

Supplementary materials

Mining the Metabolome and the Agricultural and Pharmaceutical Potential of Sea Foam-Derived Fungi

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Table S1. Identification of six foam-derived fungal strains isolated from Windebyer Noor (Nov, 2015) according to the two new BLAST searches.

Query		BLAST ALL			BLAST TYPE			
Isolate name	Acc. No	Name	Acc. No.	Similarity (%)	Name	Acc. No.	Similarity (%)	Taxonomic assignment
B2F1B	MH791266.1 <i>Plectosphaerella</i> sp.	Fungal sp. strain S255T	KU839553.1	99.79	<i>Plectosphaerella niemeijerorum</i> CBS 143233 T	NR_156677.1	99.59	<i>Plectosphaerella</i> sp.
						MG386080.1	99.59	
		Fungal sp. strain S255S	KU839552.1	99.79	<i>Plectosphaerella oligotrophica</i> CGMCC 3.15078 T	NR_155632.1	98.35	
		Fungal sp. strain S255O	KU839551.1	99.79	<i>Plectosphaerella oligotrophica</i> strain LC1990	JX508810.1	98.35	
		Fungal sp. strain S255N	KU839550.1	99.79	<i>Plectosphaerella delsorboi</i> CBS 116708T	NR_155694.1	96.30	
						KY499237.1	96.30	
		Fungal sp. strain S255L	KU839549.1	99.79	<i>Plectosphaerella populi</i> CBS 139623	NR_138002.1	95.08	
25.88F1C	MH791280.1 <i>Emericellopsis</i> sp.	<i>Emericellopsis</i> sp. s012	HQ649988.1	95.30	<i>Emericellopsis terricola</i> CBS 120.40 T	NR_159843.1	94.97	<i>Emericellopsis</i> sp.
						MH856058.1	94.97	
		<i>Emericellopsis terricola</i> CBS 120.40 T	NR_159843.1	94.97	<i>Acremonium tubakii</i> CBS 790.69	NR_163810.1	94.63	
						MH859429.1	94.63	
		<i>Emericellopsis terricola</i> strain CBS 126950	MH864349.1	94.97	<i>Emericellopsis microspora</i> strain CBS 380.62	MH858186.1	94.63	
		<i>Emericellopsis terricola</i> strain CBS 387.70A	MH859748.1	94.97	<i>Emericellopsis glabra</i> strain CBS 119.40	MH856057.1	93.55	
		<i>Emericellopsis terricola</i> strain CBS 246.70	MH859590.1	94.97	<i>Emericellopsis pusilla</i> CBS 226.62 T	NR_159878.1	94.30	

86F1C	MH791259.1 <i>Cladosporium cladosporioides</i>	<i>Cladosporium asperulatum</i> isolate LH4	MN220523.1	100	<i>Cladosporium verrucocladosporioides</i> strain CBS 126363	MH863939.1	100	<i>Cladosporium</i> sp.
		<i>Cladosporium</i> sp. strain SMY02	MK355726.1	100	<i>Cladosporium antarcticum</i> CBS 690.92 T	NR_121332.1	99.79	
						EF679334.2	99.79	
		<i>Cladosporium</i> sp. isolate 18MWF13.1	MN128233.1	100	<i>Cladosporium ossifragi</i> culture CBS:842.91	EF679381.2	99.79	
		<i>Cladosporium cladosporioides</i> isolate C11	MK813962.1	100	<i>Cladosporium iridis</i> strain CBS 138.40	EU167591.1	99.79	
<i>Cladosporium cladosporioides</i> isolate C10	MK813961.1	100	<i>Cladosporium subinflatum</i> strain CBS 121630	MH863129.1	99.79			
31.68F1B	MH791236.1 <i>Penicillium bialowiezense</i>	<i>Penicillium</i> sp. isolate P11-16A	MN186768.1	100	<i>Penicillium bialowiezense</i> strain CBS 227.28 T	NR_111323.1	100	<i>Penicillium bialowiezense</i>
						MH854996.1		
		<i>Penicillium</i> sp. isolate P8-16A	MN186762.1	100	<i>Penicillium brunneostoloniferum</i> strain CBS 317.59	KC411734.1	98.77	
		<i>Penicillium</i> sp. isolate SL71_37c_D	MN105533.1	100	<i>Penicillium brevicompactum</i> NRRL 2011T	NR_121299.1	98.77	
						AY484912.1		
<i>Penicillium</i> sp. isolate SL16_77a_G2	MN105477.1	100	<i>Penicillium brevicompactum</i> strain CBS 257.29	KF465776.1	98.77			
<i>Penicillium</i> sp. isolate SL63_46_G	MN105342.1	100	<i>Penicillium patris-mei</i> CBS 210.28 T	NR_138315.1	98.46			
36.97F1C	MH791183.1 <i>Penicillium</i> sp.	<i>Penicillium</i> sp. FA6-2	KF776918.1	98.04	<i>Penicillium limosum</i> CBS 339.97 T	NG_062729.1	97.77	<i>Penicillium</i> sp.
						EF411061.1		

	18S rRNA gene	<i>Penicillium</i> sp. FA9	KF776917.1	98.04	<i>Penicillium commune</i> CBS 343.51 T	NG_062634.1	97.77	
						AF548088.1		
		<i>Penicillium brevicompactum</i> isolate PenC	FJ717699.1	98.04	<i>Talaromyces leycettanus</i> CBS 398.68T	NG_062664.1	97.68	
		<i>Penicillium</i> sp. strain OUCMDZ-4920	KY448996.1	98.04	<i>Talaromyces wortmannii</i> CBS 391.48T	NG_062801.1	97.59	
				GU733355.1				
		<i>Penicillium brevicompactum</i> strain KR5	MH047201.1	97.86	<i>Penicillium roqueforti</i> strain ATCC 10110	GQ458035.1	97.59	
62.72F1A	MH791282.1 <i>Penicillium brevicompactum</i>	<i>Penicillium</i> sp. isolate P11-16A	MN186768.1	100	<i>Penicillium bialowiezense</i> strain CBS 227.28 T	MH854996.1	100	<i>Penicillium bialowiezense</i>
						NR_111323.1		
		<i>Penicillium</i> sp. isolate P8-16A	MN186762.1	100	<i>Penicillium brunneostoloniferum</i> strain CBS 317.59	KC411734.1	98.62	
		<i>Penicillium</i> sp. isolate SL71_37c_D	MN105533.1	100	<i>Penicillium patris-mei</i> CBS 210.28 T	NR_138315.1	98.62	
		<i>Penicillium</i> sp. isolate SL16_77a_G2	MN105477.1	100	<i>Penicillium neocrassum</i> NRRL 35639 T	NR_121321.1	98.62	
		<i>Penicillium</i> sp. isolate SL63_46_G	MN105342.1	100	<i>Penicillium brevicompactum</i> NRRL 2011	NR_121299.1	98.62	

Table S2. Putative identification of metabolites produced by foam-derived fungi. Annotations were made based on m/z $[M+H]^+$ or other adducts (specified), biological source, retention time (t_R), predicted molecular formula and fragmentation pattern. Confidence level of putative identification after Sumner et al. [1].

