

Natural Variation in Herbivore-Induced Volatiles in *Arabidopsis thaliana*

Tjeerd A.L. Snoeren, Iris F. Kappers, Colette Broekgaarden, Roland Mumm,
Marcel Dicke and Harro J. Bouwmeester

Supplementary material

Table S1. Characteristics of the identified induced volatile compounds in headspace of *Arabidopsis thaliana* leaves after *Pieris rapae* herbivory or jasmonic acid treatment.

Compound	Retention index	<i>m/z</i> fragment used for quantification
pentan-2-ol	732	45
2-heptene	735	55
(E)-3-penten-2-one	735	69
2-pentenal	758	55
1-pentanol	767	42
(E)-3-hexenal	770	41
(Z)-2-penten-1-ol	771	57
(E)-3-penten-2-ol	774	71
(Z)-3-hexenal	796	41
hexan-2-ol	801	45
(E)-3-hexen-1-ol	851	41
(E)-2-hexenal	853	41
(Z)-3-hexen-1-ol	853	67
hexanol	868	56
hexanal	878	44

1-nonene	892	56
4-methyl-3-penten-2-one	892	83
heptanal	900	70
heptan-2-ol	908	45
(E,E)-2,4-hexadienal,	910	81
α -pinene	939	93
1-octen-3-ol	943	57
benzaldehyde	965	77
trans pinane	974	55
1-octen-1-ol	977	57
6-methyl-5-hepten-2-one	989	69
β -myrcene	995	93
octanal	1000	43
(E)-hexenyl acetate	1006	43
α -phellandrene	1007	93
(Z)-3-hexenyl acetate	1010	43
3-carene	1010	93
α -terpinene	1017	93
limonene	1033	93
(Z)- β -ocimene	1040	93
(E)- β -ocimene	1052	93
γ -terpinene	1062	93
acetophenone	1068	105
hexanoic acid	1085	60
terpinolene	1093	93
linalool	1099	93
nonanal	1109	57
DMNT ¹	1114	69
allo-ocimene	1129	69

β -sesquiphellandrene	1149	69
α -terpineol	1186	59
MeSA ³	1193	92+120+150
verbenone	1204	107
decanal	1207	57
nerol	1225	69
β -citral	1239	69
pulegone	1240	81
ethylsalicylate	1248	120
geraniol	1255	69
indole	1292	117
α -copaene	1378	161
β -elemene	1379	81
(E)- β -caryophyllene	1420	93
thujopsene	1431	119
geranyl acetone	1454	69
α -himachalene	1454	93
α -humulene	1454	107
(E)- β -farnesene	1464	69
β -acoradiene	1464	119
β -chamigrene	1484	189
β -ionone	1486	121
cuparene	1506	132
(E,E)- α -farnesene	1509	93
β -bisabolene	1510	69
(Z)-nerolidol	1535	69
geranylisovalerate	1553	85
(E)-nerolidol	1564	69

TM ¹ T ²	1590	69
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¹, (*E*)-4,8-dimethyl-1,3,7-nonatriene

², (*E,E*)-4,8,12-trimethyltrideca-1,3,7,11-tetraene

³, methyl salicylate

Table S2. Sequences of *Arabidopsis thaliana*-derived primers used in quantitative Real Time-PCR analyses.

<i>Gene</i>	<i>AGI-codes</i>	<i>Forward primer (5' → 3')</i>	<i>Reverse primer (5' → 3')</i>
β -Actin	At3g46520	GGAGAAGATTGGCATCACAC	TGGCAACATACATAGCAGGAG
<i>HPL1</i>	At4g15440	ACATCGCTGAGAACGGTTG	CAAGAGGCTGAGGAACTAGC
<i>BSMT1</i>	At3g11480	TGGTCACTACTACGAAGAAGATG	GAGCATTGGTTCACTAACAGC
<i>TPS03</i>	At4g16740	GCCACCATCCTCCGTCTC	CCAAGCCACACCGATAATTCC
<i>TPS04</i>	At1g61120	TCGCAGCACACACCATTG	GAGCAGCACGGAGTTCATC
<i>CYP72A13</i>	At3g14660	GATGGCAATGACACTGATTCTAC	GATAAGAGGAGCACCGAACTG
<i>CYP82G1</i>	At3g25180	ATCAGACAGCACATCCATCAC	GCCGACACTATTATCAATCTCTC

Table S3. Results of one-way ANOVA for each compound on differences between JA and *Pieris rapae* treatment. For each accession the volatile compounds that are emitted in significantly different (ANOVA) rates between both treatments ((*Pieris rapae* infestation (Pr) and JA-treatment (JA)). F- and P-values (Sig.) are given.

