

Figure S1. Correlation matrix (Pearson's R values from -1 to 1) between the phenolic composition and the bioactivity measured by *in vitro* antioxidant and anti-hypertensive activities. ACE inhibition expressed as inverse values for the anti-hypertensive activity.

Table S1. Individual phenolic compounds quantified in the different halophyte species. Results expressed as mean \pm standard deviation (n = 2), $\mu\text{g/g}$ of fresh weight (FW). The hydroxycinnamic acids (HCA) and their derivatives were quantified as a 3-O-caffeoylquinic acid equivalent (CQAE), flavonoids were quantified as quercetin-3-glucoside equivalent (QGE) and hydroxybenzoic acids (HBA) were quantified as a gallic acid equivalent (GAE). * Phenolic compound identified with standard.

Class	Peak	Phenolic Compounds $\mu\text{g/g}$ FW	<i>I. crithmoides</i>	<i>S. fruticosa</i>	<i>S. ramosissima</i>	<i>D. crassifolium</i>	<i>C. maritimum</i>	<i>M. nodiflorum</i>	<i>M. crystallinum</i>
HCA	1	<i>p</i> -Coumaric acid derivative	ND	0.370 \pm 0.003	ND	ND	ND	ND	ND
	5	5-O-Caffeoylquinic acid	ND	0.680 \pm 0.003	4.253 \pm 0.033	ND	ND	ND	ND
	8	Caffeic acid derivative	0.482 \pm 0.004	1.138 \pm 0.028	ND	ND	ND	ND	ND
	11	Caffeic acid-O-glucoside	0.356 \pm 0.001	ND	ND	ND	0.744 \pm 0.028	ND	ND
	12	<i>p</i> -Coumaric acid-O-glucoside	ND	ND	ND	ND	10.495 \pm 0.127	0.482 \pm 0.008	0.247 \pm 0.095
	13	<i>p</i> -Coumaric acid	2.868 \pm 0.264	ND	1.538 \pm 0.003	10.457 \pm 1.670	ND	ND	ND
	15	3-O-Caffeoylquinic acid*	2.494 \pm 0.326	13.661 \pm 0.115	9.543 \pm 0.049	ND	47.754 \pm 0.141	ND	ND
	18	Ferulic acid-glucoside	ND	ND	0.839 \pm 0.001	0.208 \pm 0.057	ND	3.912 \pm 0.014	0.498 \pm 0.062
	19	Ferulic acid derivative	ND	ND	ND	ND	ND	0.408 \pm 0.003	ND
	20	Caffeic acid glucuronide	ND	ND	ND	0.302 \pm 0.006	ND	ND	ND
	21	Caffeic acid derivative	0.240 \pm 0.033	ND	ND	ND	ND	ND	ND
	22	Ferulic acid	ND	ND	1.904 \pm 0.011	ND	ND	ND	ND
	23	<i>p</i> -Coumaric acid derivative	ND	2.528 \pm 0.042	ND	ND	ND	ND	ND
	24	Ferulic acid derivative	ND	ND	ND	ND	ND	ND	0.774 \pm 0.073
	26	Sinapic acid -glucoside	ND	ND	ND	ND	ND	ND	3.354 \pm 0.356
	29	Feruloylglucaric acid	ND	ND	ND	ND	ND	ND	0.509 \pm 0.064
	30	Caffeoylquinic acid derivative	ND	2.720 \pm 0.195	ND	ND	ND	ND	ND
	32	Caffeic acid derivative	ND	0.283 \pm 0.106	ND	ND	ND	ND	ND
	34	<i>p</i> -Coumaric acid derivative	ND	-	0.175 \pm 0.001	ND	4.536 \pm 0.141	ND	ND
	38	<i>p</i> -Coumaroylquinic acid (isomer 1)	ND	11.764 \pm 0.200	3.644 \pm 0.003	ND	78.310 \pm 0.141	ND	0.623 \pm 0.078

Table S1. Cont.

Class	Peak	Phenolic Compounds	<i>I. crithmoides</i>	<i>S. fruticosa</i>	<i>S. ramosissima</i>	<i>D. crassifolium</i>	<i>C. maritimum</i>	<i>M. nodiflorum</i>	<i>M. crystallinum</i>
HCA	39	<i>p</i> -Coumaric acid derivative	ND	ND	ND	ND	ND	0.694 ± 0.003	ND
	40	Ferulic acid derivative	ND	ND	0.496 ± 0.014	ND	ND	ND	ND
	41	<i>p</i> -Coumaroylquinic acid (isomer 2)	ND	ND	1.307 ± 0.304	ND	43.242 ± 0.283	ND	ND
	43	Feruloylquinic acid	1.623 ± 0.285	ND	ND	ND	ND	ND	ND
	44	3- <i>O-p</i> -Coumaroyl-5- <i>O</i> -caffeoylquinic acid	0.132 ± 0.007	3.240 ± 0.082	ND	ND	ND	ND	ND
	47	<i>p</i> -Coumaric acid derivative	ND	ND	ND	0.991 ± 0.010	ND	ND	ND
	54	<i>p</i> -Coumaric acid derivative	ND	ND	ND	ND	ND	ND	14.490 ± 0.020
	56	<i>p</i> -Coumaric acid derivative	ND	ND	ND	ND	ND	ND	17.150 ± 0.184
	61	3,4-Dicaffeoylquinic acid	ND	ND	25.724 ± 0.368	ND	ND	ND	ND
	62	<i>p</i> -Coumaric acid glucoside derivative	ND	ND	ND	ND	ND	ND	0.141 ± 0.010
	66	<i>p</i> -Coumaric acid derivative	1.120 ± 0.032	ND	ND	ND	ND	ND	ND
	67	3,5-Dicaffeoylquinic acid	ND	ND	31.440 ± 0.557	ND	5.560 ± 0.085	ND	ND
	71	<i>p</i> -Coumaric acid derivative	ND	1.985 ± 0.205	ND	ND	ND	ND	ND
	74	4,5-Dicaffeoylquinic acid	ND	ND	18.938 ± 0.945	ND	ND	ND	ND
	75	<i>p</i> -Coumaric acid derivative	ND	1.145 ± 0.055	ND	ND	ND	ND	ND
	77	<i>p</i> -Coumaric acid derivative	ND	0.624 ± 0.065	ND	ND	ND	ND	ND
	78	<i>p</i> -Coumaric acid derivative	ND	ND	ND	ND	1.228 ± 0.057	ND	ND
	80	Caffeoylhydrocaffeoylquinic acid	ND	ND	2.661 ± 0.013	ND	ND	ND	ND
	81	Caffeic acid-glucuronide-glucoside (isomer 2)	0.163 ± 0.008	ND	ND	ND	ND	ND	ND
	84	Caffeoylquinic acid derivative	ND	ND	ND	ND	ND	0.387 ± 0.001	ND
	86	Malonyl-3,4- <i>O</i> -dicaffeoylquinic acid derivative	ND	ND	ND	0.814 ± 0.127	ND	ND	ND
	88	<i>p</i> -Coumaric acid derivative	ND	ND	ND	7.747 ± 0.242	ND	ND	ND
	89	Ferulic acid derivative	ND	ND	ND	8.577 ± 0.129	ND	ND	ND
	90	Ferulic acid derivative	ND	ND	ND	0.801 ± 0.208	ND	ND	ND
	93	<i>p</i> -Coumaric acid derivative	ND	ND	ND	0.829 ± 0.131	ND	ND	ND
	94	3,5-Diferuoylquinic acid	ND	ND	ND	ND	ND	0.167 ± 0.001	ND

