyl)-, O-methyloxime	Fructose	Sugar	C ₂₂ H ₃₈ NO ₅ Si ₅	73/999, 103/890, 217/597, 147/384, 307/158, 133/114, 218/110, 277/105, 189/95, 117/89	<i>Prunus avium</i>	nist, goim, pubchem	D	
16	20.88	D-glucose, 2,3,4,5,6-pentakis-o-(trimethylsilyl)-, o-methyloxime, (1z)-	Glucose	Sugar	C ₂₂ H ₃₈ NO ₅ Si ₅	73/999, 147/571, 205/491, 160/461, 319/229, 103/223, 217/219, 117/169, 129/109, 148/90	<i>Arabidopsis thaliana</i>	nist, goim, pubchem	A	
17	21.07	D-galactose, 2,3,4,5,6-pentakis-o-(trimethylsilyl)-, o-methyloxime, (1z)-	Galactose	Sugar	C ₂₂ H ₃₈ NO ₅ Si ₅	73/999, 147/658, 205/535, 160/401, 103/324, 319/270, 217/238, 117/180, 133/119, 129/115	<i>Arabidopsis thaliana</i>	nist, goim, pubchem	D	
18	21.36	D-Sorbitol, 6tms derivative	Sorbitol	Alcohol	C ₂₄ H ₄₂ O ₆ Si ₆	73/999, 147/622, 205/575, 103/422, 217/413, 319/304, 117/180, 129/110, 189/102,	<i>Arabidopsis thaliana</i>	nist, goim, pubchem	A	
20	22.01	D-(+)-Galacturonic acid, 5tms derivative	Galacturonic acid	Acid	C ₂₁ H ₃₀ O ₇ Si ₅	73/999, 147/461, 204/278, 217/275, 205/223, 103/169, 75/82, 117/38, 218/32, 191/29	<i>Arabidopsis thaliana</i>	nist, goim, pubchem	A	
21	23.13	Inositol, (z)-, 6tms derivative	Inositol	Alcohol	C ₂₄ H ₄₆ O ₆ Si ₆	73/999, 217/680, 147/508, 305/482, 191/192, 204/170, 318/133, 75/66, 129/46, 103/38	<i>Arabidopsis thaliana</i>	nist, goim, pubchem	A	
22	29.25	Sucrose, 8tms derivative	Sucrose	Sugar	C ₃₆ H ₆₆ O ₁₁ Si ₈	73/999, 361/977, 217/623, 147/384, 271/313, 169/292, 243/272, 362/252, 129/223,	<i>Actinidia chinensis</i>	nist, goim, pubchem	D	
23	30.43	Maltose, 8tms derivative, isomer 1	Maltose	Sugar	C ₃₆ H ₆₈ O ₁₁ Si ₈	204/999, 73/472, 205/187, 217/174, 147/157, 191/125, 129/104, 206/78, 103/55, 117/46	<i>Malus domestica</i>	nist, goim, pubchem	D	
24	31.66	Melibiose, octakis(trimethylsilyl)-	Melibiose	Sugar	C ₃₆ H ₆₆ O ₁₁ Si ₈	217/999, 73/937, 147/324, 204/292, 361/210, 103/148, 205/121, 129/88, 218/54, 75/22	<i>Solanum lycopersicum L.</i>	nist, goim, pubchem	D	
25	32.182	3- α -mannobiose, octakis(trimethylsilyl) ether (isomer 2)	Mannobiose	Sugar	C ₃₆ H ₆₈ O ₁₁ Si ₈	217/999, 73/937, 147/324, 204/292, 361/210, 103/148, 205/121, 129/88, 218/54, 75/22	<i>Bryonia laciniata</i>	nist, goim, pubchem	D	
26	32.66	D-(+)-Cellobiose, octakis(trimethylsilyl) ether, methyloxime (isomer 1)	Cellobiose	Sugar	C ₃₇ H ₆₈ O ₁₁ NSi ₈	204/999, 73/785, 217/380, 361/300, 147/282, 205/208, 160/173, 103/130, 129/117, 75/57	<i>Olea europaea</i>	nist, goim, pubchem	D	
28	33.97	D-lactose, octakis(trimethylsilyl) ether, methyloxime (isomer 2)	Lactose	Sugar	C ₃₇ H ₆₈ O ₁₁ NSi ₈	204/999, 73/667, 217/394, 205/371, 147/356, 361/228, 103/204, 129/112, 206/92, 319/89	<i>Actinidia chinensis</i>	nist, goim, pubchem	D	
29	34.43	D-Lactitol, nonakis(trimethylsilyl) ether	Lactitol	Alcohol	C ₃₉ H ₈₀ O ₁₁ Si ₉	204/999, 73/760, 217/454, 205/396, 147/395, 361/317, 103/223, 129/152, 319/135,	<i>Malus domestica</i>	nist, goim, pubchem	D	
30	40.24	3-o-coumaroul-D-Quinic acid	3-o-coumaroul-D-Quinic acid	Other Compound	C ₃₁ H ₅₈ O ₈ Si ₅	73/999, 219/836, 345/514, 255/206, 147/103, 75/4, 346/2	<i>Arabidopsis thaliana</i>	nist, goim, pubchem	D	