An-1	Treatment with higher average emission rate:		F	Sig.
(E)-2-hexenal	Pr		9.170	0.016
trans pinane	Pr		9.305	0.016
β-myrcene		JA	11.714	0.009
α-terpineol	Pr		9.712	0.014
β-acoradiene	Pr		11.409	0.010
(E)-nerolidol		JA	9.207	0.016
TMTT ¹		JA	5.632	0.045
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Col-0	Treatment with higher average emission rate:		F	Sig.
DMNT ²	Pr		10.059	0.010
MeSA ³	Pr		102.577	0.000
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C-24	Treatment with higher average emission rate:		F	Sig.
pentan-2-ol		JA	15.933	0.004
(E)-2-hexenal	Pr		52.706	0.000
4-Me-3-penten-2-one		JA	17.129	0.003
benzaldehyde		JA	8.315	0.020
(E)-β-ocimene		JA	7.592	0.025
β-sesquiphellandrene		JA	6.925	0.030
α-terpineol		JA	14.910	0.005

MeSA ³	Pr		10.119	0.013
nerol		JA	50.385	0.000
ethylsalicylate	Pr		12.938	0.007
indole		JA	448.516	0.000
β-elemene		JA	6.791	0.031
β-farnesene		JA	57.984	0.000
(E)-nerolidol		JA	10.805	0.011
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Cvi	Treatment with higher average emission rate:		F	Sig.
(E)-3-hexenal	Pr		14.663	0.005
(Z)-2-penten-1-ol		JA	6.614	0.033
(Z)-3-hexenal	Pr		7.115	0.028
(E)-2-hexenal	Pr		7.892	0.023
trans pinane	Pr		5.357	0.049
α-phellandrene	Pr		21.808	0.002
(Z)-hexenyl acetate	Pr		10.514	0.012
3-carene		JA	13.407	0.006
limonene	Pr		28.407	0.001
(Z)-β-ocimene		JA	8.471	0.020
acetophenone		JA	9.238	0.016
MeSA ³	Pr		24.951	0.001
β-farnesene	Pr		6.445	0.035
(E)-nerolidol		JA	7.072	0.029
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Eri-1	Treatment with higher average emission rate:		F	Sig.
2-pentanal	Pr		18.740	0.003
(E)-2-hexenal		JA	14.719	0.005

(Z)-3-hexen-1-ol	Pr		5.484	0.047
1-octen-1-ol	Pr		7.716	0.024
β -myrcene		JA	10.475	0.012
linalool	Pr		8.715	0.018
MeSA ³	Pr		61.541	0.000
nerol	Pr		6.343	0.036
α -farnesene		JA	30.832	0.001
β -bisbolene		JA	7.450	0.026
Kond	Treatment with higher average emission rate:		F	Sig.
(Z)-3-hexen-1-ol	Pr		17.445	0.003
hexanol	Pr		5.420	0.048
β -myrcene		JA	32.530	0.000
(Z)- β -ocimene		JA	6.367	0.036
verbenone		JA	6.500	0.034
decanal		JA	17.427	0.003
(Z)-nerolidol		JA	10.927	0.011
Kyo-1	Treatment with higher average emission rate:		F	Sig.
(Z)-nerolidol		JA	15.012	0.005
Ler	Treatment with higher average emission rate:		F	Sig.
β -myrcene	Pr		9.877	0.014
(E)-hexenyl acetate	Pr		19.068	0.002
(Z)- β -ocimene	Pr		8.886	0.018

hexanoic acid	Pr		5.999	0.040
ethylsalicylate		JA	6.290	0.036
WS				
	Treatment with higher average emission rate:		F	Sig.
(E,E)-2,4-hexadienal		JA	12.120	0.006
(Z)-hexenyl acetate		JA	5.309	0.044
linalool		JA	5.361	0.043
geranyl acteone		JA	6.643	0.028
α -farnesene		JA	5.447	0.042

¹, (E,E)-4,8,12-trimethyltrideca-1,3,7,11-tetraene

², (E)-4,8-dimethyl-1,3,7-nonatriene

³, methyl salicylate

Table S4. Volatile compounds that are emitted in significantly different rates among nine accessions of *Arabidopsis thaliana* treated with JA, infested with 20 first instar larvae of *Pieris rapae* or left untreated (control). Individual volatiles were analyzed for significant differences between accessions within one treatment, by using one-way ANOVA followed by a Dunnett T3 post-hoc analysis. Significance: * $P<0.05$; ** $P<0.01$; *** $P<0.001$.

Compound	Control	Sign.	<i>P. rapae</i>	Sign.	JA	Sign.
(E)-2-hexenal			An-1 > Ler	*		
			C-24 > Ler	*		
(E)-3-hexenal	WS > Col-0	**				
	WS > C-24	*				
(Z)-hexenyl acetate			Cvi > An-1	***		
			Cvi > C-24	***		
(Z)-2-penten-1-ol	WS > Col-0	*				
3-hexen-1-ol			An-1 > Col-0	***		
			An-1 > Cvi	*		
			Cvi > Col-0	*		
			Eri-1 > Col-0	**		
			Kond > C-24	*		
			Kond > Col-0	***		
			Kond > Cvi	**		
hexanoic acid			WS > Col-0	***		
	C-24 > Ler	***				
pentan-2-ol	C-24 > Kyo-1	*				
	Cvi > Kyo-1	*				
(E)-nerolidol	C-24 > Col-0	*	Col-0 > WS	*		
			Kond > WS	*		