Table S1. Cont.

Class	Peak	Phenolic Compounds	<i>I. crithmoides</i>	<i>S. fruticosa</i>	<i>S. ramosissima</i>	<i>D. crassifolium</i>	<i>C. maritimum</i>	<i>M. nodiflorum</i>	<i>M. crystallinum</i>
		Σ HCA	9.478 ± 0.960 ^e	40.138 ± 1.099 ^c	102.462 ± 2.302 ^b	30.726 ± 2.058 ^d	191.869 ± 1.003 ^a	6.050 ± 0.03 ^e	37.786 ± 0.942 ^c
HBA	2	Protocatechuic acid-glucoside	0.179 ± 0.005	ND	ND	ND	1.508 ± 0.048	ND	ND
	3	5-Galloylquinic acid	ND	0.269 ± 0.075	ND	ND	ND	ND	ND
	4	Gallic acid derivative	0.429 ± 0.032	ND	ND	ND	ND	ND	ND
	6	Syringic acid	0.538 ± 0.013	ND	ND	ND	ND	ND	ND
	9	Protocatechuic-acid-arabinoside	ND	1.972 ± 0.229	9.609 ± 0.202	ND	ND	ND	ND
	17	Gallic acid derivative	0.222 ± 0.009	ND	ND	ND	ND	ND	ND
	37	Syringic acid derivative	ND	ND	ND	ND	ND	0.179 ± 0.001	ND
	63	Digalloyl quinic acid rhamnoside	ND	ND	ND	ND	ND	0.253 ± 0.001	ND
	85	Galloylquinic acid derivative	ND	ND	ND	ND	ND	0.163 ± 0.001	ND
	91	Galloylquinic acid derivative	ND	ND	ND	ND	ND	0.379 ± 0.001	ND
		Σ HBA	1.368 ± 0.059 ^c	2.241 ± 0.304 ^b	9.609 ± 0.202 ^a	ND	1.508 ± 0.048 ^c	0.974 ± 0.004 ^c	ND
		Σ Phenolic acids	10.846 ± 0.078 ^f	42.379 ± 0.092 ^c	112.071 ± 0.178 ^b	30.726 ± 0.275 ^c	193.377 ± 0.112 ^a	7.024 ± 0.004 ^g	37.786 ± 0.109 ^d
Flavone	25	Apigenin 6-C-glucoside-7-O-glucoside	ND	ND	ND	ND	8.688 ± 0.113	ND	ND
	42	Chrysin-6-C-ara-8-C-glu	ND	ND	ND	ND	ND	0.643 ± 0.007	ND
	46	Acacetin 3,6-di-C-glucoside	ND	ND	ND	ND	ND	ND	0.449 ± 0.038
	50	Chrysin-6-C-glucosyl-8-C-arabinoside	ND	ND	ND	ND	ND	ND	7.357 ± 0.480
	60	Apigenin 6-C-glucoside	ND	ND	ND	ND	1.288 ± 0.057	ND	ND
	64	Luteolin derivative	ND	ND	ND	1.684 ± 0.105	ND	ND	ND
	68	Diosmetin 7-O-rutinoside	ND	ND	ND	ND	0.887 ± 0.004	ND	ND
	76	Vitexin (apigenin-8-C-glu)	0.319 ± 0.010	ND	ND	ND	ND	ND	ND
	87	Vitexin derivative	ND	ND	ND	ND	ND	0.160 ± 0.003	ND
		Σ Flavone	0.319 ± 0.010 ^e	ND	ND	1.684 ± 0.105 ^c	10.863 ± 0.174 ^a	0.803 ± 0.010 ^d	7.806 ± 0.518 ^b

Table S1. Cont.