Compound I.D	t_R	Parent mass $[M+H]^+$	Putative ID	Chemical class/ family	Molecular formula m/z $[M+H]^+$	Δ ppm	MS ²	Confidence level	Isolate	Reference
1	0.6	293.1001	italicic acid	polyketide	C ₁₅ H ₁₇ O ₆	5.0	281.0570; 267.0149; 253.0010; 237.0129; 202.0589; 187.9807; 163.0728; 145.0589; 127.0410; 118.0874; 99.0874; 81.0292	3	62.72F1A	[2]
2	3.7	235.0962	citreopyrone E	pyranone	C ₁₃ H ₁₅ O ₄	-3.4	217.0848; 191.0696; 179.1068; 151.0409; 135.0788; 126.1260; 112.9818; 86.0965; 67.0180	3	31.68F1B	[3]
3	3.9	271.1441	quinolactacin A1	alkaloid	C ₁₆ H ₁₉ N ₂ O ₂	-1.1	214.0740; 233.0798; 203.1697; 189.9731; 185.4548; 171.0642; 161.0518; 146.8997; 129.0546; 123.1155; 112.9648; 97.0940; 86.0966; 84.9576	3	31.68F1B	[4,5]
4	5.9	445.2588	preaustinoid A	meroterpenoid	C ₂₆ H ₃₇ O ₆	-0.7	385.2373; 413.2239; 367.2273; 243.1746	2	31.68F1B, 36.97F1C, 62.72F1A	[6,7]
5	5.2	461.2432	preaustinoid A1	meroterpenoid	C ₂₆ H ₃₇ O ₇	-2.0	443.2457; 411.2151; 383.2230; 365.2193; 213.1639	3	31.68F1B, 62.72F1A	[7]

6	6.0	207.0687	7-Hydroxy-2-(hydroxymethyl-5-methyl-4H-1-benzopyran-4-one	benzopyranone	C ₁₁ H ₁₁ O ₄	0	189.0543; 177.0549; 151.0384; 131.0485; 117.0320; 53.0011; 77.0384; 91.0539	3	31.68F1B	[8]
7	6.0	321.1381	mycophenolic acid	meroterpenoid	C ₁₇ H ₂₁ O ₆	2.5	303.1240; 275.1284; 285.1128; 207.0660; 195.0655; 159.0447; 177.0548; 159.0447; 109.0644; 81.0686; 77.0412; 69.9658	2	31.68F1B, 62.72F1A, 36.97F1C	[9,10]
8	4.7	359.1101 [M+Na]	4-hydroxy-mycophenolic acid	meroterpenoid	C ₁₇ H ₂₀ O ₇ Na	1.7	-	2	36.97F1C,	[9,10]
9	7.5	335.1489	mycophenolic acid methyl ester	meroterpenoid	C ₁₈ H ₂₃ O ₆	-1.8	303.1239; 275.1292; 207.0660	2	31.68F1B, 36.97F1C	[10]
10	6.3	582.31	nodulisporic acid D ₃	indole-diterpenoid	C ₃₈ H ₄₈ NO ₄	-4.1	570.3046; 564.3406; 536.3499; 519.3196; 483.2853; 453.3004; 385.1725; 367.1622; 349.1508; 253.1283; 202.0804; 170.1904; 60.0423	3	31.68F1B	[11]
11	6.0	179.0743	7-Hydroxy-4,6-dimethyl-1(3H)-isobenzofuranone	polyketide	C ₁₀ H ₁₁ O ₃	-3.4	-	3	31.68F1B, 62.72F1A, 36.97F1C	[12,13]
12	6.7	211.0971	canadensolide	polyketide	C ₁₁ H ₁₅ O ₄	0	193.0877; 175.0769; 165.0925; 141.0197	3	62.72F1A	[14,15]

13	7.1	523.2237	asperphenamate analogue; Tyr sub Phe	phenylalanine derivatives	C ₃₂ H ₃₀ N ₂ O ₅	1.0	505.3368; 256.1347; 238.1242; 134.0967	2	31.68F1B, 62.72F1A, 36.97F1C	[16]
14	7.5	259.0606	alternariol	benzocoumarin	C ₁₄ H ₁₁ O ₅	0	-	3	31.68F1B	[17]
15	7.5	607.1083	xanthoepocin	polyketide	C ₃₀ H ₂₃ O ₁₄	0	589.0986; 571.0874; 511.0656; 469.0561; 380.0502; 313.0381; 287.0562; 275.0553; 259.0594; 243.0286; 75.7144	2	31.68F1B	[18]
16	7.5	487.2709	andrastin A	meroterpenoid	C ₂₈ H ₃₉ O ₇	-2.1	427.2476; 409.2382; 395.2220; 377.2119; 243.1754; 183.0657	2	31.68F1B	[19]
17	7.6	489.2671	andrastin B	meroterpenoid	C ₂₈ H ₄₁ O ₇	1.7	449.2322; 428.2501; 410.2356; 396.2200; 243.1831; 225.1616; 183.0701	3	31.68F1B	[19]
18	7.7	429.2430	andrastin D	meroterpenoid	C ₂₆ H ₃₇ O ₅	1.0	410.2400; 396.2190; 378.2111; 244.1802; 151.0384	3	31.68F1B	[19]
19	7.7	519.1899	rubratoxin B	nanodride polyketide	C ₂₆ H ₃₁ O ₁₁	4.4	292.0955; 282.0752; 250.0851	3	31.68F1B, 36.97F1C	[20]
20	7.7	625.3936	citrinadin A	indole alkaloid	C ₃₅ H ₅₃ N ₄ O ₆	-3.9	593.0722; 551.0601; 367.2100; 281.1600	3	31.68F1B	[21]
21	8.2	507.232	asperphenamate	phenylalanine derivative	C ₃₂ H ₃₁ N ₂ O ₄	0.4	377.3205; 256.1341; 238.1238; 224.1077; 134.0963	2	31.68F1B, 36.97F1C, 62.72F1A	[16]