(E)-β-ocimene			Kond > C-24	***	An-1 > Col-0	*
			Kond > Col-0	**	An-1 > Kyo-1	**
			WS > Col-0	*	C-24 > Kyo-1	*
			WS > C-24	**	Cvi > Kyo-1	*
					Eri-1 > Kyo-1	*
					Kond > Col-0	***
					Kond > Cvi	*
					Kond > Eri-1	***
					Kond > Kyo-1	***
					WS > Col-0	***
					WS > Eri-1	**
					WS > Kyo-1	***
(Z)-β-ocimene	Kond > An-1	**	An-1 > Col-0	*	An-1 > Kyo-1	*
	Kond > Cvi	*	Kond > Col-0	*	C-24 > Col-0	*
			Ler > C-24	*	C-24 > Kyo-1	*
			Ler > Col-0	**	Cvi > Col-0	*
			WS > Col-0	*	Cvi > Kyo-1	**
			WS > C-24	*	Kond > Col-0	**
					Kond > Eri-1	*
					Kond > Kyo-1	***
					Kond > Ler	*
					WS > Col-0	*
					WS > Kyo-1	**
3-carene	An-1 > Col-0	***				

β-acoradiene			Cvi > Kyo-1	*		
DMNT	C-24 > Col-0	***			C-24 > Col-0	**
	Kond > Col-0	***			Kond > Col-0	*
	WS > Col-0	**			WS > Col-0	***
β-bisabolene			An-1 > Kond	***	C-24 > WS	*
			C-24 > Kond	***	Eri-1 > WS	*
			Kyo-1 >Kond	***		
β-farnesene	Col-0 > Ler	**	Col-0 > C-24	***		
	Kond > Ler	*	Cvi > C-24	*		
			Cvi > C-24	*		
			Kond > C-24	**		
			Kyo-1 > C-24	***		
			WS > C-24	*		
β-myrcene			Cvi > An-1	*	Cvi > An-1	*
			Eri-1 > An-1	*	Cvi > Col-0	*
			Kond > An-1	*	Cvi > Ler	*
			Ler > An-1	*	Eri-1 > Col-0	*
β-myrcene					Eri-1 > Ler	*
					Kond > An-1	***
					Kond > C-24	**
					Kond > Col-0	***
					Kond > Ler	**
					Kond > WS	*
β-sesquiphellandrene			An-1 > Eri-1	*	An-1 > Cvi	*
			An-1 > Col-0	*	An-1 > Eri-1	*

			An-1 > C-24	*	Kond > Cvi	*
					Kond > Eri-1	*
TMTT			Col-0 > C-24	**	An-1 > Cvi	*
			Col-0 > Cvi	**	Kond > Cvi	**
			Kond > C-24	*	Kyo-1 > Cvi	***
			Kond > Cvi	*	Ler > Cvi	**
			WS > Cvi	**		
			WS > C-24	*		
α-farnesene			An-1 > Cvi	**	Eri-1 > C-24	**
			An-1 > Col-0	*	Eri-1 > WS	**
			An-1 > C-24	*	Kond > C-24	*
			Eri-1 > Cvi	**		
			Kyo-1 > C-24	*		
			Kyo-1 > Col-0	*		
			Kyo-1 > Cvi	***		
			Kyo-1 > Ler	*		
α-humulene			C-24 > Kyo-1	**		
α-phellandrene	Cvi > Col-0	*				
	Kond > Col-0	*				
α-pinene					Kond > An-1	*
					Ler > An-1	*
benzaldehyde			An-1 > C-24	**		
			Col-0 > C-24	*		
			Kond > C-24	*		

			<i>Ler</i> > C-24	**		
decanal			Col-0 > An-1	*		
			Col-0 > Cvi	*		
			Eri-1 > An-1	*		
			Eri-1 > Cvi	*		
			WS > An-1	***		
			WS > Cvi	**		
ethylsalicylate					Cvi > An-1	***
					Cvi > Eri-1	***
					<i>Ler</i> > An-1	***
					<i>Ler</i> > Eri-1	***
geranyl acetone			Kond > Col-0	*	WS > Eri-1	*
indole			Cvi > C-24	***		
MeSA	An-1 > WS	*	An-1 > Eri-1	*	Col-0 > Cvi	***
	Eri-1 > WS	*	An-1 > Kond	**	Col-0 > Eri-1	*
	Kyo-1 > WS	**	An-1 > Kyo-1	*	Kond > Cvi	**
			An-1 > Cvi	***	Kyo-1 > Cvi	**
			C-24 > Cvi	***	<i>Ler</i> > Cvi	***
			Col-0 > Cvi	***	<i>Ler</i> > Eri-1	**
			Col-0 > Eri-1	*	<i>Ler</i> > Kond	*
			Col-0 > Kond	**	<i>Ler</i> > Kyo-1	*
			Col-0 > Kyo-1	*	WS > Cvi	***
			Eri-1 > C-24	***	Ws > Eri-1	**
			Kyo-1 > Cvi	**		
			<i>Ler</i> > Cvi	***		

			Ler > Kond	*		
			WS > Cvi	***		

Table S5. Significant differences in headspace among accessions after either treatment or left non-treated. One-way ANOVA followed by Dunnett T3 results for 9 *Arabidopsis thaliana* accessions for each volatile compound.