Class	Peak	Phenolic Compounds	<i>I. crithmoides</i>	<i>S. fruticosa</i>	<i>S. ramosissima</i>	<i>D. crassifolium</i>	<i>C. maritimum</i>	<i>M. nodiflorum</i>	<i>M. crystallinum</i>
Flavanol	10	Gallocatechin	0.167 ± 0.024	0.239 ± 0.004	0.299 ± 0.001	ND	ND	3.026 ± 0.003	0.182 ± 0.023
	14	Epigallocatechin	ND	ND	ND	0.862 ± 0.107	ND	0.510 ± 0.003	0.519 ± 0.047
	51	Epicatechin hydrate	ND	ND	ND	ND	ND	1.064 ± 0.011	ND
	52	Gallocatechin derivative	0.549 ± 0.035	ND	ND	ND	ND	ND	ND
	55	Epicatechin derivative	ND	11.534 ± 0.164	ND	ND	ND	ND	ND
	Σ Flavanol		0.716 ± 0.059 ^c	11.773 ± 0.168 ^a	0.299 ± 0.001 ^d	0.862 ± 0.107 ^c	ND	4.600 ± 0.017 ^b	0.701 ± 0.007 ^c
Flavanone	16	Pinobanksin-3-O-pentanoate	ND	ND	ND	ND	0.731 ± 0.048	ND	1.571 ± 0.273
	31	Pinobanksin-5-methyl ether-3-O-acetate (isomer)	6.382 ± 0.176	ND	ND	ND	ND	ND	ND
	33	Pinocembrin derivative	ND	ND	ND	ND	ND	0.338 ± 0.057	ND
	48	Eriodictyol-O-hexoside	ND	7.071 ± 0.114	ND	ND	ND	2.606 ± 0.011	ND
	49	Erydictiol	ND	ND	ND	ND	ND	8.568 ± 0.090	ND
	92	Pinobanksin-5-methyl ether-3-O-acetate (isomer)	ND	ND	ND	ND	ND	1.553 ± 0.004	ND
	Σ Flavanone		6.382 ± 0.176 ^c	7.071 ± 0.114 ^b	ND	ND	0.731 ± 0.048 ^e	13.065 ± 0.162 ^a	1.571 ± 0.273 ^d
Flavonol	7	Rhamnetin	0.194 ± 0.082	ND	ND	ND	ND	ND	ND
	27	Avicularin	ND	ND	ND	ND	ND	2.422 ± 0.003	ND
	35	Kaempferol derivative	ND	ND	ND	ND	ND	ND	0.576 ± 0.014
	45	Quercetin-3-O-rutinoside	ND	ND	ND	ND	11.854 ± 0.141	ND	ND
	53	Quercetin 3-O-glucoside*	ND	ND	3.338 ± 0.009	ND	5.932 ± 0.042	ND	ND
	58	Isorhamnetin-rutinoside derivative	ND	ND	ND	13.825 ± 0.100	ND	ND	ND
	59	Rhamnetin hexosyl pentoside	ND	27.388 ± 0.127	ND	ND	ND	ND	ND
	65	Isorhamnetin 3-O-robinobioside	ND	20.871 ± 0.143	ND	ND	ND	ND	ND
	69	Quercetin dipentoside	ND	ND	ND	ND	ND	0.133 ± 0.001	ND
	70	Quercetin-3-O-glucoside derivative	ND	ND	ND	ND	ND	ND	0.687 ± 0.066
	72	Quercetin derivative	ND	ND	ND	ND	ND	0.316 ± 0.006	ND
	73	Isorhamnetin-glucoside derivative	ND	ND	ND	7.281 ± 0.103	ND	ND	ND

Table S1. <i>Cont.</i>									
Class	Peak	Phenolic Compounds	<i>I. crithmoides</i>	<i>S. fruticosa</i>	<i>S. ramosissima</i>	<i>D. crassifolium</i>	<i>C. maritimum</i>	<i>M. nodiflorum</i>	<i>M. crystallinum</i>
Flavonols	82	Isorhamnetin-3-O-glucoside (dimer)	ND	ND	2.828 ± 0.450	ND	ND	ND	ND
	83	Kaempferol	ND	ND	ND	ND	ND	ND	0.594 ± 0.062
		3-(2''-[glucosyl-(1->3)-rhamnosyl]-6''-rhamnosyl galactoside)							
		Σ Flavonol	0.194 ± 0.082 ^s	48.259 ± 0.270 ^a	6.166 ± 0.459 ^d	21.106 ± 0.203 ^b	17.786 ± 0.183 ^c	2.871 ± 0.010 ^e	1.857 ± 0.142 ^f
Flavanonol	36	Dihydroquercetin	ND	6.404 ± 0.792	ND	ND	ND	ND	ND
		Σ Flavanonol	ND	6.404 ± 0.792	ND	ND	ND	ND	ND
		Σ Flavonoids	7.611 ± 0.065 ^f	73.507 ± 0.224 ^a	6.465 ± 0.153 ^s	23.652 ± 0.103 ^c	29.38 ± 0.068 ^b	21.339 ± 0.015 ^d	11.935 ± 0.125 ^e
Coumarin	28	Coumarin glycoside ester	ND	ND	ND	1.458 ± 0.215	ND	ND	ND
Monocarboxylic acid	57	Piscidic acid derivative	1.828 ± 0.219	ND	ND	ND	ND	ND	ND
Gallotannin	79	Hexahydroxydiphenoyl-Glucose	ND	ND	ND	ND	ND	0.177 ± 0.001	ND
		Σ Phenolic compounds	18.457 ± 0.074 ^f	115.886 ± 0.137 ^b	118.536 ± 0.174 ^b	55.836 ± 0.225 ^c	222.757 ± 0.094 ^a	28.363 ± 0.010 ^e	49.721 ± 0.114 ^d

The letters (a-g) indicate significant differences between the halophytes using the Tukey’s test (p < 0.05). ND: Not Detected; LOQ = 0.01 µg compound/g FW.

