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LAMPIRAN

Lampiran 1. Senyawa rumput laut *Caulerpa racemosa*

No	Senyawa	Area %	Berat molekul (MW)	Smiles	PubChem ID
1	Methane, (Methylsulfinyl)(Methylthio)-	1.71	124	CS(CS(=O)C)C	99129
2	Silane, Dimethoxymethyl-	0.62	106	CO[SiH](C)OC	85623
3	2-Butanon, 3-Hydroxy	0.16	88	CC(C(=O)C)O	179
4	2,2-Dimethoxybutane	0.35	118	CCC(C)(OC)OC	137941
5	5-[(Propylsulfanyl)Methyl]-2,4-Imidazolidinedione #	0.02	188	CCCS(CC1C(=O)NC(=O)N1)C	560576
6	Hexanal	0.36	100	CCCCCC=O	6184
7	(Z)-3-Chloropropenamide	0.16	105	C(CCl)N=C=O	16035
8	3-Methoxy-1,1-Dimethylsilacyclopentane	0.35	144	CCC=CO[Si](C)(C)C	88325
9	1,1-Dieethoxy-2-Ethylhexane	0.19	202	CCCC(C)C(OCC)OCC	12504289
10	2-Decenal, (E)-	0.09	154	CCCCCCCC=CC=O	5283345
11	Tetradecane	0.21	198	CCCCCCCCCCCCCCC	12389
12	1s-Neomenthyl Acetate	0.01	198	CC1CCC(C(C1)OC(=O)C)C(C)C	75699
13	Nonanoic Acid, 9-Oxo-, Methyl Ester	0.08	186	COC(=O)CCCCCCCC=O	74732
14	3-Buten-2-One, 4-(2,6,6-Trimethyl-2-Cyclohexen-1-Yl)-	0.05	192	CC1=CCCC(C1C=CC(=O)C)C(C)C	5282108
15	R-(+)-Methyl-3-Isopropyl-6-Oxoheptanoate	0.58	200	CC(C)C(CCC(=O)C)CC(=O)OC	11390031
16	4-(2,6,6-Trimethyl-2-Cyclohexen-1-Yl)-2-Pentanone	0.24	208	CC1=CCCC(C1C(C)CC(=O)C)C(C)C	579168
17	Phenol, 3,5-Bis(1,1-Dimethylethyl)-	0.17	206	CC(C)(C)C1=CC(=CC(=C1O)C)C(C)C	70825

18	<i>4-(2,2,6-Trimethyl-7-Oxa-Bicyclo[4.1.0]Hept-1-Yl)-But-3-En-2-One</i>	0.41	208	CCC(=O)OC1CC(CC=C1C)C(=C)C	7336
19	<i>Phenol, 2,4-Bis(1,1-Dimethylethyl)-</i>	0.58	206	CC(C)(C)C1=CC(=C(C=C1O)C(C)C)	7311
20	<i>Dodecanoic Acid, Methyl Ester</i>	0.18	214	CCCCCCCCCC(=O)OC	8139
21	<i>2(4h)-Benzofuranone, 5,6,7,7a-Tetrahydro-4,4,7a-Trimethyl-</i>	0.40	180	CC1(CCCC2(C1=CC(=O)O2)C)C	27209
22	<i>1,8(2H,5H)-Naphthalenedione, Hexahydro-8a-Methyl-, Cis</i>	0.52	180	CC12C(CCCC1=O)CCCC2=O	580218
23	<i>2(4h)-Benzofuranone, 5,6,7,7a-Tetrahydro-4,4,7a-Trimethyl-</i>	1.60	180	CC1(CCCC2(C1=CC(=O)O2)C)C	27209
24	<i>Octanoic Acid, 3-Pentadecyl Ester</i>	0.29	354	CCCCCCCCCC(=O)OC	581291
25	<i>Hexadecane</i>	0.39	226	CCCCCCCCCC	11006
26	<i>7-Hexadecenoic Acid, Methyl Ester, (Z)-</i>	0.46	268	CCCCCCCC=CCCCC(=O)OC	5364431
27	<i>Heneicosane</i>	0.45	296	CCCCCCCCCC	12403
28	<i>4-(1,5-Dihydroxy-2,6,6-Trimethyl-2-Cyclohexen-1-Yl)-3-Buten-2-One</i>	0.37	224	CC1=CCC(C(C1=CC(=O)C)O)(C)O	5363693
29	<i>Heneicosane</i>	0.21	296	CCCCCCCCCC	12403
30	<i>Tridecanol, 2-Ethyl-2-Methyl-</i>	0.22	242	CCCCCCCC(C)CO	545928
31	<i>As-Indacene, Dodecahydro-4-(1-Octylnonyl)-</i>	0.59	402	CCCCCCCC(CCCCCC)C1CC2CCCC2C3C1CCC3	296574
32	<i>Heptadecane</i>	0.70	240	CCCCCCCCCC	12398
33	<i>4-(4-Hydroxy-2,2,6-Trimethyl-7-Oxa-Bicyclo[4.1.0]Hept-1-Yl)-But-3-En-2-One</i>	0.79	224	CC(=O)C=CC12C(CC1(O2)C)O(C)C	636523
34	<i>Tetradecanoic Acid, Methyl Ester</i>	0.98	242	CCCCCCCCCC(=O)OC	31284