<i>Pieris rapae</i>					
Geno I	>Geno J	Compound	Mean¹ Difference (I-J)	Std. Error	Sign.
An-1	> Col-0	3-hexen-1ol	1.945	0.120	<0.001
		(Z)-β-ocimene	0.789	0.183	0.045
		β-sesquiphellandrene	1.149	0.256	0.042
		α-farnesene	1.116	0.243	0.038
	> C24	benzaldehyde	1.448	0.226	0.006
		β-sesquiphellandrene	1.262	0.285	0.047
		α-farnesene	1.314	0.297	0.049
	> Cvi	3-hexen-1ol	0.900	0.174	0.024
		MeSA	1.704	0.128	<0.001
		α-farnesene	1.724	0.225	0.003
	> Eri-1	β-sesquiphellandrene	1.280	0.240	0.022
		MeSA	0.524	0.103	0.026
	> Kond	MeSA	1.152	0.149	0.002
		β-bisbolene	7.448	0.226	<0.001
	> Kyo-1	MeSA	0.808	0.140	0.012
	> Ler	(E)-2-hexenal	7.080	1.226	0.045
Col-0	> An-1	decanal	1.352	0.256	0.027
	> Cvi	MeSA	1.602	0.106	<0.001
		decanal	1.145	0.285	0.019
	> Eri-1	MeSA	0.422	0.074	0.013
	> C24	benzaldehyde	1.155	0.192	0.024

		β -farnesene	1.043	0.067	<0.001
		TMTT	1.454	0.252	0.008
	> Cvi	TMTT	1.558	0.240	0.005
	> Kond	MeSA	1.050	0.130	0.008
	> Kyo-1	MeSA	0.706	0.120	0.030
	> WS	(E)-nerolidol	6.193	1.194	0.047
C24	> Cvi	MeSA	1.182	0.135	0.001
	> Kond	β -bisbolene	7.166	0.312	<0.001
	> Kyo-1	α -humulene	1.294	0.170	0.002
	> Ler	(E)-2-hexenal	7.148	1.202	0.047
Cvi	> An-1	β -myrcene	0.836	0.160	0.019
		(Z)-hexenyl acetate	1.442	0.126	<0.001
	> Col-0	3-hexen-1ol	1.045	0.155	0.012
	> C24	(Z)-hexenyl acetate	1.620	0.163	0.001
		indole	6.170	0.312	0.001
		β -farnesene	1.104	0.144	0.018
		β -farnesene	1.104	0.144	0.018
	> Kyo-1	β -acoradiene	0.564	0.119	0.032
Eri-1	> An-1	decanal	1.066	0.218	0.050
		β -myrcene	0.622	0.138	0.044
	> Col-0	3-hexen-1ol	1.899	0.180	0.002
	> C24	MeSA	1.180	0.115	0.001
	> Cvi	decanal	1.164	0.251	0.040
		α -farnesene	1.286	0.183	0.003
Kond	> An-1	β -myrcene	0.846	0.156	0.015
	> Col-0	3-hexen-1ol	2.313	0.156	<0.001
		(Z)- β -ocimene	1.005	0.196	0.016
		(E)- β -ocimene	1.202	0.183	0.007
		geranyl acetone	6.275	1.150	0.041

	> C24	3-hexen-1ol	2.334	0.426	0.041
		benzaldehyde	1.158	0.213	0.020
		(E)-β-ocimene	1.274	0.108	<0.001
		β-farnesene	1.096	0.084	0.002
		TMTT	1.188	0.231	0.022
	> Cvi	3-hexen-1ol	1.268	0.201	0.005
		TMTT	1.292	0.218	0.012
	> WS	(E)-nerolidol	6.190	1.187	0.048
Kyo-1	> Col-0	α-farnesene	1.106	0.205	0.011
	> C24	β-farnesene	1.048	0.687	0.001
		α-farnesene	1.304	0.266	0.040
	> Cvi	MeSA	0.896	0.149	0.008
		α-farnesene	1.714	0.183	<0.001
	> Kond	β-bisbolene	5.790	0.343	0.001
	> Ler	α-farnesene	1.096	0.233	0.040
Ler	> An-1	β-myrcene	0.834	0.151	0.013
	> Col-0	(Z)-β-ocimene	0.955	0.157	0.009
	> C24	benzaldehyde	1.380	0.207	0.008
	> Cvi	(Z)-β-ocimene	0.524	0.096	0.014
		MeSA	1.400	0.128	<0.001
	> Kond	MeSA	0.848	0.149	0.015
WS	> An-1	decanal	1.492	0.129	<0.001
	> Col-0	3-hexen-1ol	1.962	0.169	<0.001
	> C24	(Z)-β-ocimene	0.870	0.149	0.017
		(E)-β-ocimene	1.008	0.210	0.021
	> Cvi	(Z)-β-ocimene	0.439	0.082	0.015
		(E)-β-ocimene	1.080	0.149	0.002
		β-farnesene	0.918	0.132	0.014
		TMTT	0.917	0.196	0.030