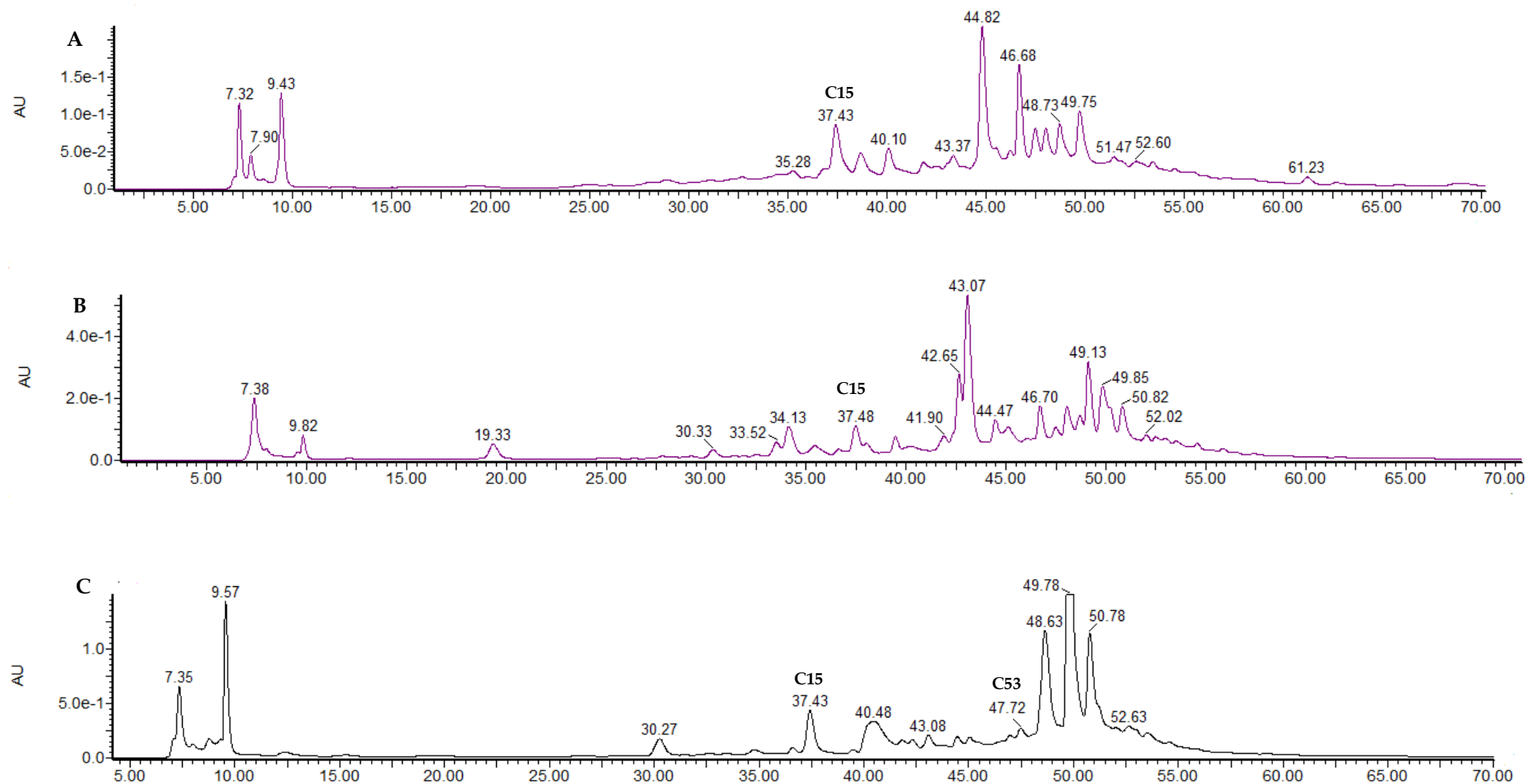


Figure S2. Chromatographic profile at 280 nm by HPLC-DAD-ESI-MS/MS of the halophytes species. *I. crithmoides* (A); *S. fruticosa* (B); *S. ramosissima* (C); *D. crassifolium*(D); *C. maritimum* (E); *M. nodiflorum* (F) and *M. crystallinum* (G).

Figure S2. Cont.