22	4.9	366.1812	brevianamide A	indole alkaloid	C ₂₁ H ₂₄ N ₃ O ₃	3.5	348.1708; 338.1867; 321.1603; 235.1445; 176.1068; 54.0934	3	62.72F1A, 36.97F1C	[22]
23	5.1	390.18	roquefortine C	diketopiperazine	C ₂₂ H ₂₄ N ₅ O ₂	-0.3	322.1290; 193.0723	3	62.72F1A	[23]
24	7.7	493.2815	fungisporin: cyclo-(Phe-Phe-Val-Val)	cyclic tetrapeptide	C ₂₈ H ₃₇ N ₄ O ₄	0.6	465.2788; 394.2130; 346.2133; 295.1447; 247.1454; 219.1501; 199.1445; 120.0810	2	62.72F1A	[24]
25	6.5	509.2759	cyclo-(Tyr-Phe-Val-Val)	cyclic tetrapeptide	C ₂₈ H ₃₇ N ₄ O ₅	-1.0	410.2088; 365.1879; 346.2130; 311.1393; 283.1457; 247.1457; 219.1501; 199.1463; 136.0762; 120.0913	2	62.72F1A, 36.97F1C	[24]
26	7.4	532	cyclo-(Phe-Trp-Val-Val)	cyclic tetrapeptide	C ₃₀ H ₃₈ N ₅ O ₄	-0.6	504.2950; 433.2282; 385.2233; 334.1559; 286.1519; 258.1613; 247.1449; 199.1477; 159.0923; 120.0829	2	62.72F1A, 31.68F1B	[24]
27	6.9	562	cyclo-(Tyr-Trp-Val-Ile)	cyclic tetrapeptide	C ₃₁ H ₄₀ N ₅ O ₅	-0.2	534.3136; 463.2361; 449.2158; 399.2369; 350.1498; 322.1569; 300.1726; 286.1556; 258.1624; 213.1596; 159.0920; 136.0750	2	31.68F1B	[24]
28	7.5	575.0833	xanthomegnin	binaphthoquinone	C ₃₀ H ₂₃ O ₁₂	22*	557.0644; 529.0737; 497.0429; 459.0681; 243.0282	3	36.97F1C, 31.68F1B	[25]
29	6.5	471.2722	communesin G	indole alkaloid	C ₂₉ H ₃₅ N ₄ O ₂	-6.8	439.2459; 261.1839; 233.0794	3	31.68F1B	[26]

30	2.5	209.0474	viticolin C	tropolone	C ₁₀ H ₉ O ₅	1.9	191.0344; 181.0497; 163.0390; 147.0440	3	36.97F1C	[27]
31	6.2	446.1970	fumiquinazoline A	imidazoindole	C ₂₄ H ₂₄ N ₅ O ₄	10	228.1018; 200.1068; 132.0487; 175.0364	3	36.97F1C, 62.72F1A	[28]
32	5.7	596.3093	JBIR-113	cyclodespsipeptide	C ₃₁ H ₄₂ N ₅ O ₇	1.5	485.2384; 467.2200; 409.2340; 356.1597; 338.2077; 280.1658; 209.1296; 140.0705	2	36.97F1C	[29]
33	10.7	263.2365 (M+H- H ₂ O)	conjugated linoleic acid (9E,11E) - primary metabolite	fatty acid	C ₁₈ H ₃₁ O	2.3	245.2311; 189.1600; 175.1515; 163.1522; 149.1302; 123.1240	3	36.97F1C, 31.68F1B	GNPS library
34	4.8	368.1857	brevianamide E	indole alkaloid	C ₂₁ H ₂₆ N ₃ O ₃	0.5	348.1715; 338.1873; 321.1608; 310.1924; 235.1451; 176.1075	3	36.97F1C	[22]
35	5	390.1927	tryhistatin	diketopiperazine	C ₂₂ H ₂₄ N ₅ O ₂	0.5	322.1304; 193.0731	3	62.72F1A	[30]
36	1.5	173.0570	4-(3-furanyl)- benzaldehyde	polyketide	C ₁₁ H ₈ O ₂	11*	130.9707; 112.1531; 96.9574	3	86F1C, 25.88F1C	[31]
37	2.4	167.0703	2',4'-dihydroxy-3'- methylacetophenone	polyketide	C ₉ H ₁₁ O ₃	-1.2	149.0591; 137.9860; 130.0076; 125.0595; 114.9765; 108.9610; 96.9611; 79.0490	3	86F1C, 25.88F1C	[32]
38	7.7	515.2657	lucidenic acid D	triterpenoid	C ₂₉ H ₃₉ O ₈	-1.0	419.2328; 392.1592; 363.1694; 307.1118; 216.9408; 184.0787; 172.1043;	3	86F1C, 25.88F1C	[33]
39	11.2	475.2885	methyl lucidenate N	triterpenoid	C ₂₈ H ₄₃ O ₆	-14.3*	419.2301; 363.1702; 307.1100; 250.1831	3	86F1C, 25.88F1C	[33]

40	9.0	791.2353	cladochrome D	perylenequinone	C ₄₄ H ₃₉ O ₁₄	0.5	663.1892; 637.2188; 601.2106; 552.1516; 515.1664; 483.1454; 473.1285; 443.1058; 404.1078; 359.6608; 317.2110; 301.5874; 286.8502; 253.0100; 195.1283; 163.0770; 105.0340;	3	86F1C	[34]
41	12.3	593.4031	5 α ,8 α -epidioxy- 24(R)- methylcholesta-6,22- dien-3 β -D- glucopyranoside	sterol	C ₃₄ H ₅₇ O ₈	-5.1	495.2724; 451.3066; 393.1087; 355.0716; 307.1878; 251.0862; 223.0799; 165.0840; 147.0609	3	86F1C	[35]
42	7.7	493.2807	preaustinoid D	meroterpenoid	C ₂₇ H ₄₁ O ₈	1.2	394.2118; 346.2113; 295.1448; 267.1487; 247.1447; 219.1497; 120.0802	3	86F1C, 25.88F1C	[36]
43	12.3	685.4373	icosalide A3	cyclic peptide	C ₃₄ H ₆₁ N ₄ O ₁₀	2.9	555.2990; 527.1161; 478.3198; 442.3124; 378.3267; 341.0226; 195.1078; 133.0927	3	86F1C, 25.88F1C	[27]
44	7.5	470.3380	cytochalasin R	indole alkaloid	C ₂₈ H ₄₀ NO ₅	18	452.3407; 236.1507	3	B2F1D	[37]
45	7.8	279.0750	betulinan C	pyranone	C ₁₈ H ₁₅ O ₃	-29.7*	219.9331; 195.9081; 149.0201	3	B2F1D	[38]
46	1.8	184.0530	3-acetyl-5-isopropyl- pyrrolidine-2,4-dione	pyrrolidine alkaloid	C ₉ H ₁₄ NO ₃	11	142.9600; 98.9577	3	B2F1D	[39]
47	9.8	444.2747	stachybotrin A	alkaloid	C ₂₆ H ₃₈ NO ₅	0.0	426.2615; 416.2809; 402.2657; 360.1798;	3	B2F1D	[40]

							334.1642; 233.2270; 226.1070; 212.0551; 194.0448; 175.1481			
48	4.7	235.0969	alternariphent A1	phenol derivative	C ₁₃ H ₁₅ O ₄	0.9	207.1034; 194.9547; 191.1058; 163.1119; 154.9656; 86.0951	3	B2F1D	[41,42]
49	6.6	583.3277	6-O-propionyl-6,16-O-dideacetylhelvolic acid	nortriterpenoid	C ₃₄ H ₄₇ O ₈	0.0	543.1447; 421.2726; 391.2717; 281.1542; 229.1223; 123.1101	3	B2F1D	[43]

*Peak ions of low intensity (molecular formula prediction with high ppm errors)

Table S3: Bioactivity (% inhibition at 100 µg /ml concentration) of the foam-derived fungi. Mean values are based on triplicate measurements. Plant pathogens: Pss, *P. syringae*; Xc, *X. campestris*, Ea, *E. amylovora*; Rs, *R. solanacearum*; Pi, *P. infestans*; Mo, *M. oryzae*; Bc, *B. cinerea*. Human pathogens include the ESKAPE panel Efm, *E. faecium*; MRSA, methicillin-resistant *S. aureus*; Kp, *K. Pneumoniae*; Ab, *A. baumannii*; Psa, *P. aeruginosa*; Ec, *E. coli*; Ca, *C. albicans*; Cn, *C. neoformans*. 0.5% DMSO was used as a solvent control. (+) control see Table S4.