		3-hexen-1ol	0.917	0.211	0.042
		MeSA	1.273	0.163	0.001
		decanal	1.590	0.179	0.002
jasmonic acid					
Geno I	>Geno J	Compound	Mean ¹ Difference (I-J)	Std. Error	Sign.
An-1	> Col-0	(E)-β-ocimene	0.957	0.192	0.031
	> Cvi	β-sesquiphellandrene	1.068	0.224	0.031
		TMTT	0.828	0.148	0.027
	> Eri-1	β-sesquiphellandrene	0.892	0.192	0.043
	> Kyo-1	(Z)-β-ocimene	1.052	0.143	0.012
		(E)-β-ocimene	1.392	0.205	0.004
Col-0	> Cvi	MeSA	1.505	0.573	<0.001
	> Eri-1	MeSA	1.253	0.180	0.020
C24	> Col-0	(Z)-β-ocimene	0.980	0.229	0.049
		DMNT	1.672	0.257	0.004
	> Kyo-1	(Z)-β-ocimene	1.082	0.145	0.011
		(E)-β-ocimene	1.098	0.227	0.037
	> WS	β-bisbolene	5.974	1.154	0.039
Cvi	> An-1	β-myrcene	0.638	0.126	0.025
		ethylsalicylate	5.564	0.343	0.001
	> Col-0	β-myrcene	0.885	0.168	0.014
		(Z)-β-ocimene	1.176	0.231	0.017
	> Eri-1	ethylsalicylate	5.564	0.343	0.001
	> Kyo-1	(Z)-β-ocimene	1.278	0.148	0.006
	> Kyo-1	(E)-β-ocimene	1.108	0.199	0.013
	> Ler	β-myrcene	0.734	0.147	0.024
Eri-1	> Col-0	β-myrcene	0.969	0.198	0.022

	> C24	α -farnesene	1.478	0.215	0.003	
	> Kyo-1	(<i>E</i>)- β -ocimene	0.816	0.154	0.023	
	> <i>Ler</i>	β -myrcene	0.818	0.180	0.047	
Kond	> WS	α -farnesene	1.205	0.209	0.007	
		β -bisbolene	7.034	1.140	0.019	
Kond	> An-1	α -pinene	0.686	0.115	0.021	
		β -myrcene	1.110	0.091	<0.001	
	> Col-0	β -myrcene	1.357	0.144	0.001	
		(<i>Z</i>)- β -ocimene	1.682	0.197	0.002	
		(<i>E</i>)- β -ocimene	1.531	0.115	<0.001	
		DMNT	0.950	0.189	0.023	
	> C24	β -myrcene	1.250	0.149	0.007	
		α -farnesene	1.266	0.244	0.019	
	> Cvi	(<i>E</i>)- β -ocimene	0.858	0.160	0.049	
		β -sesquiphellandrene	1.208	0.240	0.023	
		MeSA	1.074	0.099	0.002	
		TMTT	1.386	0.179	0.008	
Kyo-1	> Eri-1	(<i>Z</i>)- β -ocimene	0.954	0.165	0.024	
		(<i>E</i>)- β -ocimene	1.150	0.099	<0.001	
		β -sesquiphellandrene	1.032	0.210	0.037	
	> Kyo-1	(<i>Z</i>)- β -ocimene	1.784	0.085	<0.001	
	> Kyo-1	(<i>E</i>)- β -ocimene	1.966	0.136	<0.001	
Ler		β -myrcene	1.206	0.118	0.002	
		(<i>Z</i>)- β -ocimene	0.962	0.175	0.033	
		β -myrcene	0.857	0.143	0.017	
Kyo-1	> Cvi	MeSA	1.078	0.139	0.012	
		TMTT	1.108	0.099	<0.001	
Ler	> An-1	α -pinene	0.732	0.134	0.039	
		ethylsalicylate	6.228	0.153	<0.001	

	> Cvi	MeSA	1.900	0.120	<0.001
		TMTT	1.592	0.201	0.009
> Eri-1		MeSA	1.648	0.208	0.002
		ethylsalicylate	6.228	0.153	<0.001
> Kond		MeSA	0.826	0.147	0.013
> Kyo-1		MeSA	0.822	0.176	0.037
WS	>Col-0	(Z)-β-ocimene	1.257	0.242	0.011
		(E)-β-ocimene	1.298	0.154	<0.001
		DMNT	1.412	0.196	0.001
	> Cvi	MeSA	1.585	0.136	0.001
	> Eri-1	(E)-β-ocimene	0.918	0.142	0.003
		MeSA	1.333	0.218	0.007
		geranyl acetone	6.608	1.108	0.039
	> Kyo-1	(Z)-β-ocimene	1.359	0.164	0.004
		(E)-β-ocimene	1.734	0.170	<0.001
Control					
Geno I	> Geno J	Compound	Mean ¹ Difference (I-J)	Std. Error	Sign.
An-1	> Col-0	3-carene	5.740	0.199	<0.001
	> WS	MeSA	1.031	0.201	0.036
Col-0	> Ler	β-farnesene	0.922	0.155	0.007
C24	> Col-0	DMNT	1.030	0.155	0.001
		(E)-nerolidol	5.887	1.125	0.039
	> Kyo-1	pentan-2-ol	5.232	1.133	0.042
	> Ler	hexanoic acid	6.540	0.087	<0.001
Cvi	> Col-0	α-phellandrene	5.557	1.042	0.045
	> Kyo-1	pentan-2-ol	5.518	1.122	0.032
Eri-1	> WS	MeSA	0.789	0.171	0.045