Full ripe sweet cherries

	1st-year-old spur					2nd-year-old spur						
	Control		1M Ca		t-test	Control		1M Ca		t-test		
	Mean	SD	Mean	SD		Mean	SD	Mean	SD			
Oxalic acid	0.009	0.012	0.000	0.000	0.265		0.004	0.003	0.007	0.009	0.678	
Phosphoric acid	0.895	0.082	1.027	0.119	0.190		1.003	0.083	1.144	0.020	0.045	*
Proline	0.013	0.011	0.016	0.007	0.729		0.012	0.005	0.009	0.008	0.527	
Succinic acid	0.094	0.037	0.116	0.019	0.420		0.120	0.014	0.122	0.014	0.836	
Malic acid	6.878	0.530	8.261	0.277	0.016	*	7.754	0.244	7.775	0.229	0.917	
Oxoproline	0.003	0.000	0.001	0.000	0.001	**	0.004	0.001	0.004	0.003	0.997	
GABA	0.009	0.006	0.008	0.003	0.697		0.010	0.005	0.008	0.004	0.562	
Ribose	0.193	0.008	0.141	0.010	0.002	**	0.168	0.005	0.152	0.008	0.040	*
Arabinose	1.256	0.034	0.999	0.061	0.003	**	1.147	0.066	1.023	0.054	0.067	
Xylose	0.014	0.002	0.009	0.001	0.010	*	0.013	0.002	0.011	0.002	0.226	
Arabitol	0.062	0.009	0.071	0.003	0.148		0.075	0.011	0.069	0.002	0.411	
Rhamnose	0.004	0.002	0.036	0.020	0.042	*	0.021	0.007	0.024	0.009	0.656	
Talose	3.481	2.291	2.639	0.658	0.574		2.287	0.480	2.487	0.441	0.624	
Gulonic acid	0.304	0.526	0.425	0.424	0.771		0.244	0.221	0.000	0.000	0.129	
Quinic acid	0.322	0.053	0.693	0.101	0.005	**	0.494	0.099	0.506	0.099	0.892	
Fructose	132.644	3.680	141.678	7.517	0.135		128.288	1.784	124.462	5.111	0.288	
Glucose	162.730	6.673	176.427	18.584	0.296		159.489	11.711	151.230	3.236	0.304	
Galactose	22.630	1.676	16.634	2.413	0.024	*	17.483	1.233	16.632	0.823	0.376	
Sorbitol	83.091	4.620	84.469	11.091	0.852		74.686	5.169	70.020	2.240	0.225	
Galacturonic acid	0.356	0.109	0.354	0.108	0.977		0.424	0.093	0.415	0.021	0.881	
Inositol	0.130	0.014	0.111	0.026	0.319		0.077	0.066	0.060	0.103	0.822	
Sucrose	15.441	0.567	16.814	0.923	0.093		12.222	0.380	14.140	0.913	0.028	*
Maltose	1.362	0.051	1.589	0.067	0.009	**	1.346	0.141	1.394	0.055	0.611	
Melibiose	1.117	0.100	1.201	0.124	0.410		1.043	0.182	1.077	0.055	0.771	
Mannobiose	0.109	0.022	0.136	0.029	0.269		0.122	0.029	0.158	0.037	0.261	
Cellobiose	1.328	0.086	1.172	0.167	0.223		1.020	0.106	1.107	0.094	0.350	
Lactose	1.510	0.097	1.694	0.066	0.054		1.316	0.155	1.390	0.044	0.470	
Lactitol	2.477	0.220	2.957	0.078	0.023	*	2.171	0.225	2.270	0.060	0.506	
3-o-coumaroul-D-Quinic acid	0.003	0.004	0.050	0.061	0.248		0.001	0.001	0.048	0.057	0.227	
Sugars	344.122	12.127	361.596	28.196	0.380		326.208	13.524	315.286	4.826	0.258	
Alcohols	85.761	4.404	87.609	11.141	0.803		77.008	5.046	72.418	2.180	0.222	
Acids	7.660	0.699	9.425	0.466	0.022	*	8.796	0.154	8.826	0.326	0.895	
Other compounds	0.898	0.081	1.077	0.179	0.189		1.004	0.083	1.192	0.069	0.039	*
Amino acids	0.026	0.017	0.024	0.010	0.896		0.027	0.011	0.021	0.008	0.474	

Table S4. PCA extraction variables.