Sea foam derived fungus	Anti-microbial activity against plant pathogens							Anti-microbial activity against human pathogens								Anti-cancer Activity		
	Anti-bacterial				Anti-fungal			Anti-ESKAPE					Anti-fungal					
	Pss	Xc	Ea	Rs	Pi	Mo	Bc	Efm	MRSA	Kp	Ab	Psa	Ec	Ca	Cn	HaCaT	A549	MB231
<i>Plectospaerella</i> sp.	-	77	-	-	44	-	-	-	50	-	-	-	-	-	-	-	-	-
<i>Cladosporium</i> sp.	-	67	25	-	-	-	-	-	54	-	-	-	-	-	-	-	-	-
<i>Emericellopsis</i> sp.	-	42	-	-	-	-	-	-	37	-	-	-	-	-	-	-	-	-
<i>Penicillium</i> :31.68F1B	-	47	27	-	97	98	74	99	99	-	-	-	-	80	100	-	-	-

<i>Penicillium</i> :36.97F1C	-	29	-	-	86	100	41	100	33	-	20	-	-	66	82	-	-	-
<i>Penicillium</i> :62.72F1A	-	33	-	-	80	-	30	-	-	-	-	-	-	-	-	-	-	26
(+) control	95	95	66	99	96	97	98	99	93	99	96	99	88	85	100	41	32	62
(-) control (solv./PDA)	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

Table S4. Test pathogens and their specific conditions for bioactivity testing

Human pathogen	Strain	Medium	Optical density	Preculture incubation temperature (°C)	Positive control	Incubation temperature (°C)	Incubation time (h)
<i>Enterococcus faecium</i>	DSM 20478	M92	0.01	37	Ampicillin	37	4
Methicillin-resistant <i>Staphylococcus aureus</i> (MRSA)	DSM 18827	TSB12	0.01	28	Chloramphenicol	37	5
<i>Klebsiella pneumoniae</i>	DSM 30104	TSB12	0.01	28	Chloramphenicol	37	5
<i>Acinetobacter baumannii</i>	DSM 30007	TSB12	0.01	28	Ampicillin	37	5

<i>Pseudomonas aeruginosa</i>	DSM 1128	TSB12	0.01	37	Chloramphenicol	37	5
<i>Escherichia coli</i>	DSM 1576	TSB12	0.01	28	Chloramphenicol	37	5
<i>Candida albicans</i>	DSM 1386	M186	0.03	28	Nystatin	37	5
<i>Cryptococcus neoformans</i>	DSM 6973	M186	0.03	28	Amphotericin B	28	6-7
Plant pathogens							
<i>Pseudomonas syringae</i> pv. <i>aptata</i>	DSM 50252	TSB12	0.03	28	Chloramphenicol	28	9
<i>Xanthomonas campestris</i>	DSM 2405	TSB12	0.03	28	Chloramphenicol	28	7
<i>Erwinia amylovora</i>	DSM 50901	TSB12	0.03	28	Chloramphenicol	28	9
<i>Ralstonia solanacearum</i>	DSM 9544	M1	0.03	28	Tetracycline	28	9
<i>Phytophthora infestans</i>	CBS 120920	PB*	(10 ⁴ spores)	22	Cycloheximide	22	74-96
<i>Magnaporthe oryzae</i>	DSM 62938	SA*	(10 ⁴ spores)	22	Nystatin	22	74-96
<i>Botrytis cinerea</i>	DSM 5145	MA	(5×10 ⁴ spores)	22	Boscalid	25	72

M92: 30 g tryptic soy broth, 3 g yeast extract, 1000 mL H₂O, pH 7.0-7.2

TSB12: 12 g tryptic soy broth, 5 g NaCl, 1000 mL H₂O

M186: 10 g glucose, 5 g peptone from soybeans, 3 g yeast extract, 3 g malt extract, 1000 mL H₂O