Kond	> An-1	(Z)- β -ocimene	1.073	0.175	0.004
	> Col-0	α -phellandrene	5.479	1.042	0.048
		DMNT	0.913	0.133	0.001
	> Cvi	(Z)- β -ocimene	1.620	0.346	0.050
	> Ler	β -farnesene	0.748	0.163	0.029
Kyo-1	> WS	MeSA	0.707	0.131	0.005
WS	> Col-0	(E)-3-hexanal	1.757	0.280	0.009
		(Z) 2-penten-1-ol	6.405	1.199	0.036
		DMNT	0.902	0.153	0.005
	> C24	(E)-3-hexanal	1.598	0.278	0.015

Table S6. Results of t-tests for each volatile compound on differences between non-treated plants and *P. rapae*-infested plants, and non-treated plants and JA-treated plants.

	P. rapae	F	Sig.2tail	Jasmonic acid		F	Sig.2tail
An-1	(Z)- β -ocimene	0,023	2,83E-06	An-1	(Z)- β -ocimene	0,358	7,96E-06
An-1	MeSA	5,825	4,80E-04	An-1	(E)- β -ocimene	0,009	4,28E-04
An-1	α -farnesene	0,014	0,001	An-1	α -farnesene	0,000	0,002
An-1	geraniol	2,570	0,003	An-1	TMTT	1,503	0,004
An-1	α -terpineol	4,333	0,005	An-1	β -sesquiphellandrene	2,442	0,012
An-1	DMNT	4,379	0,010	An-1	geraniol	0,379	0,023
An-1	β -sesquiphellandrene	1,620	0,011	An-1	(Z)-nerolidol	2,488	0,044
An-1	(E)- β -ocimene	0,622	0,011	An-1	ethylsalicylate	44,108	0,075
An-1	(E)-2-hexenal	0,148	0,015	An-1	DMNT	2,074	0,078
An-1	(Z)-3-hexen-1-ol	85,282	0,044				
An-1	octanal	0,337	0,059				
An-1	(Z)-2-penten-1-ol	44,425	0,060				
An-1	nerol	4,382	0,067				
An-1	3-carene	56,757	0,072				
An-1	geranylisovalerate	1,025	0,084				
An-1	β -elemene	15,297	0,084				
An-1	allo ocimene	5,898	0,085				
An-1	A-terpinene	5,164	0,088				
An-1	β -bisbolene	62,390	0,089				
Col-0	MeSA	4,207	2,32E-07	Col-0	MeSA	3,887	7,57E-06
Col-0	(E)-nerolidol	3,933	2,02E-04	Col-0	(Z)-2-penten-1-ol	3,806	1,60E-04
Col-0	TMTT	2,574	0,001	Col-0	(E)-nerolidol	4,810	2,10E-04
Col-0	(Z)- β -ocimene	0,045	0,002	Col-0	α -phellandrene	0,000	0,018
Col-0	DMNT	0,471	0,004	Col-0	TMTT	1,359	0,045
Col-0	linalool	1,299	0,052	Col-0	α -farnesene	297,446	0,048

Col-0	(Z)-3-hexen-1-ol	458,342	0,069	Col-0	(Z)- β -ocimene	0,864	0,051
Col-0	decanal	0,418	0,069	Col-0	3-carene	2.939,156	0,076
Col-0	α -farnesene	586,641	0,072				
Col-0	(Z)-2-penten-1-ol	2,332	0,077				
Col-0	(E)-hexenyl acetate	25,916	0,094				
C-24	pentan-2-ol	6,816	4,41E-08	C-24	(E)- β -ocimene	1,934	8,03E-06
C-24	MeSA	6,603	3,90E-06	C-24	(Z)- β -ocimene	0,594	1,49E-04
C-24	(E)- β -ocimene	0,214	8,48E-06	C-24	MeSA	0,165	0,002
C-24	hexanoic acid	3,265	5,23E-05	C-24	DMNT	0,156	0,014
C-24	(Z)- β -ocimene	3,796	3,48E-04	C-24	ethylsalicylate	0,112	0,016
C-24	(E)-2-hexenal	0,496	4,28E-04	C-24	(E)-hexenyl acetate	0,158	0,027
C-24	4-me-3-penten-2-one	3,446	0,002	C-24	6-me-5-hepten-2-one	1,724	0,038
C-24	1-pentanol	3,614	0,008	C-24	β -farnesene	1,500	0,061
C-24	indole	16,298	0,009	C-24	α -humulene	0,296	0,098
C-24	11-octen-1-ol	0,014	0,011				
C-24	α -terpineol	3,710	0,013				
C-24	benzaldehyde	0,103	0,016				
C-24	6-me-5-hepten-2-one	0,082	0,016				
C-24	(E)-nerolidol	0,089	0,016				
C-24	α -humulene	0,165	0,024				
C-24	β -myrcene	0,041	0,027				
C-24	β -chamigrene	1,980	0,031				
C-24	β -farnesene	9,643	0,035				
C-24	(E)- β -caryophyllene	0,100	0,040				
C-24	thujopsene	31,742	0,042				
C-24	(Z)-3-hexen-1-ol	0,540	0,049				
C-24	(E)-hexenyl acetate	0,205	0,064				
C-24	β -elemene	1,867	0,076				
Cvi	(Z)-hexenyl acetate	1,238	8,15E-07	Cvi	(Z)- β -ocimene	1,580	7,33E-05

