

SUPPLEMENTARY MATERIAL

Table S1: Experimental Refractive Indices, n_D , and Refractive Index Deviation, Δn_D , of the Binary Mixtures Isomeric Chlorobutane (1) + Butyl Ethyl Ether (2) at Several Temperatures

ϕ_1	n_D	Δn_D	ϕ_1	n_D	Δn_D
$T = 283.15 \text{ K}$					
1-Chlorobutane (1) + butyl ethyl ether (2)			2-Chlorobutane (1) + butyl ethyl ether (2)		
0.0400	1.388076	0.000042	0.0413	1.387878	0.000034
0.0798	1.388927	0.000086	0.0775	1.388451	0.000063
0.1616	1.390669	0.000169	0.1629	1.389796	0.000124
0.2476	1.392475	0.000230	0.2525	1.391181	0.000165
0.3385	1.394356	0.000268	0.3416	1.392558	0.000199
0.4354	1.396341	0.000288	0.4622	1.394388	0.000216
0.5373	1.398416	0.000296	0.5402	1.395567	0.000218
0.6418	1.400527	0.000288	0.6455	1.397140	0.000213
0.7546	1.402789	0.000262	0.7560	1.398785	0.000197
0.8742	1.405134	0.000181	0.8747	1.400518	0.000145
0.9360	1.406328	0.000122	0.9369	1.401405	0.000091
1-Chloro-2-methylpropane(1) + butyl ethyl ether (2)			2-Chloro-2-methylpropane(1) + butyl ethyl ether (2)		
0.0386	1.387881	0.000028	0.0417	1.387415	0.000028
0.0795	1.388582	0.000062	0.0851	1.387640	0.000083
0.1623	1.390000	0.000128	0.1687	1.388016	0.000130
0.2502	1.391481	0.000175	0.2563	1.388380	0.000150
0.3424	1.393016	0.000205	0.3505	1.388738	0.000138
0.4388	1.394608	0.000224	0.4440	1.389077	0.000109
0.5381	1.396231	0.000227	0.5466	1.389445	0.000074
0.6429	1.397927	0.000213	0.6517	1.389826	0.000042
0.7552	1.399725	0.000178	0.7628	1.390239	0.000018
0.8736	1.401584	0.000105	0.8783	1.390684	0.000009
0.9352	1.402544	0.000059	0.9376	1.390914	0.000006
$T = 288.15 \text{ K}$					
1-Chlorobutane (1) + butyl ethyl ether (2)			2-Chlorobutane (1) + butyl ethyl ether (2)		
0.0400	1.385478	0.000033	0.0413	1.385286	0.000033
0.0798	1.386328	0.000080	0.0775	1.385848	0.000056
0.1616	1.388055	0.000159	0.1629	1.387182	0.000116
0.2476	1.389855	0.000222	0.2525	1.388556	0.000157
0.3385	1.391725	0.000260	0.3416	1.389918	0.000192
0.4354	1.393699	0.000277	0.4622	1.391730	0.000208
0.5373	1.395764	0.000286	0.5402	1.392899	0.000213
0.6418	1.397864	0.000278	0.6455	1.394460	0.000207
0.7546	1.400113	0.000251	0.7560	1.396087	0.000187
0.8742	1.402451	0.000176	0.8747	1.397807	0.000140
0.9360	1.403634	0.000112	0.9369	1.398683	0.000090

(Table S1). Continued.