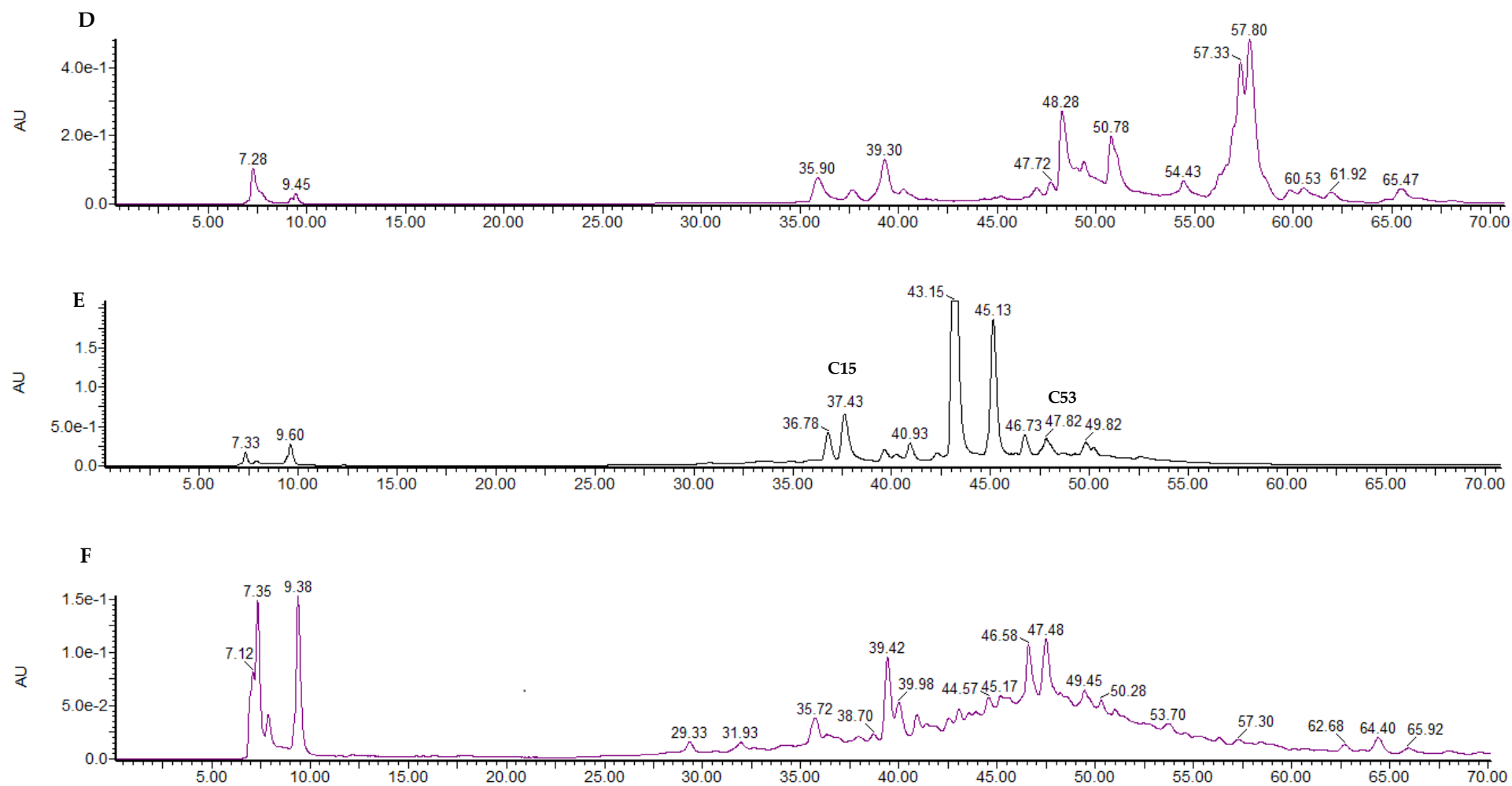


Figure S2. Chromatographic profile at 280 nm by HPLC-DAD-ESI-MS/MS of the halophytes species. *I. crithmoides* (A); *S. fruticosa* (B); *S. ramosissima* (C); *D. crassifolium* (D); *C. maritimum* (E); *M. nodiflorum* (F) and *M. crystallinum* (G).

Figure S2. Cont.

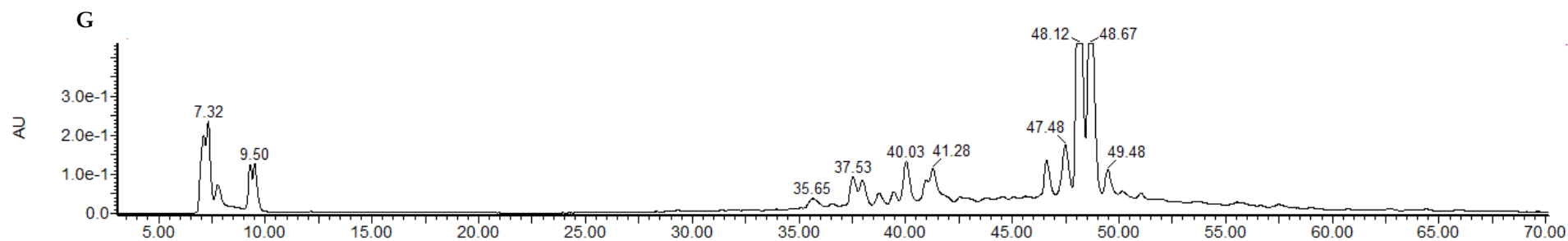


Figure S2. Chromatographic profile at 280 nm by HPLC-DAD-ESI-MS/MS of the halophytes species. *I. crithmoides* (A); *S. fruticosa* (B); *S. ramosissima* (C); *D. crassifolium*(D); *C. maritimum* (E); *M. nodiflorum* (F) and *M. crystallinum* (G).

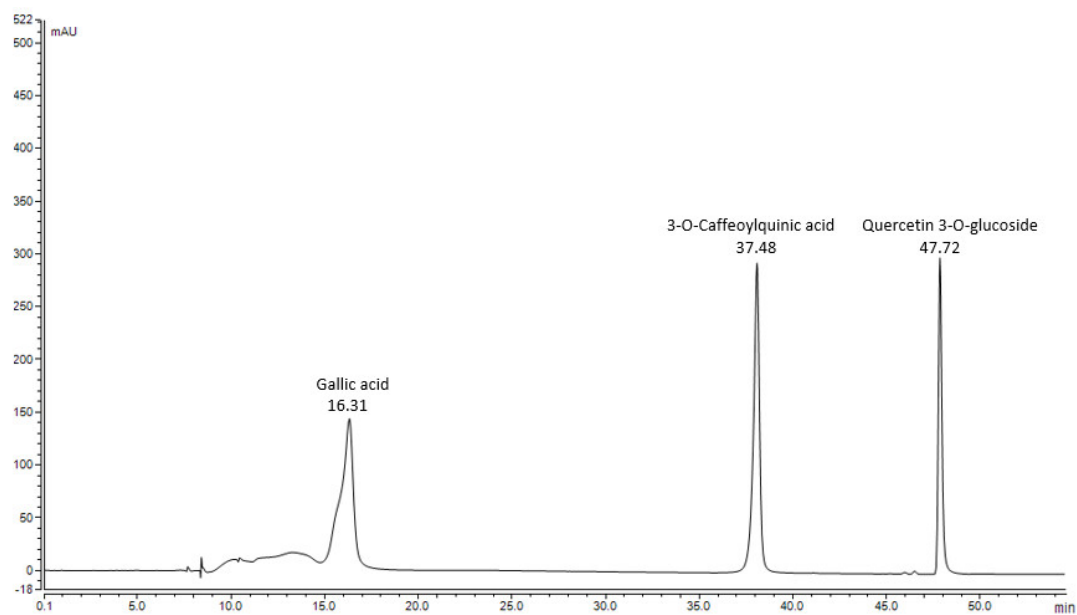


Figure S3. Chromatographic profile at 280 nm by HPLC-DAD of the standards of phenolic compounds: gallic acid, chlorogenic acid (3-O-caffeoylquinic acid) and quercetin-3-glucoside.