M1: 5 g peptone, 3 g peptone from meat, 1000 mL H₂O

PB: 150 g pea, 5 g glucose, 0.1 mg thiamine HCl, 1000 mL H₂O

SA: 10 g peptone, 20 g glucose, 15 g agar, 1000 mL H₂O, pH 5.6 ± 0.02

MA: 10 g malt extract, 15 g agar, 1000 mL H₂O

*: In addition to broth, 15 g agar was added for pre-cultivation of fungi

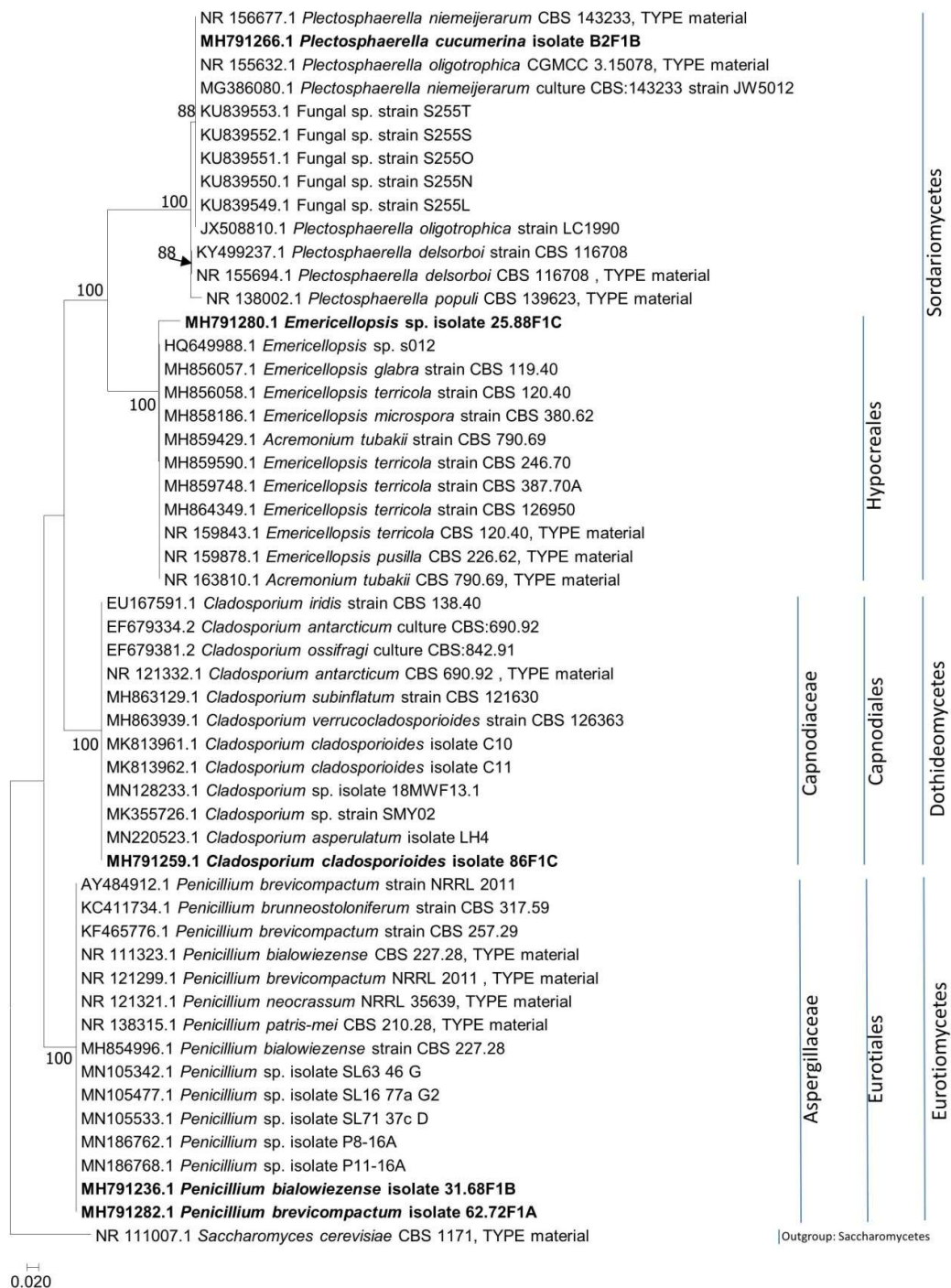


Figure S1. Phylogenetic analysis of selected foam-derived fungal strains based on ITS region. Reference sequences were downloaded from the NCBI database. *Penicillium* sp. isolate 36.97 F1C (only identified by 18S rRNA) is located in another phylogenetic tree (Figure S2).

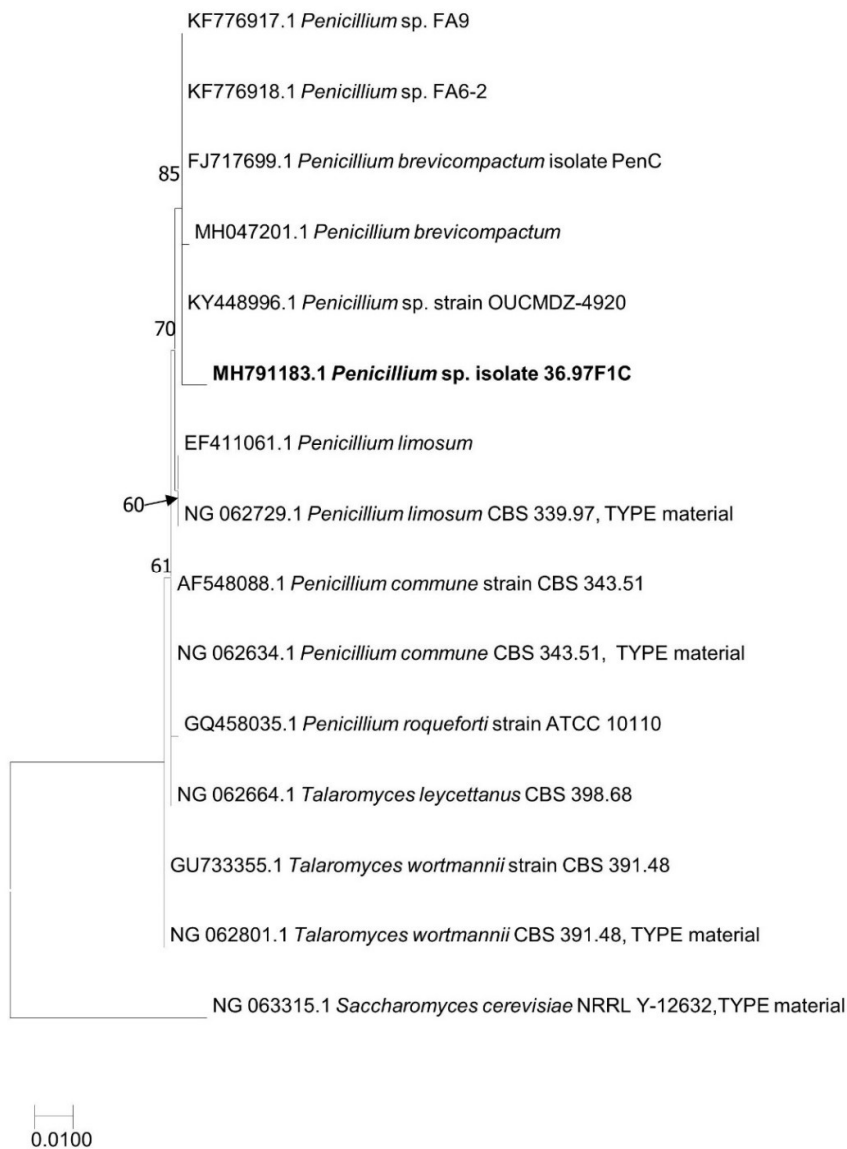


Figure S2. Phylogenetic analysis of selected strain *Penicillium* sp. isolate 36.97 F1C based on 18SrRNA gene. Reference sequences were downloaded from the NCBI database.

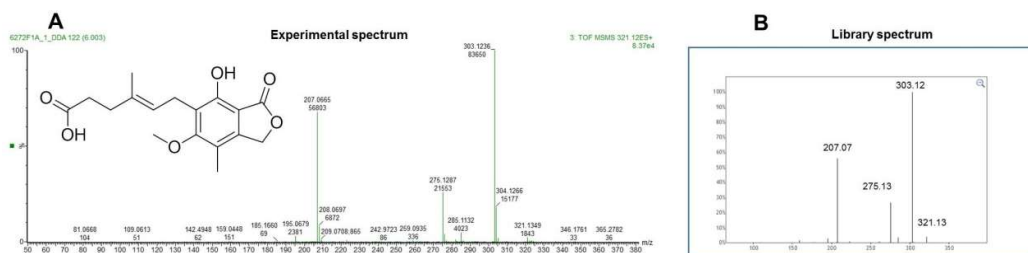


Figure S3: MS/MS spectrum of (A) node m/z [M+H]⁺ 321.1342 annotated as mycophenolic acid and (B) library MS/MS spectrum of mycophenolic acid.