Cvi	(Z)- β -ocimene	0,957	0,001	Cvi	(E)- β -ocimene	0,082	1,17E-04
Cvi	trans pinane	2,702	0,003	Cvi	β -myrcene	2,246	4,22E-04
Cvi	β -myrcene	0,834	0,004	Cvi	linalool	0,365	0,004
Cvi	α -phellandrene	0,097	0,005	Cvi	(Z)-hexenyl acetate	1,243	0,005
Cvi	linalool	0,785	0,007	Cvi	1-pentanol	2,419	0,020
Cvi	1-pentanol	0,466	0,014	Cvi	(E)-nerolidol	133,242	0,029
Cvi	(E)- β -ocimene	1,801	0,016	Cvi	α -phellandrene	2,858	0,030
Cvi	β -acoradiene	2,343	0,028	Cvi	acetophenone	1,304	0,035
Cvi	β -farnesene	0,002	0,037	Cvi	(E)-3-hexenal	855,557	0,051
Cvi	geranyl acteone	61,472	0,042	Cvi	3-carene	206,499	0,052
Cvi	(E)- β -caryophyllene	1,701	0,059	Cvi	(Z)-2-penten-1-ol	26,958	0,055
Cvi	benzaldehyde	3,059	0,064	Cvi	g terpinene	1,559	0,057
Cvi	α -terpinene	0,065	0,065	Cvi	geranyl acteone	116,002	0,071
Cvi	MeSA	2,537	0,068				
Cvi	α -farnesene	447,986	0,069				
Cvi	hexan-2-ol	0,202	0,074				
Cvi	pentan-2-ol	92,570	0,085				
Cvi	(E)-2-hexanal	3,809	0,094				
Cvi	verbenone	0,148	0,096				
Cvi	(E)-3-hexanal	1.117,670	0,096				
Cvi	(E)-nerolidol	42,628	0,100				
Eri-1	MeSA	2,004	8,59E-05	Eri-1	α -farnesene	0,004	4,23E-06
Eri-1	(Z)- β -ocimene	0,895	3,26E-04	Eri-1	(E)- β -ocimene	2,622	0,002
Eri-1	(E)- β -ocimene	0,023	0,001	Eri-1	(Z)- β -ocimene	0,250	0,003
Eri-1	α -farnesene	0,059	0,001	Eri-1	α -humulene	0,336	0,005
Eri-1	DMNT	2,589	0,005	Eri-1	DMNT	1,995	0,006
Eri-1	linalool	0,778	0,011	Eri-1	TMTT	0,012	0,038
Eri-1	TMTT	0,124	0,022	Eri-1	β -bisbolene	47,132	0,063
Eri-1	α -humulene	0,088	0,038	Eri-1	β -farnesene	0,141	0,068

Eri-1	(Z)-3-hexen-1-ol	68,315	0,050	Eri-1	ethylsalicylate	60,631	0,073
Eri-1	octanal	74,681	0,071	Eri-1	(E)-2-hexenal	47,663	0,081
Eri-1	heptanal	0,250	0,076	Eri-1	heptanal	0,018	0,087
				Eri-1	β-myrcene	78,719	0,093
Kond	(E)-β-ocimene	1,150	6,73E-06	Kond	(E)-β-ocimene	3,553	1,63E-06
Kond	(Z)-β-ocimene	0,022	2,57E-04	Kond	(Z)-β-ocimene	2,020	5,03E-06
Kond	(Z)-nerolidol	0,220	0,001	Kond	(Z)-nerolidol	0,100	2,55E-05
Kond	β-bisbolene	4,718	0,002	Kond	β-myrcene	3,208	4,78E-04
Kond	2-pentanal	0,182	0,007	Kond	MeSA	1,399	0,005
Kond	MeSA	0,428	0,007	Kond	α-pinene	0,581	0,025
Kond	β-elemene	0,085	0,017	Kond	β-farnesene	0,020	0,027
Kond	(E)-hexenyl acetate	1,274	0,020	Kond	β-bisbolene	0,011	0,029
Kond	1-nonene	4,542	0,032	Kond	(E)-3-hexen-1-ol	0,108	0,043
Kond	β-myrcene	0,966	0,035	Kond	β-sesquiphellandrene	0,020	0,045
Kond	TMTT	3,853	0,065	Kond	α-farnesene	3,265	0,062
Kond	(Z)-3-hexen-1-ol	26,140	0,065	Kond	TMTT	3,626	0,066
Kond	A-pinene	0,004	0,073	Kond	α-phellandrene	0,147	0,088
Kond	(E)-3-hexenal	21,423	0,076				
Kyo-1	MeSA	0,001	5,30E-05	Kyo-1	MeSA	0,046	0,003
Kyo-1	α-farnesene	46,460	0,012	Kyo-1	(Z)-nerolidol	0,420	0,005
Kyo-1	α-himachalene	155,770	0,019	Kyo-1	α-farnesene	37,732	0,009
Kyo-1	limonene	39,446	0,020	Kyo-1	(E,E)-2,4-hexadienal	0,902	0,019
Kyo-1	hexanoic acid	76,421	0,041	Kyo-1	α-himachalene	116,646	0,031
Kyo-1	trans pinane	422,776	0,048	Kyo-1	trans pinane	1.675,678	0,055
Kyo-1	terpinolene	0,120	0,064	Kyo-1	(E)-β-ocimene	4,216	0,057
				Kyo-1	β-myrcene	8,423	0,061
				Kyo-1	terpinolene	0,033	0,076
Ler	hexanoic acid	16,710	1,38E-04	Ler	MeSA	5,625	0,001
Ler	(E)-β-ocimene	0,800	0,001	Ler	(E)-β-ocimene	0,037	0,005