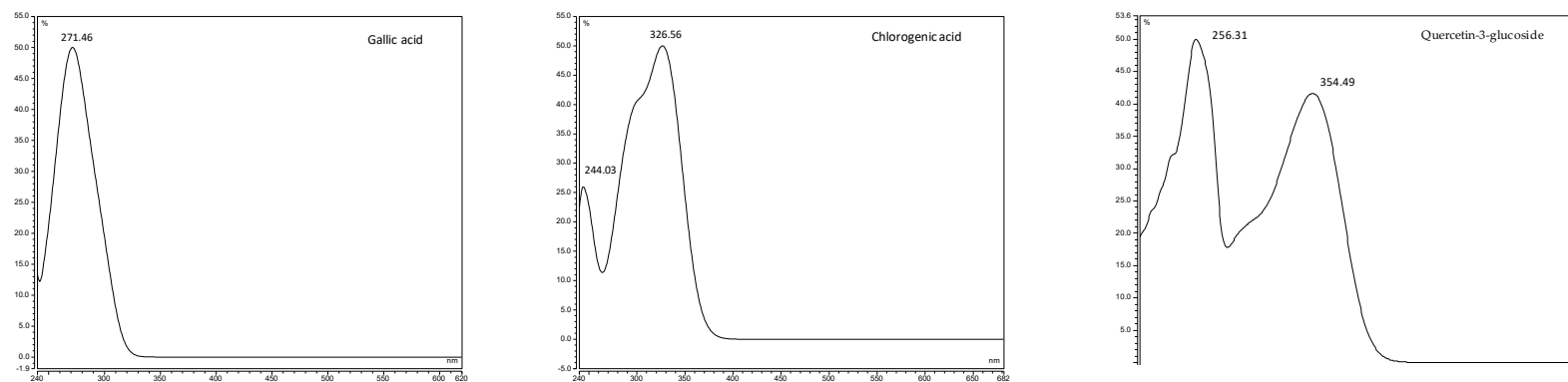


Figure S4. UV-Visible spectra of the phenolic compounds standards: gallic acid, chlorogenic acid (3-O-caffeoylquinic acid) and quercetin-3-glucoside.

Table S2. Characterization parameters (retention time, RT, % of total peak area, tentative identification, chemical classes, odor, calculated/theoretical linear retention index, LRI) of volatile compounds in halophyte species using SPME-GC–MS. The references used to support the tentative identifications were indicated.

Species	Peak	RT (min) ^a	Area%	Tentative Identification	Chemical Classes	Odor Description ^b	Calculated LRI ^c	Theoretical LRI	References
<i>C. maritimum</i>	1	7.24	0.91	3-hexen-1-ol	Alcohol	green, marine, seaweed	860	861	[153-155]
	2	9.55	0.71	alpha-thujene	Terpene	herbal, green, weak earthy	927	934	[156-158]
	3	9.742	1.71	alpha-pinene	Terpene	oily, green	932	933	[159,160]
	4	11.283	17.24	alpha-fenchene	Terpene	fruity, fresh	973	962	[161,162]
	5	11.992	1.01	beta-myrcene	Terpene	herbaceous, sweet	992	992	[158-163]
	6	12.817	0.40	alpha-terpinene	Terpene	resinous	1016	1017	[158,164]
	7	13.125	13.41	<i>p</i> -cymene	Terpene	green, fruity, aromatic	1025	1031	[159,165-167]
	8	13.225	0.40	beta-phellandrene	Terpene	NF	1028	1032	[168,169]
	9	13.65	8.87	(<i>Z</i>)-beta-ocimene	Terpene	floral, herbaceous	1041	1037	[168,169]
	10	13.992	0.30	(<i>E</i>)-beta-ocimene	Terpene	floral, herbaceous	1051	1047	[168,169]
	11	14.358	30.44	<i>m</i> -mentha-4,8-diene	Terpene	NF	1062	1081	[168]
	12	15.25	0.30	alpha-terpinolene	Terpene	woody, herbaceous	1088	1086	[164,170]
	13	16.617	0.20	allo-ocimene	Terpene	fresh	1131	1131	[171,172]
	14	18.075	0.20	4-terpineol	Terpene	roasted, woody	1178	1178	[173,174]
	15	19.65	23.59	thymol methyl ether	Terpene	oregano-like, thyme-like, rose- mary-like, medicinal	1239	1237	[175-176]
	16	25.242	0.20	alpha-bergamotene	Terpene	wood, warm, tea	1439	1438	[174,178]
	17	27.092	0.10	beta-bisabolene	Terpene	NF	1513	1512	[179]

Table S2. *Cont.*

Species	Peak	RT (min) ^a	Area%	Tentative Identification	Chemical Classes	Odor Description ^b	Calculated LRI ^c	Theoretical LRI	References
<i>I. crithmoides</i>	1	7.00	0.42	3-hexen-1-ol	Alcohol	green, marine, seaweed	860	861	[153-155]
	2	7.53	0.73	1-hexanol	Alcohol	woody, sweet, green, fruity	873	872	[180-182]
	3	9.74	0.52	alpha-pinene	Terpene	oily, green	932	933	[159,160]
	4	10.26	8.34	camphene	Terpene	sweet	946	947	[159,160,183]
	5	11.31	0.73	beta-pinene	Terpene	woody, green, pine-like	974	979	[158,14,183]
	6	12.01	0.21	beta-myrcene	Terpene	herbaceous, sweet	993	992	[158-162]
	7	12.40	17.41	alpha-phellandrene	Terpene	fresh, green	1004	1005	[164,181,183]
	8	12.63	0.10	3-hexen-1-ol acetate	Ester	fruity, floral	1010	1007	[180,184]
	9	12.87	0.31	hexyl acetate	Ester	fruit, herb	1018	1019	[185,186]
	10	13.15	22.94	<i>p</i> -cymene	Terpene	green, fruity, aromatic	1026	1031	[159,165-167]
	11	13.24	2.29	limonene	Terpene	pine/chemical, floral/fresh	1029	1029	[187,188]
	12	14.28	0.21	gamma-terpinene	Terpene	green, woody	1059	1059	[159,170,189]
	13	15.25	1.15	alpha-terpinolene	Terpene	woody, herbaceous	1088	1086	[164,170]
	14	17.71	0.31	borneol	Terpene	camphoraceous, earthy	1167	1165	[190,191]
	15	18.08	0.21	4-terpineol	Terpene	roasted, woody	1179	1178	[173,174]
	16	19.78	3.96	thymol methyl ether	Terpene	oregano-like, thyme-like, rose- mary-like, medicinal	1236	1237	[175-177]
	17	20.10	0.73	allyl hexanoate	Ester	sweet, fruity	1248	-	[192]
	18	20.25	17.52	2,4,6-trimethylacetophenone	Ketone	not found	1253	1252	[193]
	19	21.01	0.63	cyclooctyl acetate	Ester	not found	1279	1279	[194]
	20	21.26	19.71	isobornyl acetate	Ester	herb, woody, sweet, minty	1288	1285	[191,192]
	21	21.56	0.21	thymol	Terpene	thyme-like, spicy	1299	1293	[157,195]
	22	23.80	1.15	cadinene	Terpene	herb, burnt, sulphur	1382	1362	[196,197]
	23	24.92	0.21	2-tert-Butyl-1,4-dimethoxybenzene	Terpene	not found	1426	1427	[198]