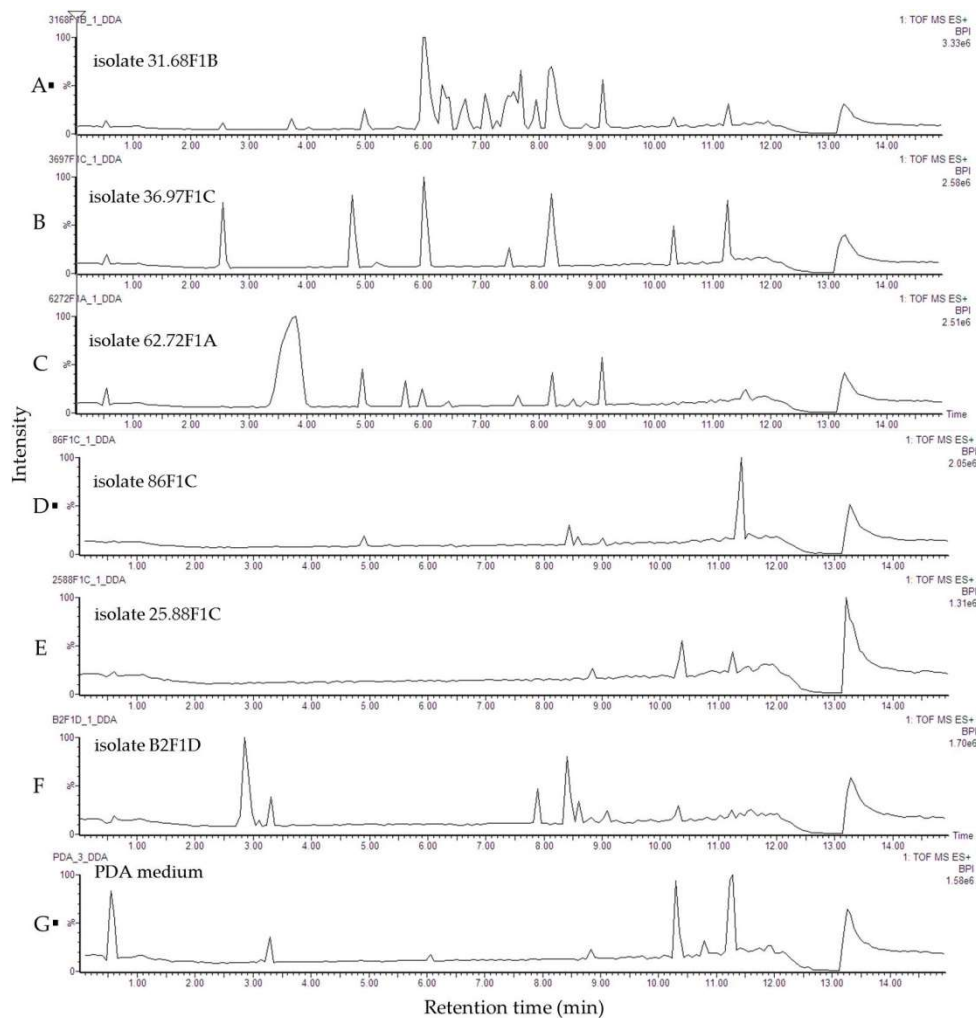


Figure S4. Base peak chromatograms (acquired by an UPLC-QToF-MS/MS system in positive mode) of extract of (A) *Penicillium* sp. 31.68F1B, (B) *Penicillium* sp. 36.97F1C, (C) *Penicillium* sp. 62.72F1A, (D) *Cladosporium* sp. 86F1C, (E) *Emericellopsis* sp. 25.88F1C, (F) *Plectosphaerella* sp. B2F1D, and (G) PDA medium

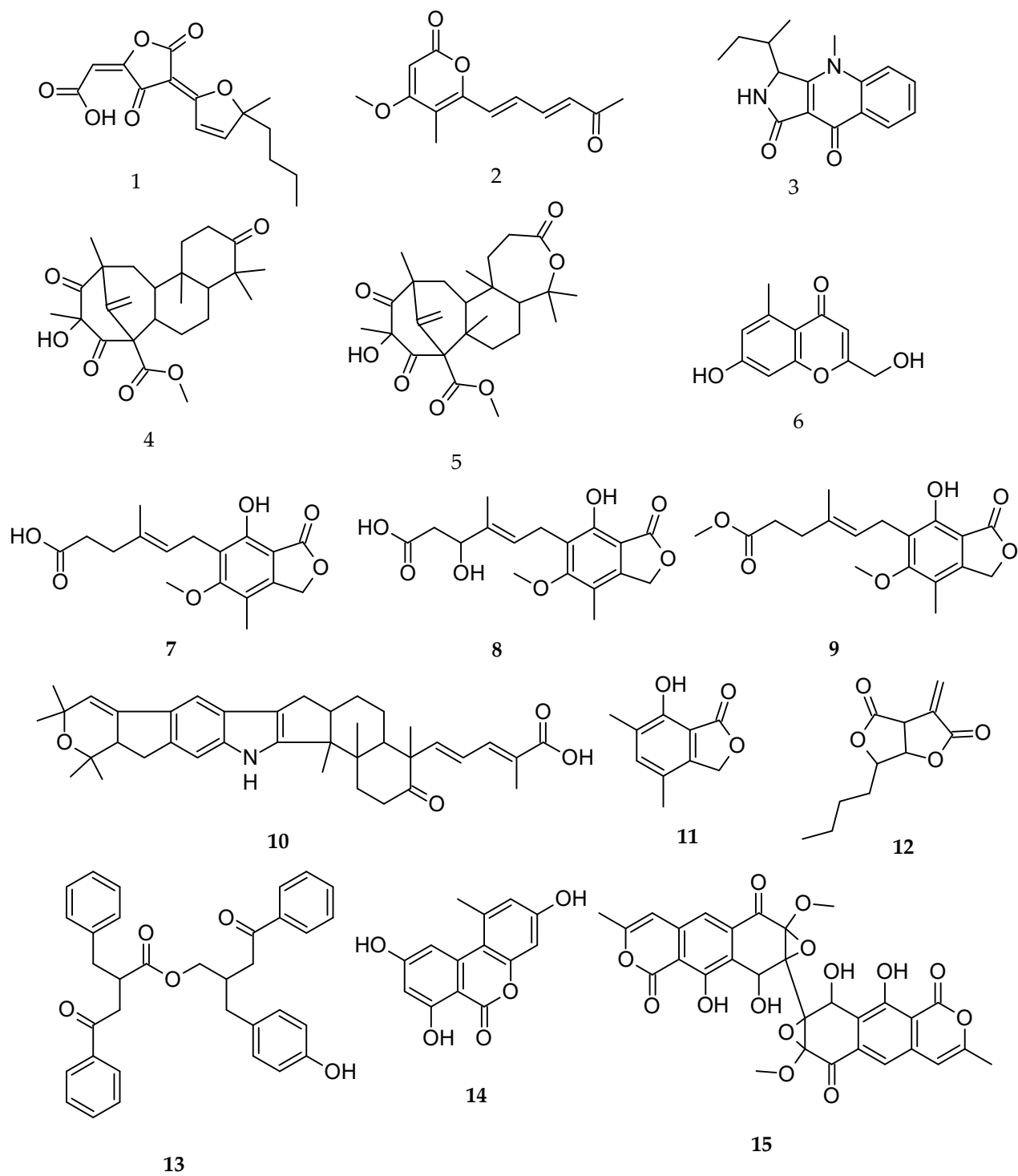
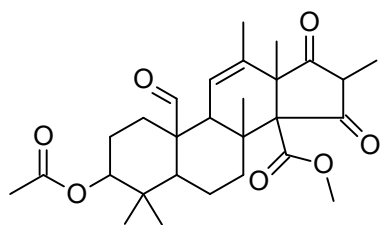
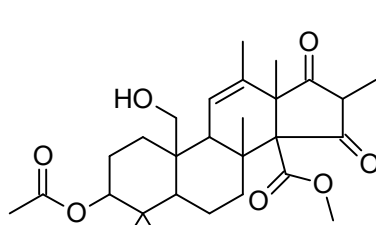