35	<i>Hexadecane, 2,6,10,14-Tetramethyl-</i>	0.38	282	CCC(C)CCCC(C)CCCC(C)CCCC(C)C	12523
36	<i>Hexadecane, 2-Methyl-</i>	0.19	240	CCCCCCCCCC(C)CCCCCCC	15266
37	<i>Pentadecane, 8-Hexyl-</i>	0.22	296	CCCCCCCC(CCCCC)CCCCCCC	300518
38	<i>Heptadecane, 2-Methyl-</i>	0.22	254	CCCCCCCCCCCCCCCC(C)C	15265
39	<i>3-Methylheptadecane</i>	0.45	254	CCCCCCCCCCCCCCCC(C)CC	94321
40	<i>2(4h)-Benzofuranone, 5,6,7,7a-Tetrahydro-6-Hydroxy-4,4,7a-Trimethyl-, (6s-Cis)-</i>	0.44	196	CC1(CC(CC2(C1=CC(=O)O2)C)O)C	100332
41	<i>1-Octadecanethiol</i>	0.29	286	CCCCCCCCCCCCCCCCCS	17905
42	<i>Octadecane</i>	0.50	254	CCCCCCCCCCCCCCCCCCC	11635
43	<i>2(4h)-Benzofuranone, 5,6,7,7a-Tetrahydro-6-Hydroxy-4,4,7a-Trimethyl-, (6s-Cis)-</i>	1.35	196	CC1(CC(CC2(C1=CC(=O)O2)C)O)C	100332
44	<i>Pentadecanoic Acid, Methyl Ester</i>	1.02	256	CCCCCCCCCCCCC(=O)OC	23518
45	<i>2,6,10-Trimethyl,14-Ethylene-14-Pentadecne</i>	2.69	278	CC(C)CCCC(C)CCCC(C)CCCC(=C)C=C	10446
46	<i>2-Pentadecanone, 6,10,14-Trimethyl-</i>	1.87	268	CC(C)CCCC(C)CCCC(C)CCCC(=O)C	10408
47	<i>2,6,10-Trimethyl,14-Ethylene-14-Pentadecne</i>	1.46	278	CC(C)CCCC(C)CCCC(C)CCCC(=C)C=C	10446
48	<i>1-Butyl 2-(8-Methylnonyl) Phthalate #</i>	0.30	362	CCCCOC(=O)C1=CC=CC=C1C(=O)OCCCCCCCC(C)C	39180
49	<i>2,6,10-Trimethyl,14-Ethylene-14-Pentadecne</i>	1.93	278	CC(C)CCCC(C)CCCC(C)CCCC(=C)C=C	10446
50	<i>9-Hexadecenoic Acid, Methyl Ester, (Z)-</i>	1.97	268	CCCCCCC=CCCCCCCC(=O)OC	643801
51	<i>Hexadecanoic Acid, Methyl Ester</i>	9.83	270	CCCCCCCCCCCCCCCC(=O)OC	8181
52	<i>3,7,11,15-Tetramethylhexadec-1-En-3-OI</i>	0.03	296	CC(C)CCCC(C)CCCC(C)CCCC(C)(C=C)O	10453

53	2,3-Dimethyl-1-Undecen-3-OI	0.07	198	CCCCCCCCCC(C)(C(=C)C)O	559021
54	1,2-Benzenedicarboxylic Acid, Butyl 8-Methylnonyl Ester	0.08	362	CCCCOC(=O)C1=CC=CC=C1C(=O)OCCCCCC(C)C	39180
55	Hexadecanoic Acid, Ethyl Ester	1.29	284	CCCCCCCCCC(C)CCCCCCC(=O)OCC	12366
56	Cis-9-Hexadecenal	0.65	238	CCCCCCCC=CCCCCCCC=O	5364643
57	Heneicosanoic Acid, Methyl Ester	0.82	340	CCCCCCCCCCCCCCCCCCCC(=O)OC	22434
58	Methyl 2-Ethylhexyl Phthalate	0.45	292	CCCC(CC)COC(=O)C1=CC=CC=C1C(=O)OC	596027
59	2-Norpinanol, 3,6,6-Trimethyl-	0.52	154	CC1CC2CC(C1O)C2(C)C	558905
60	8,9,12-Octadecadienoic Acid (Z,Z)-, Methyl Ester	21.3 1	294	CCCCC=CCC=CCCCCCCC(=O)OC	528441
61	Palmitaldehyde, Diallyl Acetal	1.45	338	CCCCCCCCCCCCCCC(OCC=C) OCC=C	319417400
62	Ethyl (9z,12z)-9,12-Octadecadienoate #	4.55	308	CCCCC=CCC=CCCCCCCC(=O)OCC	5282184
63	10-Nonadecenoic Acid, Methyl Ester	0.68	310	CCCCCCCC=CCCCCCCC(=O)OC	5364486
64	6,9,12-Octadecatrienoic Acid, Methyl Ester	3.29	292	CCCCC=CCC=CCC=CCCCC(=O)OC	5362805
65	Linolensaeuremethylester	3.57	292	CCC=CCC=CCC=CCCCCCCC(=O)OC	5367462
66	9-Octadecynoic Acid, Methyl Ester	3.61	294	CCCCCCCC#CCCCCCCC(=O)OC	534587
67	4,8,12,16-Tetramethylheptadecan-4-Olide	2.08	324	CC(C)CCCC(C)CCCC(C)CCCC1(CC(=O)O1)C	567149