Ler	MeSA	8,118	0,001	Ler	β -sesquiphellandrene	3,207	0,014
Ler	β -sesquiphellandrene	4,187	0,008	Ler	ethylsalicylate	183,936	0,042
Ler	DMNT	9,416	0,015	Ler	(Z)- β -ocimene	24,405	0,050
Ler	β -farnesene	1,505	0,024	Ler	decanal	1.021,292	0,052
Ler	(E)-2-hexenal	0,003	0,026	Ler	pulegone	178,818	0,053
Ler	α -copaene	0,011	0,029	Ler	3-carene	974,396	0,076
Ler	(Z)- β -ocimene	27,146	0,032	Ler	geraniol	352,990	0,077
Ler	acetophenone	178,215	0,047	Ler	(Z)-3-hexenal	0,959	0,079
Ler	(E)-3-hexanal	25,167	0,057	Ler	allo ocimene	149,587	0,084
Ler	pulegone	84,486	0,067	Ler	TMTT	6,776	0,089
Ler	allo ocimene	673,440	0,074	Ler	β -farnesene	3,541	0,090
Ler	3-carene	974,396	0,076				
Ler	(E)-hexenyl acetate	22,177	0,090				
WS	MeSA	0,443	2,13E-07	WS	MeSA	0,541	5,86E-07
WS	(Z)- β -ocimene	2,285	1,39E-05	WS	(Z)- β -ocimene	0,033	5,42E-05
WS	TMTT	2,205	3,09E-05	WS	nonanal	0,254	0,009
WS	nonanal	0,520	0,002	WS	DMNT	0,011	0,009
WS	(E)-hexenyl acetate	3,563	0,003	WS	linalool	1,124	0,014
WS	β -farnesene	0,148	0,015	WS	(E)-hexenyl acetate	6,100	0,019
WS	2-pentanal	35,049	0,018	WS	β -farnesene	0,307	0,025
WS	(E)- β -ocimene	4,715	0,029	WS	geranyl acteone	293,039	0,038
WS	β -myrcene	33,865	0,056	WS	(E)- β -ocimene	5,071	0,045
WS	(Z)-3-hexen-1-ol	4,346	0,079	WS	β -myrcene	30,728	0,050
WS	cuparene	1,679	0,097	WS	β -sesquiphellandrene	4,477	0,054
				WS	α -farnesene	408,025	0,060
				WS	pulegone	4.725,005	0,061
				WS	hexan-2-ol	0,491	0,063
				WS	hexanal	4,296	0,063
				WS	(E)-2-hexanal	2.309,289	0,065

				WS	TMTT	41,050	0,076
				WS	ethylsalicylate	4.771,445	0,083

Table S7. Total number of compounds that are significantly different between accessions. In bold are the number of compounds depicted that are significantly different after herbivory between two accessions, these numbers are given in italics for the JA-treated plants.

	An-1	C-24	Col-0	Cvi	Eri-1	Kond	Kyo-1	Ler	WS
An-1		3	5	5	4	3	1	2	1
C-24	<i>0</i>		2	5	1	6	3	3	4
Col-0	<i>1</i>	3		4	2	5	2	1	4
Cvi	4	<i>0</i>	3		2	2	3	1	4
Eri-1	<i>1</i>	1	2	<i>1</i>		0	0	0	0
Kond	2	2	4	4	3		0	1	2
Kyo-1	2	2	<i>0</i>	4	<i>1</i>	2		1	0
Ler	2	<i>0</i>	<i>0</i>	3	3	3	<i>1</i>		0
WS	<i>0</i>	1	3	<i>1</i>	5	<i>1</i>	2	<i>0</i>	