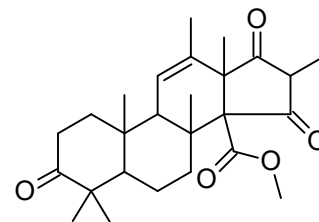
Figure S5. Chemical structures of metabolites annotated in the selected sea foam –derived fungi. Numbers correspond to putatively annotated known compounds, which are reported in Table S2.



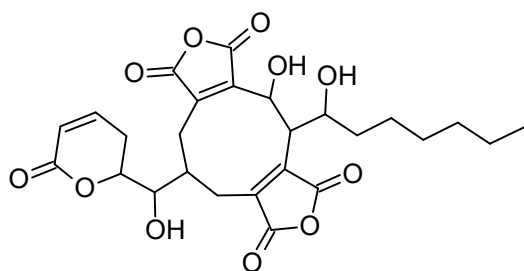
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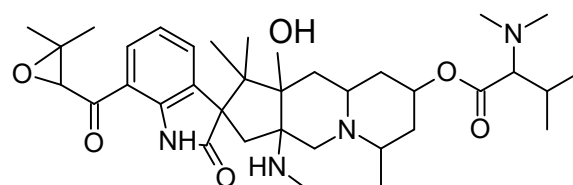
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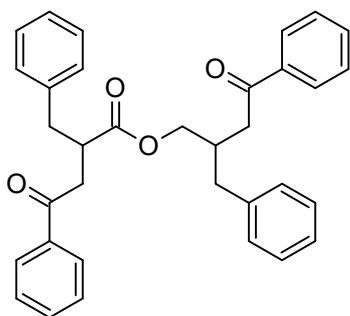
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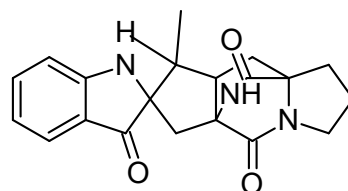
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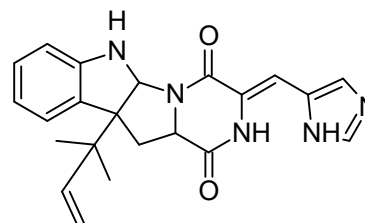
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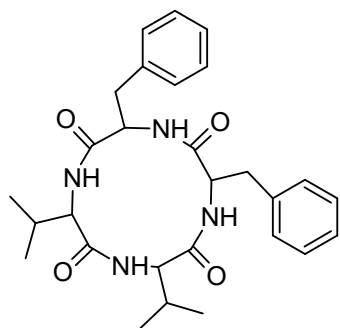
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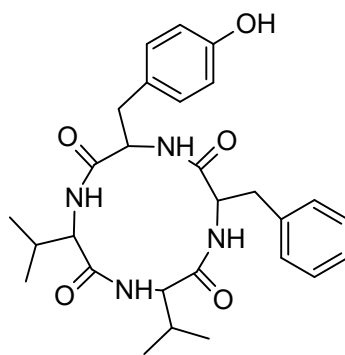
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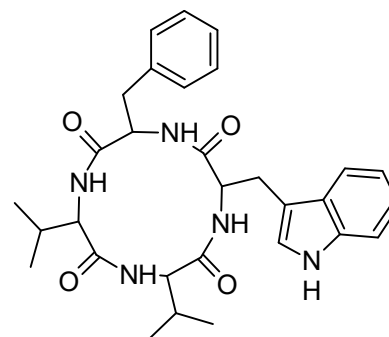
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Figure S5. (continued)

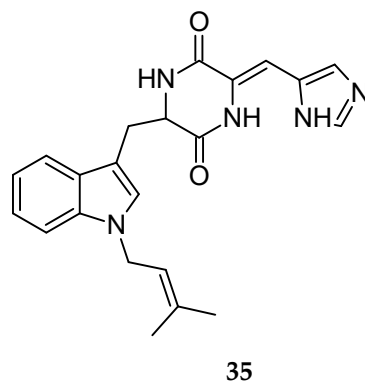
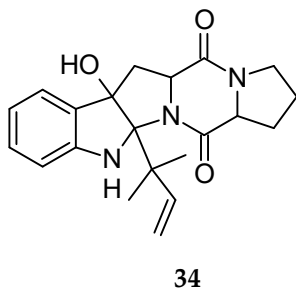
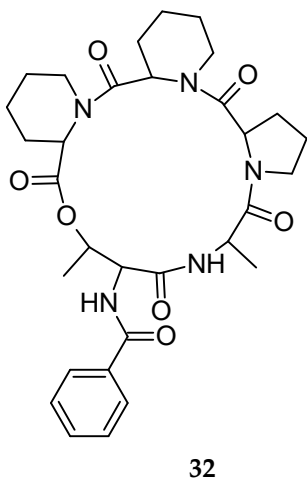
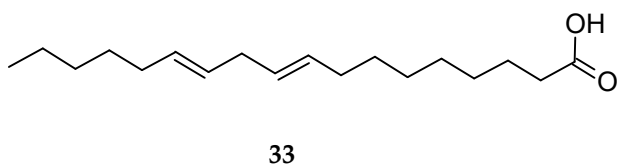
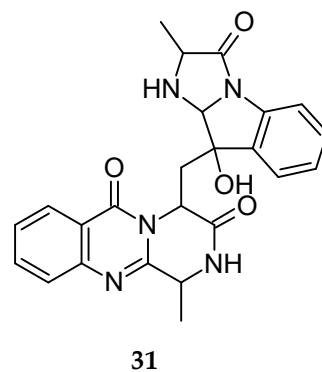
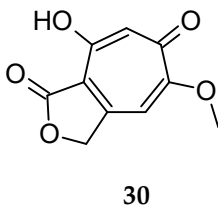
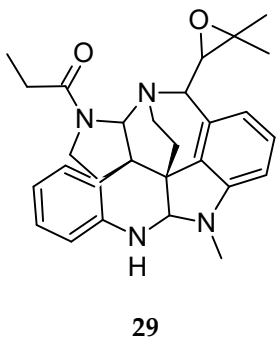
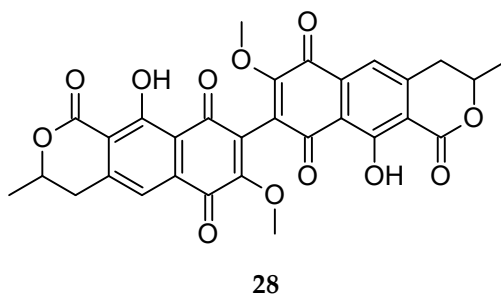
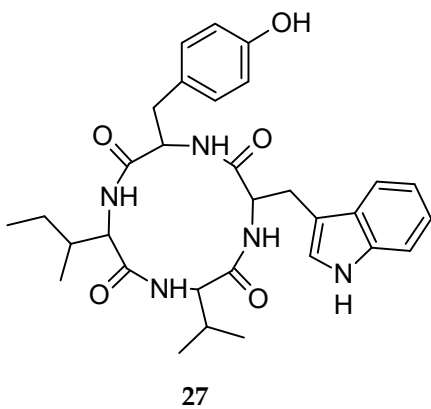
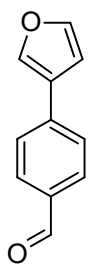
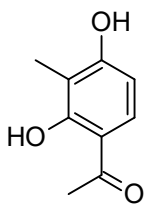