68	<i>Methyl Eicosa-5,8,11,14,17-Pentaenoate</i>	3.59	316	CCC=CCC=CCC=CCC=CCC=CCC CC(=O)OC	5367433
69	<i>7,9-Dimethyl-8-Nitrobicyclo[4.3.1]Decan-10-One</i>	0.68	225	CC1C2CCCC(C2=O)C(C1[N+](=O)[O-])C	534696
70	<i>1,2-Benzenedicarboxylic Acid, Ditridecyl Ester</i>	0.38	530	CCCCCCCCCCCOOC(=O)C1=C C=CC=C1C(=O)OCCCCCCCCCCC CCC	8379
71	<i>Docosanoic Acid, Methyl Ester</i>	0.72	354	CCCCCCCCCCCCCCCCCCCCCCCC (=O)OC	13584
72	<i>1,2-Benzenedicarboxylic Acid</i>	0.83	390	CCCC(CC)COC(=O)C1=CC=CC= C1C(=O)OCC(CC)CCCC	8343
73	<i>Ethyl Docosanoate</i>	0.03	368	CCCCCCCCCCCCCCCCCCCCCCCC (=O)OCC	22199
74	<i>Tricosanoic Acid, Methyl Ester</i>	0.06	368	CCCCCCCCCCCCCCCCCCCCCCCC C(=O)OC	75519
75	<i>Dotriacontane</i>	0.08	450	CCCCCCCCCCCCCCCCCCCCCCCC CCCCCCCCCCCC	11008
76	<i>Tetracosanoic Acid, Methyl Ester</i>	0.18	382	CCCCCCCCCCCCCCCCCCCCCCCC CC(=O)OC	75546
77	<i>N-[1-(1-Adamantan-1-Yl-Propyl)-2,5-Dioxo-4-Trifluoromethyl-Imidazolidin-4-Yl]-4-Methoxy-Benzamide</i>	0.04	493	CCC(C12CC3CC(C1)CC(C3)C2)N 4C(=O)C(NC4=O)(C(F)(F)F)NC(=O))C5=CC=C(C=C5)OC	4213120
78	<i>17-Pentatriacontene</i>	0.05	490	CCCCCCCCCCCCCCCCCCC=CCC CCCCCCCCCCCCCCC	5365022
79	<i>2,6,10,14,18,22-Tetracosahexaene, 2,6,10,15,19,23-Hexamethyl-, (All-E)-</i>	0.18	410	CC(=CCCC(=CCCC(=CCCC=C(C) CCC=C(C)CCC=C(C)C)C)C)C	638072

80	<i>1,6,10-Dodecatrien-3-Ol, 3,7,11-Trimethyl-, [S-(Z)]-</i>	0.03	222	CC(=CCCC(=CCCC(C)(C=C)O)C)C	3520128
81	<i>2H-1-Benzopyran-6-Ol, 3,4-Dihydro-2,8-Dimethyl-2-(4,8,12-Trimethyltridecyl)-, [2R-[2R@(4R@,8R@)]]-</i>	0.57	402	CC1=CC(=CC2=C1OC(CC2)(C)CC CC(C)CCCC(C)CCCC(C)C)O	586537
82	<i>2H-1-Benzopyran-6-Ol, 3,4-Dihydro-2,8-Dimethyl-2-(4,8,12-Trimethyltridecyl)-, [2R-[2R@(4R@,8R@)]]-</i>	0.76	402	CC1=CC(=CC2=C1OC(CC2)(C)CC CC(C)CCCC(C)CCCC(C)C)O	586537
83	<i>Hexatriacontane</i>	0.16	506	CCCCCCCCCCCCCCCCCCCCCC	12412
84	<i>1-(4-Ethoxy-2-Hydroxyphenyl)-1-Octadecanone #</i>	0.06	404	CCCCCCCCCCCCCCCC(=O)C1=C(C=C(C=C1)OCC)O	601846
85	<i>.Beta.-Tocopherol</i>	0.15	416	CC1=CC(=C(C2=C1OC(CC2)(C)CC(C)CCCC(C)CCCC(C)C)O	6857447
86	<i>Tetracontane</i>	0.08	562	CCCCCCCCCCCCCCCCCCCCCC	20159
87	<i>Cholesta-4,6-Dien-3-Ol, Benzoate, (3.Beta.)-</i>	0.07	488	CC(C)CCCC(C)C1CCC2C1(CCC3C2C=CC4=CC(CCC34C)OC(=O)C5=CC=CC=C5)C	33010
88	<i>Stigmast-5-En-3-Ol, Oleate</i>	0.10	678	CCCCCCCC=CCCCCCCC(=O)OC1CCC2(C3CCCC4(C(C3CC=C2C1)CCC4C(C)CCC(CC)C(C)C)C)	20831071
89	<i>.Alpha.-Tocopherol-.Beta.-D-Mannoside</i>	0.12	592	CC1=C(C(=C(C2=C1OC(CC2)(C)CC(C)CCCC(C)CCCC(C)C)OC3C(C(C(O3)C(CO)O)O)O)C	570057

90	<i>Stigmasta-5,23-Dien-3-Ol, (3.Beta.)-</i>	0.31	412	CCC(CCC(C)C1CCC2C1(CCC3C2=CC=C4C3(CCC(C4)O)C)C)C(C)C	10740
91	<i>Stigmast-5-En-3-Ol, (3.Beta.)-</i>	0.61	414	CCC(CCC(C)C1CCC2C1(CCC3C2CC=C4C3(CCC(C4)O)C)C)C(C)C	222284
92	<i>Lupeol</i>	0.15	426	CC(=C)C1CCC2(C1C3CCC4C5(CC(C(C5CCC4(C3(CC2)C)C)(C)C)O)C)C	259846

