

Supplementary Information

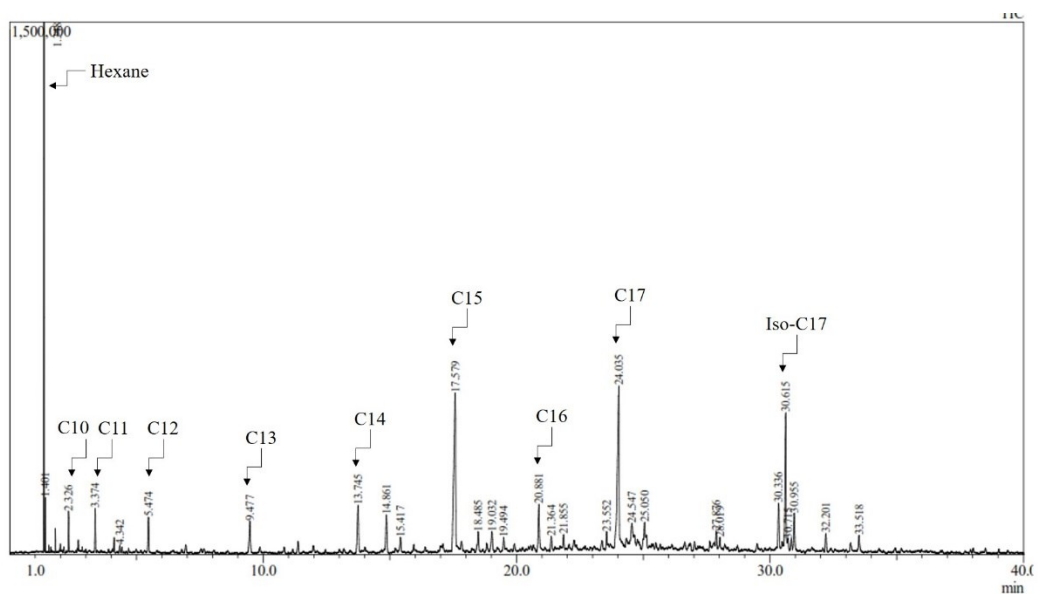


Fig. S1 GC-MS spectra of deoxygenated liquid product using 10%Ni/Al-MCM-41 catalyst

Table S1. Composition of liquid products of *Reutealis trisperma* oil deoxygenation using Ni/Al-MCM-41 by GC-MS analysis

Retention Time	Compound name	Selectivity (%)				
		Al-MCM-41	5%Ni/Al-MCM-41	10%Ni/Al-MCM-41	15%Ni/Al-MCM-41	20%Ni/Al-MCM-41
Alkane						
1.539	Octane	0.57	0.11	-	-	0.47
1.801	Nonane	1.98	0.27	-	0.25	1.61
2.326	Decane	3.48	0.50	1.17	0.48	2.31
3.388	Undecane	4.74	0.83	1.90	0.74	2.61
5.512	Dodecane	6.26	0.87	2.42	1.02	2.30
9.546	Tridecane	10.72	3.93	4.21	2.1	4.37
13.802	Tetradecane	8.72	5.36	6.13	4.99	7.45
17.589	Pentadecane	17.03	32.09	20.95	19.85	39.52
20.895	Hexadecane	0.90	5.83	5.78	6.58	-
23.979	Heptadecane	2.81	7.64	21.34	19.98	-
26.802	Octadecane	-	0.82	-	3.00	-
30.615	Iso-heptadecane	1.42	11.85	20.53	18.71	4.87
Alkene						
1.639	1-Octene	-	0.26	-	-	0.21
1.994	1-Nonene	1.01	0.45	-	-	0.66
2.711	1-Decene	1.71	0.64	-	-	1.43
4.141	1-Dodecene	4.29	2.86	-	0.69	2.00
4.45	Undecene	-	0.63	-	-	-
10.64	Tetradecene	-	1.14	-	-	-
11.421	1-Tridecene	1.93	1.72	-	0.50	0.52
18.814	3-Octadecene	-	-	-	0.67	-
19.031	1-Hexadecene	-	0.96	1.22	-	-
19.494	1-Pentadecene	-	0.35	0.85	0.94	-
24.769	1-Heptadecene	-	2.21	2.93	6.75	-
Cyclic						
1.396	Methylcyclopentane	0.40	0.64	1.06	0.58	3.66
2.245	Propylcyclohexane	-	-	-	-	0.43
4.447	Cyclopropane, 1,2-dibutyl	0.52	-	-	-	2.53
13.036	Cyclohexane, (4-methylpentyl)	-	0.33	-	-	-