Table S2. *Cont.*

Species	Peak	RT (min) ^a	Area%	Tentative Identification	Chemical Classes	Odor Description ^b	Calculated LRI ^c	Theoretical LRI	References
<i>D. crassifolium</i>	1	5.23	6.35	hexanal	Aldehyde	herbal, grassy, green	807	801	[154,180,184,199]
	2	7.42	6.69	3-hexen-1-ol	Alcohol	green, marine, seaweed	866	861	[153-155]
	3	7.88	10.37	1-hexanol	Alcohol	woody, sweet, green, fruity	878	872	[180-182]
	4	8.19	1.34	<i>n</i> -butyl ether	Ether	not found	887	860	[200]
	5	9.27	6.35	oxime-methoxy-phenyl-	Ester	fishy	916	910	[201,202]
	6	11.97	1.67	2,2,4,6,6-pentamethylheptane	Terpene	stony, dusty	990	991	[203,204]
	7	13.24	9.03	<i>p</i> -cymene	Terpene	green, fruity, aromatic	1027	1031	[159,165-167]
	8	15.78	1.00	undecane	Hydrocarbon	not found	1103	1100	[201]
	9	18.83	3.34	dodecane	Hydrocarbon	alkane-like, chemical	1202	1200	[201,205]
	10	19.81	20.74	thymol methyl ether	Terpene	oregano-like, thyme-like, rosemary-like, medicinal	1236	1237	[175-177]
	11	20.23	9.70	2,4,6-trimethylacetophenone	Ketone	not found	1251	1252	[193]
	12	21.52	19.73	isobornyl acetate	Ester	herb, woody, sweet, minty	1287	1285	[191,192]
	13	23.80	3.68	alpha-gurjunene	Terpene	green, herbal	1382	1407	[190,206]

Table S2. *Cont.*

Species	Peak	RT (min) ^a	Area%	Tentative Identification	Chemical Classes	Odor Description ^b	Calculated LRI ^c	Theoretical LRI	References
<i>M. nodiflorum</i>	1	4.342	1.258149	1-pentanol	Alcohol	balsamic	<800	766	[207,208]
	2	5.108	1.372527	octane	Hydrocarbon	Alkane-like	803	800	[209]
	3	7.108	8.578291	3-hexen-1-ol	Alcohol	green, marine, seaweed	858	861	[153-155]
	4	7.683	77.77651	1-hexanol	Alcohol	woody, sweet, green, fruity	873	872	[180-182]
	5	9.542	0.228754	alpha-thujene	Terpene	herbal, green, weak earthy	924	924	[190,210]
	6	11.342	0.343132	1-heptanol	Alcohol	Metallic	973	967	[211]
	7	11.925	2.85943	2-ethylbutanal	Aldehyde	green	989	-	[210]
	8	12.667	0.228754	3-hexen-1-ol acetate	Ester	fruity, floral	1010	1007	[180,184]
	9	12.883	0.800641	hexyl acetate	Ester	fruit, herb	1016	1019	[185,186]
	10	12.983	1.830035	trans-2-hexenyl acetate	Ester	fruity, sweet, herbaceous	1019	1018	[213,214]
	11	13.258	0.343132	limonene	Terpene	pine, chemical, floral, fresh	1028	1029	[188,215]
	12	14.908	0.125815	1-octanol	Alcohol	citrus-like, herbal	1077	1070	[216,217]
	13	15.742	0.137253	linalool	Terpene	pleasant scent, floral	1102	1097	[166,180]
	14	15.9	0.343132	2,6-dimethylcyclohexanol	Alcohol	earthy	1107	1108	[218,219]
	15	18.025	0.457509	1-nonanol	Alcohol	fatty	1176	1176	[220,221]
	16	19.358	0.343132	beta-cyclocitral	Terpene	sweet-tobacco, grape	1221	1220	[215,222]
	17	19.8	1.258149	thymol methyl ether	Terpene	oregano-like, thyme-like, rosemary-like, medicinal	1236	1237	[175-177]
	18	20.225	0.571886	2,4,6-trimethylacetophenone	Ketone	not found	1251	1252	[193]
	19	21.233	0.915018	isobornyl acetate	Ester	herb, woody, sweet, minty	1286	1285	[190,191]
	20	23.8	0.228754	cadinene	Terpene	herb, burnt, sulphur	1382	1362	[196,197]

Table S2. Cont.