Lampiran 2. Senyawa rumput laut *Sargassum polycystum*

No.	Senyawa	Area %	Berat molekul (MW)	Smiles	PubChem ID
1	1,3,5-Triazine-2,4-Diamine, 6-Chloro-N-Ethyl-	12.85	173	CCNC1=NC(=NC(=N1)N)Cl	13878
2	1,6-Octadien-3-Ol, 3,7-Dimethyl-	2.31	154	CC(=CCCC(C)(C=C)O)C	6549
3	1-Decanol	2.40	158	CCCCCCCO	8174
4	Undecane	6.54	156	CCCCCC	14257
5	N-Formylmorpholine	1.59	115	C1COCCN1C=O	20417
6	Cyclohexasiloxane, Dodecamethyl-	0.99	444	C[Si]1(O[Si](O[Si](O[Si](O[Si](O[Si](O[Si](O[Si](O[Si](O[Si](O[Si](O1)(C)C)(C)C)(C)C)(C)C)(C)C)C)C)C)C)C	10911
7	1,3-Dioxane, 4,6-Dimethyl-	5.00	116	CC1CC(OCO1)C	136893
8	Methenamine	6.73	140	C1N2CN3CN1CN(C2)C3	4101
9	Azulene, 1,2,3,5,6,7,8,8a-Octahydro-1,4-Dimethyl-7-(1-Methylethenyl)-, [1S-(1.Alpha.,7.AI	2.07	204	CC1CCCC2=CCC(CC12C)C(=C)C	2237
10	2(4h)-Benzofuranone, 5,6,7,7a-Tetrahydro-6-Hydroxy-4,4,7a-Trimethyl-, (6S-CIS)-	1.05	196	CC(C)COC(=O)CCC1=CC=CO1	100332
11	2,6,10-Trimethyl,14-Ethylene-14-Pentadecne	2.07	278	CC(C)CCCC(C)CCCC(C)CCCC(=C)C=C	10446
12	2,6,10-Trimethyl,14-Ethylene-14-Pentadecne	1.33	278	CC(C)CCCC(C)CCCC(C)CCCC(=C)C=C	10446
13	Hexadecanoic Acid, Methyl Ester	6.66	270	CCCCCCCCCCCCCCCC(=O)OC	8181
14	Hexadecanoic Acid, Ethyl Ester	1.08	284	CCCCCCCCCCCCCCCC(=O)O	5281

15	<i>9-Octadecenoic Acid (Z)-, Methyl Ester</i>	3.11	296	CCCCCCCCCC=CCCCCCCC(=O)OC	8202
16	<i>6,8-Dioxabicyclo(3.2.1)Octan-4.Beta.-Ol</i>	1.59	130	CCC(C)C(=O)C(=O)O	47
17	<i>Octadecanoic Acid, Methyl Ester</i>	1.30	298	CCCCCCCCCCCCCCCCCC(=O)OC	8201
18	<i>9,19-Cycloergost-24(28)-En-3-Ol, 4,14-Dimethyl-, Acetate, (3.Beta.,4.Alpha.,5.Alpha.)-</i>	3.62	468	CC(=O)OC1CCC2(C(C1(C)C)CCC3(C2CC=C4C3(CCC5(C4CC(CC5)(C)C)C)C)C)	92156
19	<i>Dotriacontane</i>	7.90	450	CCCCCCCCCCCCCCCCCCCCCCCC	11008
20	<i>Heneicosane, 11-Cyclopentyl-</i>	2.60	364	CCCCCCCCCCCCCCCCCCCC1CCCC1	20511
21	<i>1,2-Benzenedicarboxylic Acid</i>	2.05	390	CCCC(C)COC(=O)C1=CC=CC=C1C(=O)OCC(CC)CCCC	8343
22	<i>Triacontane</i>	1.99	422	CC(C)CCCC(C)CCCC(C)CCCC(C)CCCC(C)CCCC(C)C	8089
23	<i>Tetrapentacontane</i>	1.68	758	CCCCCCCCCCCCCCCCCCCCCCCC CCCCCCCCCCCCCCCCCCCCCCCC CCCCCCCCCCCC	521846
24	<i>Triacontane</i>	3.64	422	CC(C)CCCC(C)CCCC(C)CCCC(C)CCCC(C)C	8089
25	<i>1-Hentetracontanol</i>	4.08	592	CCCCCCCCCCCCCCCCCCCC CCCCCCCCCCCCCCCCCCCCO	38627
26	<i>Cyclohexane, Nonadecyl-</i>	2.04	350	CCCCCCCCCCCCCCCCCCCC1C CCCC1	89671

27	<i>Hexacontane</i>	1.73	842	CCCCCCCCCCCCCCCCCCCCCC CCCCCCCCCCCCCCCCCCCCCC CCCCCCCCCCCCCCCCCC	24318
28	<i>Tetracontane</i>	3.68	562	CCCCCCCCCCCCCCCCCCCCCC CCCCCCCCCCCCCCCCCCCCCC	20149
29	<i>Tetracontane</i>	4.80	562	CCCCCCCCCCCCCCCCCCCCCC CCCCCCCCCCCCCCCCCCCCCC	20149
30	<i>Cholest-5-En-3-Ol (3.Beta.)-</i>	1.51	386	CC(C)CCCC(C)C1CCC2C1(CCC3 C2CC=C4C3(CCC(C4)O)C)C	5997