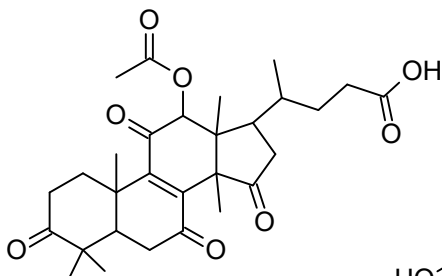
Figure S5. (continued)



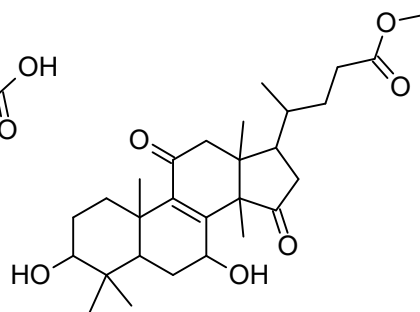
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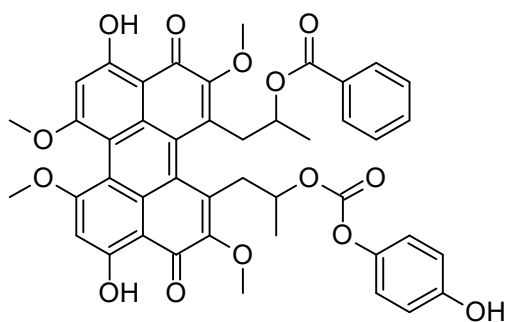
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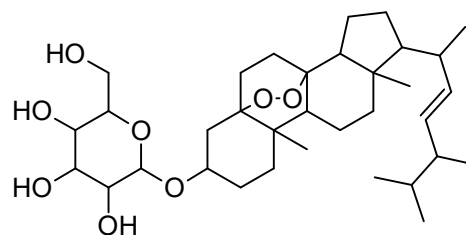
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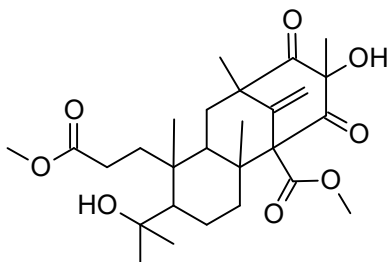
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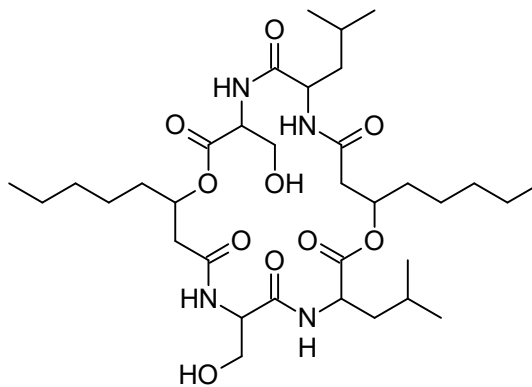
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Figure S5. (continued)

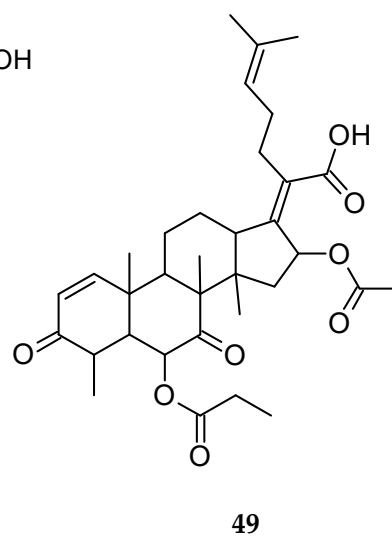
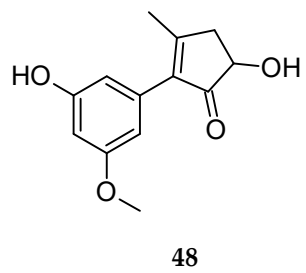
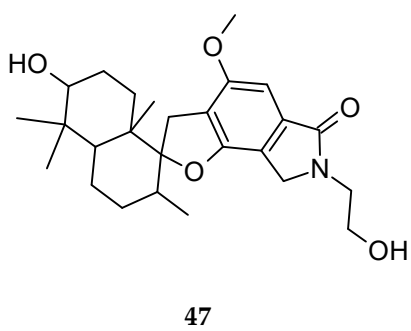
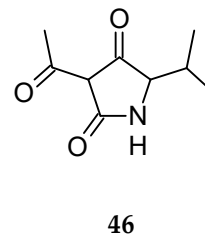
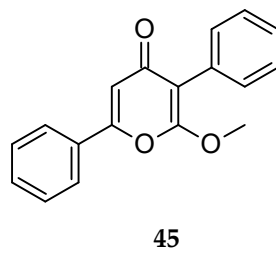
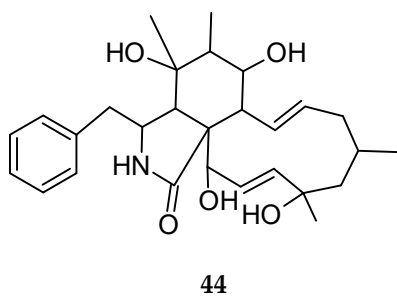


Figure S5. (continued)

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