Retention Time	Compound name	Selectivity (%)				
		Al-MCM-41	5%Ni/Al-MCM-41	10%Ni/Al-MCM-41	15%Ni/Al-MCM-41	20%Ni/Al-MCM-41
20.655	Undecylcyclohexane	-	0.20	-	0.40	-
23.363	Cyclodecane-octyl	-	-	-	0.85	-
23.689	Cyclohexane, 1,2-dimethyl-3-pentyl-4-propyl-	-	-	-	0.25	-
24.335	Cyclohexane, hexaethyl	-	-	-	1.05	-
25.398	Cyclooctane,tetradecanyl	-	-	-	0.44	-
25.492	Cyclohexane, 1,2-dimethyl-3-pentyl-4-propyl-	-	-	-	0.54	-
27.014	Undecylcyclohexane	-	-	-	0.60	-
Aromatics						
5.192	o-xylene	0.66	0.23	-	-	-
9.919	n-butylbenzene	1.09	0.81	-	-	0.72
10.871	2-propyltoluene	1.17	0.86	-	-	0.81
14.058	Pentylbenzene	0.78	0.45	-	-	-
14.846	Benzene, (1,2-dimethylpropyl)	-	-	-	-	6.18
14.902	2-Phenylpentane	8.23	2.95	3.12	1.46	-
17.066	Benzene, 1-ethyl-4-(2-methylpropyl)-	-	0.65	-	0.57	-
17.133	p-cymene	1.95	-	-	-	-
17.863	1-Phenylhexane	2.46	1.39	-	0.34	3.12
18.475	Benzene, (1,3-dimethylbutyl)	-	-	-	-	1.85
18.523	2-Methyl-4-phenylpentane	3.36	2.50	1.25	1.36	-
21.884	1-Methyl-2-N-hexylbenzene	1.52	-	0.98	0.61	0.85
26.628	Benzene, 1,1'-(1,1,3,3-tetramethyl-1,3-propanediyl)bis-	-	-	-	0.54	-
28.019	Hexan,2-phenyl-3-propyl	-	-	0.74	0.81	0.94
30.715	2-Naphthalenamine, 1,2,3,4-tetrahydro-N-(1-methylethyl)	-	-	0.93	-	-
32.201	Benzene, 1,1'-(1,1,3,3-tetramethyl-1,3-propanediyl)bis-	-	-	1.62	-	-

Retention Time	Compound name	Selectivity (%)				
		Al-MCM-41	5%Ni/Al-MCM-41	10%Ni/Al-MCM-41	15%Ni/Al-MCM-41	20%Ni/Al-MCM-41
Carboxylic acid						
24.15	Benzenepentanoic acid	-	-	-	0.26	-
25.075	Butanoic acid	3.06	-	-	-	-
28.051	Hexanoic acid	2.33	1.51	-	-	-
30.846	Heptanoic acid	2.38	1.64	-	-	1.28
33.492	Octanoic acid	1.50	1.56	-	-	1.66
35.97	Nonanoic acid	-	0.57	-	-	1.21
Alcohol						
14.81	1-Dodecanol	-	0.59	-	-	-
15.403	1-Tetradecanol	-	-	-	-	0.58
15.417	1-Tridecanol	-	0.08	0.84	0.56	-
19.014	Pentadecanol	-	0.31	-	1.05	-
25.038	2-phenyl-1-propanol	-	-	-	-	3.82
Ketone						
18.465	Hydroxybenzophenon	-	0.57	-	-	-
25.16	2,2-dimethyl-5-phenyl-hexone	0.99	-	-	-	-
30.7	5,9,9-Trimethylspiro[3.6]deca-5,7-dien-1-one	-	-	-	0.45	-
Aldehyde						
18.364	9-Octadecenal	-	0.81	-	-	-