Lampiran 3. Senyawa rumput laut *Halymenia durvillei*

No	Senyawa	Area %	Berat molekul (MW)	Smiles	PubChem ID
1	Benzene, Ethyl-	6.27	106	CCC1=CC=CC=C1	7500
2	Benzene, 1,2-Dimethyl-	5.50	106	CC1=CC=CC=C1C	7237
3	Benzene, Ethenyl-	4.96	104	C=CC1=CC=CC=C1	7501
4	Ethanol, 2-Butoxy-	10.3 1	118	CCCCOCCO	8133
5	2-Butoxyethyl Acetate	0.25	160	CCCCOCCOC(=O)C	8160
6	1-Dodecanol	0.61	186	CCCCCCCCCCCO	8193
7	1-Tetradecanol	2.79	214	CCCCCCCCCCCCCO	8209
8	Tetradecane	0.35	198	CCCCCCCCCCCC	12389
9	Dodecanal	0.30	184	CCCCCCCCCCCC=O	8194
10	Tetradecanal	2.79	212	CCCCCCCCCCCCCCC=O	31291
11	1-Octadecene	1.62	252	CCCCCCCCCCCCCCCC=C	8217
12	Hexadecane	0.38	226	CCCCCCCCCCCCCCC	11006
13	Tetradecanal	0.87	192	CCCCCCCCCCCCCCC=O	31291
14	1-Tetradecanol	0.46	212	CCCCCCCCCCCCCO	8209
15	Heneicosane	8.25	214	CCCCCCCCCCCCCCCCCCC	12403
16	Octadecanal	3.42	268	CCCCCCCCCCCCCCCCCCC=O	12533
17	Phenol, 4-(2,2,3,3-Tetramethylbutyl)-	0.29	206	CC(C)(C)C(C)(C)CC1=CC=C(C=C1)O	41234
18	Tetradecanoic Acid	1.06	228	CCCCCCCCCCCCC(=O)O	11005
19	9-Octadecenal, (Z)-	0.76	226	CCCCCCCC=CCCCCCCC=O	5364492
20	5-Octadecene, (E)-	0.33	252	CCCCCCCCCCCC=CCCCC	5364598

21	1-Octadecene	1.47	252	CCCCCCCCCC=CCCCCCCC=	8217
22	Heneicosane	0.87	214	CCCCCCCCCC=CCCCCCCC=	12403
23	13-Methyl-13-Tetradecene-1,12-Diol	0.33	242	CCCCCCCCCC(=O)O	13849
24	Cis-1-Chloro-9-Octadecene	0.40	286	CCCCCCCC=CCCCCCCCCI	5367784
25	15-Isopropenylloxacyclopentadecan-2-One	0.30	266	CC(=C)C1CCCCCCCCCCC(=O)O1	557450
26	2,6,10-Trimethyl,14-Ethylene-14-Pentadecene	0.24	278	CCCCCCCC#CCCCCCCC	557019
27	2-Pentadecanone, 6,10,14-Trimethyl-	0.47	268	CC(C)CCCC(C)CCCC(C)CCCC(=O)C	10408
28	3,7,11,15-Tetramethyl-2-Hexadecen-1-OI	0.23	296	CC(C)CCCC(C)CCCC(C)CCCC(=C CO)C	5366244
29	E,E,Z-1,3,12-Nonadecatriene-5,14-Diol	2.30	242	CCCCCCCCCC(=O)O	13849
30	1-Octadecanol	0.64	270	CCCCCCCCCCCCCCCCCO	8221
31	E,E,Z-1,3,12-Nonadecatriene-5,14-Diol	2.30	294	CCCCC(C=CCCCCCCC(C=CC=C)O)O	5364768
32	1,2-Tetradecanediol	1.49	230	CCCCCCCCCCCC(CO)O	89436
33	Cis-1-Chloro-9-Octadecene	4.28	286	CCCCCCCC=CCCCCCCCCI	5367784
34	Cis-9-Hexadecenal	4.24	238	CCCCCCC=CCCCCCCC=O	5364643
35	E,E,Z-1,3,12-Nonadecatriene-5,14-Diol	1.25	294	CCCCC(C=CCCCCCCC(C=CC=C)O)O	5364768
36	N-Hexadecanoicacid	10.0 7	256	CCCCCCCCCCCCCCCC(=O)O	985
37	1-Nonadecene	1.83	266	CCCCCCCCCCCCCCCC=	29075
38	1H-3a,7-Methanoazulene, Octahydro-1,4,9,9-Tetramethyl-	0.86	206	CC1CCC23C1CC(C2(C)C)CCC3C	29408
39	Naphthalene, Decahydro-1,8a-Dimethyl-7-(1-Methylethyl)-, [1R-(1.Alpha.,4a.Beta.,7.Beta.	1.00	208	CC1CCCC2C1(CC(CC2)C(C)C)C	27255