Variables	Extraction												
P	0.864	P	0.865	P	0.858	P	0.866	P	0.867	P	0.884	P	0.891
K	0.227	K	0.231	K	0.232	K	0.212	K	0.212	K	0.648	K	0.638
Ca	0.69	Ca	0.685	Ca	0.682	Ca	0.659	Ca	0.659	Ca	0.84	Ca	0.825
Mg	0.808	Mg	0.811	Mg	0.836	Mg	0.834	Mg	0.831	Mg	0.531	Mg	0.529
Na	0.368	Na	0.372	Na	0.412	Na	0.486	Na	0.523	Na	0.531	Na	0.529
Zn	0.547	Zn	0.545	Zn	0.529	Zn	0.521	Zn	0.547	Zn	0.544	Zn	0.571
Cu	0.651	Cu	0.675	Cu	0.717	Cu	0.738	Cu	0.731	Cu	0.737	Cu	0.738
Mn	0.783	Mn	0.782	Mn	0.784	Mn	0.788	Mn	0.812	Mn	0.815	Mn	0.8
Fe	0.439	Fe	0.437	Fe	0.421	Fe	0.431	Fe	0.46	Fe	0.45	Fe	0.466
Anthocyanins	0.745	Anthocyanins	0.741	Anthocyanins	0.743	Anthocyanins	0.744	Anthocyanins	0.747	Anthocyanins	0.758	Anthocyanins	0.751
Flavonols	0.061	Hydroxycinnamic acid	0.181	Polyphenols	0.461	Polyphenols	0.464	Polyphenols	0.471	Polyphenols	0.475	Polyphenols	0.471
Hydroxycinnamic acid	0.204	Polyphenols	0.442	Christensen water abs	0.298	Christensen water abs	0.262						
Polyphenols	0.44	Christensen water abs	0.303	Christensen cracking	0.392	Christensen cracking	0.433	Christensen cracking	0.424	Christensen cracking	0.424	Christensen cracking	0.424
Christensen water abs	0.305	Christensen cracking	0.381	Waterfall water abs	0.395	Waterfall water abs	0.418	Waterfall water abs	0.415	Waterfall water abs	0.415	Waterfall water abs	0.415
Christensen cracking	0.375	Waterfall water abs	0.396	Waterfall cracking	0.441	Waterfall cracking	0.452	Waterfall cracking	0.448	Waterfall cracking	0.448	Waterfall cracking	0.448
Waterfall water abs	0.39	Waterfall cracking	0.445	Cracking in the field	0.744	Cracking in the field	0.742	Cracking in the field	0.744	Cracking in the field	0.744	Cracking in the field	0.744
Waterfall cracking	0.437	Cracking in the field	0.739	Dry weight	0.649	Dry weight	0.656	Dry weight	0.628	Dry weight	0.631	Dry weight	0.611
Cracking in the field	0.726	Dry weight	0.65	Penetration	0.417	Penetration	0.402	Penetration	0.409	Penetration	0.388	Penetration	0.386
Dry weight	0.653	Penetration	0.416	Respiration rate	0.849	Respiration rate	0.796	Respiration rate	0.754	Respiration rate	0.757	Respiration rate	0.757
Penetration	0.414	Respiration rate	0.857	Ratio flesh to fruit	0.291	Ratio flesh to fruit	0.291	Ratio flesh to fruit	0.272	Ratio flesh to fruit	0.274	Ratio flesh to fruit	0.274
Respiration rate	0.844	Ratio flesh to fruit	0.282	TSS	0.394	TSS	0.404	Ratio flesh to fruit	0.272	TSS	0.4	TSS	0.386
Ratio flesh to fruit	0.28	TSS	0.402	TA	0.201	TA	0.201	TSS	0.42	TA	0.274	TA	0.274
TSS	0.411	TA	0.252										
TA	0.274												

The lowest variable extraction was rejected from the model until all remaining variables to have extraction higher than 0.5.