Species	Peak	RT (min) ^a	Area%	Tentative Identification	Chemical Classes	Odor Description ^b	Calculated LRI ^c	Theoretical LRI	References
<i>S. fruticosa</i>	1	7.41	27.56	2-hexenal	Aldehyde	floral, herbal	856	861	[153,180,184]
	2	7.76	44.04	1-hexanol	Alcohol	woody, sweet, green, fruity	875	872	[180-182]
	3	10.28	1.14	camphene	Terpene	sweet	944	947	[159,160,183]
	4	11.07	4.15	2-hexenoic acid, methyl ester	Ester	not found	966	972	[223]
	5	11.81	0.31	<i>n</i> -caproic acid vinyl ester	Ester	not found	986	964	[205]
	6	12.52	3.83	alpha-phellandrene	Terpene	fresh, green	1006	1005	[164,181,183]
	7	12.81	0.21	4-carene	Terpene	sweet, pungent	1014	-	[166]
	8	12.92	0.10	<i>trans</i> -2-hexenyl acetate	Ester	fruity, sweet, herbaceous	1017	1018	[213,214]
	9	13.05	2.80	<i>p</i> -cymene	Terpene	green, fruity, aromatic	1022	1031	[159,165-167]
	10	13.19	1.66	limonene	Terpene	pine, chemical, floral, fresh	1026	1029	[185,188]
	11	13.81	0.62	ethyl 2-hexenoate	Ester	citrus-like, fatty	1045	1053	[224,225]
	12	15.20	0.41	alpha-terpinolene	Terpene	woody, herbaceous	1085	1086	[164,170]
	13	15.46	0.52	1-nonen-4-ol	Alcohol	cardboard	1093	1094	[226]
	14	20.14	4.46	<i>p</i> -(1-butenyl)anisole	Hydrocarbon	not found	1248	1259	[227]
	15	20.92	0.41	3-octen-1-ol, acetate	Ester	toasted nut, smoky, dusty	1275	1279	[184]
	16	21.18	6.63	isobornyl acetate	Ester	herb, woody, sweet, minty	1284	1285	[190,191]
	17	23.71	1.14	cadinene	Terpene	herb, burnt, sulphur	1378	1362	[196,197]

Table S2. *Cont.*

Species	Peak	RT (min) ^a	Area%	Tentative Identification	Chemical Classes	Odor Description ^b	Calculated LRI ^c	Theoretical LRI	References
<i>M. crystallinum</i>	1	4.28	1.78	1-pentanol	Alcohol	balsamic	766	766	[207,208]
	2	5.07	9.89	hexanal	Aldehyde	herbal, grassy, green	802	801	[154,180,184,199,208]
	3	6.93	36.66	2-hexenal	Aldehyde	floral, herbal	853	861	[153,28,32]
	4	7.63	40.33	1-hexanol	Alcohol	woody, sweet, green, fruity	872	872	[180-182]
	5	8.76	0.11	heptanal	Aldehyde	penetrating oily, harsh	902	902	[182,199,228]
	6	9.57	0.67	methoxy-phenyl-oxime	Ester	fishy	924	916	[201,202]
	7	11.55	0.11	3,5,5-trimethyl-2-hexene	Hydrocarbon	not found	979	977	[201]
	8	11.78	0.22	1-octen-3-ol	Alcohol	mushroom	985	980	[154,158,188]
	9	12.30	0.11	3-octanol	Alcohol	mushroom-like, herbal, nutty	994	998	[157,229]
	10	12.69	0.11	<i>cis</i> -3-hexenyl acetate	Ester	fruity, green, floral	1011	1007	[180,184,230]
	11	12.92	0.56	hexyl acetate	Ester	fruit, herb	1017	1019	[185,186]
	12	13.02	0.22	<i>trans</i> -2-hexenyl acetate	Ester	fruity, sweet, herbaceous	1020	1018	[213,214]
	13	13.13	0.22	<i>p</i> -cymene	Terpene	green, fruity, aromatic	1024	1031	[159,165-167]
	14	14.33	0.33	3-carene	Terpene	sweet, strong	1060	-	[231]
	15	14.98	0.03	1-octanol	Alcohol	citrus-like, herbal	1079	1070	[216,217]
	16	15.89	1.22	nonanal	Aldehyde	green, citrus-like	1107	1111	[153,232]
	17	18.31	0.22	octanoic acid	Carboxylic acid	fruity, faint, candy	1185	1173	[233,234]
	18	18.81	0.33	dodecane	Hydrocarbon	alkane-like, chemical	1201	1200	[201,205]
	19	19.80	2.78	thymol methyl ether	Terpene	oregano-like, thyme-like, rose- mary-like, medicinal	1236	1237	[175-177]
	20	20.22	1.22	2,4,6-trimethylacetophenone	Ketone	not found	1251	1252	[193]
	21	21.05	0.44	nonanoic acid	Carboxylic acid	fatty, cheesy	1280	1275	[157]
	22	21.23	2.33	isobornyl acetate	Ester	herb, woody, sweet, minty	1286	1285	[190,191]
	23	22.99	0.09	alpha-cubebene	Terpene	herbaceous, waxy	1351	1352	[235,236]

Table S2. Cont.

Species	Peak	RT (min) ^a	Area%	Tentative Identification	Chemical Classes	Odor Description ^b	Calculated LRI ^c	Theoretical LRI	References
<i>S. ramosissima</i>	1	4.99	6.30	hexanal	Aldehyde	herbal, grassy, green	800	801	[154,180,184,199,208]
	2	7.04	45.65	2-hexenal	Aldehyde	floral, herbal	856	861	[153,180,184]
	3	7.78	47.68	octanal	Aldehyde	green, citrusy, lemon	976	1001	[208,237,238]
	4	11.58	0.06	<i>n</i> -caproic acid vinyl ester	Ester	not found	979	964	[205]
	5	13.81	0.10	benzeneacetaldehyde	Aldehyde	fresh, floral	1044	1043	[232,239]
	6	15.57	0.10	1-nonen-4-ol	Alcohol	cardboard	1096	1094	[226,240]
	7	15.88	0.10	2,6-dimethyl-cyclohexanol	Alcohol	earthy	1106	1108	[218,219]

^aMean of the retention times in the chromatograms of the fresh halophyte species (n=2). ^bDescribed odor descriptions. ^cMean calculated of the linear retention indexes. NF – Not found.

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