40	<i>Z</i> -2-Tridecen-1-Ol	0.54	198	CCCCCCCCCC=CCO	5364998
41	1-Hepten-1-Ol, Acetate	0.36	156	CCCCC=COC(=O)C	5363146
42	Cyclopentadecanone, 2-Hydroxy-	0.66	240	C1CCCCCCC(C(=O)CCCCC1)O	543400
43	Heptadecanoic Acid	0.24	270	CCCCCCCCCCCCCCCCC(=O)O	10465
44	5.Alpha.-Lanostane-7,11-Dione, 3.Beta.-Hydroxy-, Acetate	0.30	500	CC(C)CCCC(C)C1CCC2(C1(CC(=O)C3C2C(=O)CC4C3(CCC(C4(C)C)OC(=O)C)C)C	242326
45	Hexadecanal	1.08	240	CCCCCCCCCCCCCCCC=O	984
46	9-Octadecenoic Acid, (<i>E</i>)-	2.56	282	CCCCCCCC=CCCCCCCC(=O)O	637517
47	9-Octadecenoic Acid (<i>Z</i>)-	0.70	282	CCCCCCCC=CCCCCCCC(=O)O	445639
48	Hexadecanamide	0.33	255	CCCCCCCCCCCCCCCC(=O)N	69421
49	1-Tetracosanol	0.74	354	CCCCCCCCCCCCCCCCCCCCCCCCCO	10472
50	<i>Z</i> -(13,14-Epoxy)Tetradec-11-En-1-Ol Acetate	0.49	268	CC(=O)OCCCCCCCCCCC=CC1CO1	5363633
51	9-Octadecenamide, (<i>Z</i>)-	0.90	281	CCCCCCCC=CCCCCCCC(=O)N	5283387
52	1-Triacontanol	0.53	438	CCCCCCCCCCCCCCCCCCCCCCCCCCCCCO	68972
53	1,2-Benzenedicarboxylic Acid	0.30	390	C1=CC=C(C(=C1)C(=O)O)C(=O)O	1017
54	1-Triacontanol	0.35	438	CCCCCCCCCCCCCCCCCCCCCCCCCCCCCO	68972
55	1-Triacontanol	0.23	438	CCCCCCCCCCCCCCCCCCCCCCCCCCCCCO	68972

56	<i>2,6,10,14,18,22-Tetracosahexaene, 2,6,10,15,19,23-Hexamethyl-, (All-E)-</i>	0.25	410	CC(=CCCC(=CCCC(=CCCC=C(C)CCC=C(C)C)C)C)	638072
57	<i>5a,8a-Epidioxycholesta-6,9(11)-Dien-3b-Ol</i>	0.35	414	CC1CCC2(C(C3C(O2)CC4C3(CCC5C4CC=C6C5(CCC(C6)O)C)C)O)C1	99474
58	<i>15-Methyl-Z-11-Hexadecenal</i>	0.46	252	CC(C)CCC=CCCCCC=O	5369022
59	<i>9-Octadecenoic Acid, 1,2,3-Propanetriyl Ester, (E,E,E)-</i>	0.97	884	CCCCCC=CCCCCC(=O)OCC(COC(=O)CCCCCC=CCC)OC(=O)CCCCCC=CCCCCCC	5364673
60	<i>Cholest-5-En-3-Ol (3.Beta.)-</i>	1.51	386	CC(C)CCCC(C)C1CCC2C1(CCC3C2CC=C4C3(CCC(C4)O)C)C	5997

Lampiran 4. Nilai *binding affinity* senyawa rumput laut *Caulerpa racemosa*

	Senyawa	Binding affinity
1	Methane, (Methylsulfinyl)(Methylthio)-	-3,1
2	2-Butanon, 3-Hydroxy	-3,9
3	2,2-Dimethoxybutane	-3,8
4	5-[(Propylsulfanyl)Methyl]-2,4-Imidazolidinedione #	-5,1
5	Hexanal	-3,7
6	(Z)-3-Chloropropenamide	-3,4
7	1,1-Diethoxy-2-Ethylhexane	-4,9
8	2-Decenal, (E)-	-4,6
9	Tetradecane	-4,5
10	1s-Neomenthyl Acetate	-5,9
11	Nonanoic Acid, 9-Oxo-, Methyl Ester	-4,9
12	3-Buten-2-One, 4-(2,6,6-Trimethyl-2-Cyclohexen-1-Yl)-	-6,3
13	R-(+)-Methyl-3-Isopropyl-6-Oxoheptanoate	-5,7
14	4-(2,6,6-Trimethyl-2-Cyclohexen-1-Yl)-2-Pentanone	-5,5
15	Phenol, 3,5-Bis(1,1-Dimethylethyl)-	-6
16	4-(2,2,6-Trimethyl-7-Oxa-Bicyclo[4.1.0]Hept-1-Yl)-But-3-En-2-One	-5,9
17	Phenol, 2,4-Bis(1,1-Dimethylethyl)-	-6,1
18	Dodecanoic Acid, Methyl Ester	-5
19	2(4h)-Benzofuranone, 5,6,7,7a-Tetrahydro-4,4,7a-Trimethyl-	-6,8
20	1,8(2H,5H)-Naphthalenedione, Hexahydro-8a-Methyl-, Cis	-6,2
21	Octanoic Acid, 3-Pentadecyl Ester	-5,7
22	Hexadecane	-4,5