Variables	Extraction	Variables	Extraction	Variables	Extraction	Variables	Extraction
P	0.883	P	0.912	P	0.929	P	0.932
Ca	0.634	Ca	0.604	Ca	0.6	Ca	0.623
Mg	0.827	Mg	0.839	Mg	0.871	Mg	0.877
Na	0.483	Na	0.48	Na	0.462	Na	0.614
Zn	0.573	Zn	0.597	Zn	0.591	Zn	0.614
Cu	0.776	Cu	0.776	Cu	0.77	Cu	0.769
Mn	0.801	Mn	0.787	Mn	0.76	Mn	0.743
Fe	0.464	Fe	0.424	Anthocyanins	0.744	Anthocyanins	0.746
Anthocyanins	0.728	Anthocyanins	0.74	Polyphenols	0.516	Polyphenols	0.509
Polyphenols	0.472	Polyphenols	0.494	Christensen cracking	0.57	Christensen cracking	0.56
Christensen cracking	0.538	Christensen cracking	0.568	Waterfall water abs	0.537	Waterfall water abs	0.542
Waterfall water abs	0.514	Waterfall water abs	0.531	Waterfall cracking	0.505	Waterfall cracking	0.515
Waterfall cracking	0.52	Waterfall cracking	0.514	Cracking in the field	0.644	Cracking in the field	0.622
Cracking in the field	0.643	Cracking in the field	0.635	Dry weight	0.608	Dry weight	0.617
Dry weight	0.588	Dry weight	0.593	Respiration rate	0.781	Respiration rate	0.84
Penetration	0.408	Respiration rate	0.77				
Respiration rate	0.78						

Table S5. Quantitative results of nutrients in fruits.

Phosphorus (P, % DW)				Potassium (K, % DW)				Magnesium (Mg, % DW)				Sodium (Na, % DW)				Iron (Fe, ppm)			
		Mean	SD			Mean	SD			Mean	SD			Mean	SD			Mean	SD
Treatment (T)	Control	0.29	0.04	Control	1.32	0.17	Control	0.06	0.01	Control	0.02	0	Control	18.98	3.11	Control	18.98	3.11	
	CaCl ₂ * 0.25M	0.3	0.05	CaCl ₂ * 0.25M	1.28	0.2	CaCl ₂ * 0.25M	0.06	0.01	CaCl ₂ * 0.25M	0.02	0	CaCl ₂ * 0.25M	19.93	3.73	CaCl ₂ * 0.25M	19.93	3.73	
	2H ₂ O 0.5M	0.29	0.03	2H ₂ O 0.5M	1.32	0.16	2H ₂ O 0.5M	0.05	0.01	2H ₂ O 0.5M	0.02	0	2H ₂ O 0.5M	17.73	3.61	2H ₂ O 0.5M	17.73	3.61	
	1M	0.29	0.03	1M	1.17	0.12	1M	0.05	0.01	1M	0.02	0	1M	18.69	1.48	1M	18.69	1.48	
Age of Flower spur (AS)	1st buds	0.29	0.04	1st buds	1.28	0.19	1st buds	0.06	0.01	1st buds	0.02	0	1st buds	18.84	3.68	1st buds	18.84	3.68	
	2nd buds	0.29	0.03	2nd buds	1.27	0.16	2nd buds	0.06	0.01	2nd buds	0.02	0	2nd buds	18.82	2.53	2nd buds	18.82	2.53	
Ripening (R)	Color Red	0.32	0.03	Color Red	1.32	0.15	Color Red	0.06	0.01	Color Red	0.02	0	Color Red	19.17	2.27	Color Red	19.17	2.27	
	Color Black	0.27	0.02	Color Black	1.22	0.18	Color Black	0.05	0	Color Black	0.02	0	Color Black	18.49	3.81	Color Black	18.49	3.81	

Factors	P-value	N
T	0.755	12
AS	0.461	24
R	<0.001 ***	24
T*AS	0.655	6
T*R	0.5	6
AS*R	0.078	12
T*AS*R	0.807	3

Factors	P-value	N
T	0.051	12
AS	0.858	24
R	0.021 *	24
T*AS	0.611	6
T*R	0.119	6
AS*R	0.03 *	12
T*AS*R	0.109	3

Factors	P-value	N
T	0.169	12
AS	0.609	24
R	<0.001 ***	24
T*AS	0.732	6
T*R	0.615	6
AS*R	0.157	12
T*AS*R	0.618	3

Factors	P-value	N
T	0.489	12
AS	0.021 *	24
R	0.58	24
T*AS	0.301	6
T*R	0.679	6
AS*R	0.758	12
T*AS*R	0.772	3

Factors	P-value	N
T	0.333	12
AS	0.978	24
R	0.419	24
T*AS	0.081	6
T*R	0.088	6
AS*R	0.576	12
T*AS*R	0.314	3

Different letters indicate significant difference based on Duncan's Multiple Range Test $P \leq 0.05$

Different letters indicate significant difference based on Duncan's Multiple Range Test $P \leq 0.05$