23	<i>7-Hexadecenoic Acid, Methyl Ester, (Z)-</i>	-5,4
24	<i>Heneicosane</i>	-5,2
25	<i>4-(1,5-Dihydroxy-2,6,6-Trimethyl-2-Cyclohexen-1-Yl)-3-Buten-2-One</i>	-6,4
26	<i>Tridecanol, 2-Ethyl-2-Methyl-</i>	-5,2
27	<i>As-Indacene, Dodecahydro-4-(1-Octylnonyl)-</i>	-7,2
28	<i>Heptadecane</i>	-4,9
29	<i>4-(4-Hydroxy-2,2,6-Trimethyl-7-Oxa-Bicyclo[4.1.0]Hept-1-Yl)-But-3-En-2-One</i>	-7
30	<i>Tetradecanoic Acid, Methyl Ester</i>	-5,3
31	<i>Hexadecane, 2,6,10,14-Tetramethyl-</i>	-5,6
32	<i>Hexadecane, 2-Methyl-</i>	-5
33	<i>Pentadecane, 8-Hexyl-</i>	-5,4
34	<i>Heptadecane, 2-Methyl-</i>	-5,1
35	<i>3-Methylheptadecane</i>	-5,1
36	<i>2(4h)-Benzofuranone, 5,6,7,7a-Tetrahydro-6-Hydroxy-4,4,7a-Trimethyl-, (6s-Cis)-</i>	-7,2
37	<i>1-Octadecanethiol</i>	-4,7
38	<i>Octadecane</i>	-4,8
39	<i>Pentadecanoic Acid, Methyl Ester</i>	-5,2
40	<i>2,6,10-Trimethyl,14-Ethylene-14-Pentadecne</i>	-5,6
41	<i>2-Pentadecanone, 6,10,14-Trimethyl-</i>	-5,6
42	<i>1-Butyl 2-(8-Methylnonyl) Phthalate #</i>	-7
43	<i>9-Hexadecenoic Acid, Methyl Ester, (Z)-</i>	-5,4
44	<i>Hexadecanoic Acid, Methyl Ester</i>	-5,2
45	<i>3,7,11,15-Tetramethylhexadec-1-En-3-Ol</i>	-5,8
46	<i>2,3-Dimethyl-1-Undecen-3-Ol</i>	-4,8
47	<i>1,2-Benzenedicarboxylic Acid, Butyl 8-Methylnonyl Ester</i>	-6,9

48	<i>Hexadecanoic Acid, Ethyl Ester</i>	-5,4
49	<i>Cis-9-Hexadecenal</i>	-5
50	<i>Heneicosanoic Acid, Methyl Ester</i>	-5,7
51	<i>Methyl 2-Ethylhexyl Phthalate</i>	-6,9
52	<i>2-Norpinanol, 3,6,6-Trimethyl-</i>	-5,8
53	<i>8,9,12-Octadecadienoic Acid (Z,Z)-, Methyl Ester</i>	-5,9
54	<i>Palmitaldehyde, Diallyl Acetal</i>	-5,7
55	<i>Ethyl (9z,12z)-9,12-Octadecadienoate #</i>	-5,7
56	<i>10-Nonadecenoic Acid, Methyl Ester</i>	-5,6
57	<i>6,9,12-Octadecatrienoic Acid, Methyl Ester</i>	-5,9
58	<i>Linolensaeuremethyleneester</i>	-6,1
59	<i>9-Octadecynoic Acid, Methyl Ester</i>	-5,8
60	<i>4,8,12,16-Tetramethylheptadecan-4-Olide</i>	-6,3
61	<i>Methyl Eicosa-5,8,11,14,17-Pentaenoate</i>	-6,6
62	<i>7,9-Dimethyl-8-Nitrobicyclo[4.3.1]Decan-10-One</i>	-6,9
63	<i>1,2-Benzenedicarboxylic Acid, Ditridecyl Ester</i>	-6,8
64	<i>Docosanoic Acid, Methyl Ester</i>	-5,7
65	<i>1,2-Benzenedicarboxylic Acid</i>	-7,3
66	<i>Ethyl Docosanoate</i>	-5,6
67	<i>Tricosanoic Acid, Methyl Ester</i>	-6
68	<i>Dotriacontane</i>	-5,8
69	<i>Tetracosanoic Acid, Methyl Ester</i>	-5,8
70	<i>N-[1-(1-Adamantan-1-Yl-Propyl)-2,5-Dioxo-4-Trifluoromethyl-Imidazolidin-4-Yl]-4-Methoxy-Benzamide</i>	-3,7
71	<i>17-Pentatriacontene</i>	-5,9
72	<i>2,6,10,14,18,22-Tetracosahexaene, 2,6,10,15,19,23-Hexamethyl-, (All-E)-</i>	-7,4

73	<i>1,6,10-Dodecatrien-3-Ol, 3,7,11-Trimethyl-, [S-(Z)]-</i>	-6
74	<i>2H-1-Benzopyran-6-Ol, 3,4-Dihydro-2,8-Dimethyl-2-(4,8,12-Trimethyltridecyl)-, [2R-[2R@(4R@,8R@)]]-</i>	-7,6
75	<i>Hexatriacontane</i>	-5,8
76	<i>1-(4-Ethoxy-2-Hydroxyphenyl)-1-Octadecanone #</i>	-6,6
77	<i>.Beta.-Tocopherol</i>	-7,1
78	<i>Tetracontane</i>	-5,3
79	<i>Cholesta-4,6-Dien-3-Ol, Benzoate, (3.Beta.)-</i>	-4,8
80	<i>Stigmast-5-En-3-Ol, Oleate</i>	-4,3
81	<i>.Alpha.-Tocopherol-.Beta.-D-Mannoside</i>	-6,8
82	<i>Stigmastera-5,23-Dien-3-Ol, (3.Beta.)-</i>	-6,3
83	<i>Stigmast-5-En-3-Ol, (3.Beta.)-</i>	-3,3
84	<i>Lupeol</i>	4,4

Lampiran 5. Nilai *binding affinity* senyawa rumput laut *Sargassum polycystum*

	Senyawa	Binding affinity
1	1,3,5-Triazine-2,4-Diamine, 6-Chloro-N-Ethyl-	-6,0
2	1,6-Octadien-3-Ol, 3,7-Dimethyl-	-5,1
3	1-Decanol	-6,0
4	Undecane	-4,0
5	N-Formylmorpholine	-4,2
6	1,3-Dioxane, 4,6-Dimethyl-	-4,3
7	Methenamine	-4,5
8	Azulene, 1,2,3,5,6,7,8,8a-Octahydro-1,4-Dimethyl-7-(1-Methylethenyl)-, [1S-(1.Alpha.,7.A)]	-6,3
9	2(4h)-Benzofuranone, 5,6,7,7a-Tetrahydro-6-Hydroxy-4,4,7a-Trimethyl	-6,8
10	2,6,10-Trimethyl,14-Ethylene-14-Pentadecne	-5,6
11	Hexadecanoic Acid, Methyl Ester	-5,2
12	Hexadecanoic Acid, Ethyl Ester	-5,4
13	9-Octadecenoic Acid (Z)-, Methyl Ester	-5,8
14	6,8-Dioxabicyclo(3.2.1)Octan-4.Beta.-Ol	-4,9
15	Octadecanoic Acid, Methyl Ester	-5,5
16	9,19-Cycloergost-24(28)-En-3-Ol, 4,14-Dimethyl-, Acetate, (3.Beta.,4.Alpha.,5.Alpha.)-	-1,5
17	Dotriacontane	-5,8
18	Heneicosane, 11-Cyclopentyl-	-6,0
19	1,2-Benzenedicarboxylic Acid	-7,3
20	Triacontane	-5,7
21	Tetrapentacontane	-5,3
22	1-Hentetracontanol	-6,0

23	<i>Cyclohexane, Nonadecyl-</i>	-6,1
24	<i>Hexacontane</i>	-4,5
25	<i>Tetracontane</i>	-5,7
26	<i>Cholest-5-En-3-Ol (3.Beta.)-</i>	-7,0

Lampiran 6. Nilai *binding affinity* senyawa rumput laut *Halymenia durvillei*

	Senyawa	Binding affinity
1	Benzene, Ethyl-	-4,8
2	Benzene, 1,2-Dimethyl-	-4,6
3	Benzene, Ethenyl-	-4,8
4	Ethanol, 2-Butoxy-	-4,2
5	2-Butoxyethyl Acetate	-4,5
6	1-Dodecanol	-5,0
7	1-Tetradecanol	-4,8
8	Tetradecane	-4,5
9	Dodecanal	-4,4
10	Tetradecanal	-4,8
11	1-Octadecene	-5,0
12	Hexadecane	-4,5
13	Heneicosane	-5,2
14	Octadecanal	-4,9
15	Phenol, 4-(2,2,3,3-Tetramethylbutyl)-	-6,2
16	Tetradecanoic Acid	-5,2
17	9-Octadecenal, (Z)-	-5,3
18	5-Octadecene, (E)-	-5,2
19	13-Methyl-13-Tetradecene-1,12-Diol	-5,5
20	Cis-1-Chloro-9-Octadecene	-5,2
21	15-Isopropenyloxacyclopentadecan-2-One	-6,5
22	2,6,10-Trimethyl,14-Ethylene-14-Pentadecene	-5,6
23	2-Pentadecanone, 6,10,14-Trimethyl-	-5,6
24	3,7,11,15-Tetramethyl-2-Hexadecen-1-Ol	-6,3
25	E,E,Z-1,3,12-Nonadecatriene-5,14-Diol	-5,9
26	1-Octadecanol	-5,1
27	1,2-Tetradecanediol	-5,1

28	<i>Cis</i> -9-Hexadecenal	-5
29	<i>N</i> -Hexadecanoic acid	-5,4
30	1-Nonadecene	-4,9
31	1 <i>H</i> -3 <i>a</i> ,7-Methanoazulene, Octahydro-1,4,9,9-Tetramethyl-	-6,4
32	Naphthalene, Decahydro-1,8 <i>a</i> -Dimethyl-7-(1-Methylethyl)-, [<i>1R</i> -(1.Alpha.,4 <i>a</i> .Beta.,7.Beta.	-6,2
33	<i>Z</i> -2-Tridecen-1-Ol	-5,2
34	1-Hepten-1-Ol, Acetate	-4,6
35	Cyclopentadecanone, 2-Hydroxy-	-6,4
36	Heptadecanoic Acid	-5,5
37	5.Alpha.-Lanostane-7,11-Dione, 3.Beta.-Hydroxy-, Acetate	-0,8
38	Hexadecanal	-4,8
39	9-Octadecenoic Acid, (<i>E</i>)-	-5,8
40	9-Octadecenoic Acid (<i>Z</i>)-	-5,8
41	Hexadecanamide	-5,6
42	1-Tetracosanol	-5,6
43	<i>Z</i> -(13,14-Epoxy)Tetradec-11-En-1-Ol Acetate	-5,6
44	9-Octadecenamide, (<i>Z</i>)-	-5,8
45	1-Triacontanol	-5,9
46	1,2-Benzenedicarboxylic Acid	-7,3
47	2,6,10,14,18,22-Tetracosahexaene, 2,6,10,15,19,23-Hexamethyl-, (<i>All-E</i>)-	-7,4
48	5 <i>a</i> ,8 <i>a</i> -Epidioxycholest-6,9(11)-Dien-3 <i>b</i> -Ol	-2,1
49	15-Methyl- <i>Z</i> -11-Hexadecenal	-5,2
50	9-Octadecenoic Acid, 1,2,3-Propanetriyl Ester, (<i>E,E,E</i>)-	-6,8
51	Cholest-5-En-3-Ol (3.Beta.)-	-7,0