ϕ_1	n_D	Δn_D	ϕ_1	n_D	Δn_D
1-Chloro-2-methylpropane(1) + butyl ethyl ether (2)			2-Chloro-2-methylpropane(1) + butyl ethyl ether (2)		
0.0386	1.385346	0.000084	0.0417	1.384831	0.000044
0.0795	1.386103	0.000180	0.0851	1.385038	0.000096
0.1623	1.387649	0.000388	0.1687	1.385391	0.000150
0.2502	1.389206	0.000525	0.2563	1.385723	0.000169
0.3424	1.390821	0.000650	0.3505	1.386042	0.000152
0.4388	1.392410	0.000681	0.4440	1.386343	0.000119
0.5380	1.393975	0.000642	0.5466	1.386670	0.000079
0.6429	1.395550	0.000523	0.6517	1.387011	0.000045
0.7552	1.397195	0.000354	0.7628	1.387384	0.000021
0.8736	1.398918	0.000163	0.8783	1.387786	0.000011
0.9352	1.399814	0.000064	0.9376	1.387995	0.000008
T = 293.15 K					
1-Chlorobutane (1) + butyl ethyl ether (2)			2-Chlorobutane (1) + butyl ethyl ether (2)		
0.0400	1.382871	0.000031	0.0413	1.382676	0.000029
0.0798	1.383711	0.000071	0.0775	1.383232	0.000051
0.1616	1.385431	0.000150	0.1629	1.384552	0.000109
0.2476	1.387221	0.000211	0.2525	1.385915	0.000150
0.3385	1.389084	0.000250	0.3416	1.387266	0.000186
0.4354	1.391051	0.000271	0.4622	1.389062	0.000202
0.5373	1.393106	0.000279	0.5402	1.390221	0.000208
0.6418	1.395197	0.000271	0.6455	1.391767	0.000200
0.7546	1.397436	0.000244	0.7560	1.393380	0.000181
0.8742	1.399758	0.000164	0.8747	1.395083	0.000133
0.9360	1.400937	0.000100	0.9369	1.395953	0.000085
1-Chloro-2-methylpropane(1) + butyl ethyl ether (2)			2-Chloro-2-methylpropane(1) + butyl ethyl ether (2)		
0.0386	1.382680	0.000026	0.0417	1.382218	0.000048
0.0795	1.383367	0.000060	0.0851	1.382415	0.000107
0.1623	1.384755	0.000124	0.1687	1.382737	0.000163
0.2502	1.386205	0.000170	0.2563	1.383036	0.000183
0.3424	1.387708	0.000199	0.3505	1.383318	0.000165
0.4389	1.389269	0.000218	0.4440	1.383579	0.000129
0.5381	1.390857	0.000221	0.5466	1.383862	0.000086
0.6429	1.392517	0.000206	0.6517	1.384160	0.000049
0.7552	1.394276	0.000171	0.7628	1.384488	0.000024
0.8737	1.396100	0.000101	0.8783	1.384843	0.000012
0.9352	1.397035	0.000054	0.9376	1.385027	0.000008
T = 298.15 K					
1-Chlorobutane (1) + butyl ethyl ether (2)			2-Chlorobutane (1) + butyl ethyl ether (2)		
0.0400	1.380265	0.000031	0.0413	1.380058	0.000024
0.0798	1.381099	0.000066	0.0775	1.380610	0.000047

(Table S1). Continued.

ϕ_1	n_D	Δn_D	ϕ_1	n_D	Δn_D
0.1616	1.382817	0.000145	0.1629	1.381912	0.000100
0.2476	1.384607	0.000209	0.2525	1.383261	0.000142
0.3385	1.386467	0.000245	0.3416	1.384600	0.000180
0.4354	1.388433	0.000266	0.4622	1.386376	0.000195
0.5373	1.390484	0.000272	0.5402	1.387522	0.000200
0.6418	1.392576	0.000267	0.6455	1.389053	0.000194
0.7546	1.394821	0.000248	0.7560	1.390649	0.000175
0.8742	1.397151	0.000178	0.8747	1.392329	0.000123
0.9360	1.398327	0.000111	0.9369	1.393191	0.000077
1-Chloro-2-methylpropane(1) + butyl ethyl ether (2)			2-Chloro-2-methylpropane(1) + butyl ethyl ether (2)		
0.0386	1.380068	0.000025	0.0417	1.379601	0.000052
0.0796	1.380752	0.000059	0.0851	1.379783	0.000111
0.1623	1.382126	0.000122	0.1687	1.380081	0.000173
0.2503	1.383565	0.000166	0.2563	1.380350	0.000194
0.3425	1.385058	0.000197	0.3505	1.380598	0.000176
0.4390	1.386607	0.000216	0.4440	1.380827	0.000141
0.5382	1.388181	0.000217	0.5466	1.381069	0.000094
0.6430	1.389828	0.000203	0.6517	1.381332	0.000060
0.7553	1.391573	0.000167	0.7628	1.381618	0.000032
0.8737	1.393383	0.000100	0.8783	1.381926	0.000015
0.9352	1.394312	0.000054	0.9376	1.382087	0.000009
T = 303.15 K					
1-Chlorobutane (1) + butyl ethyl ether (2)			2-Chlorobutane (1) + butyl ethyl ether (2)		
0.0400	1.377587	0.000030	0.0413	1.377379	0.000020
0.0798	1.378411	0.000060	0.0775	1.377924	0.000040
0.1616	1.380119	0.000137	0.1629	1.379216	0.000093
0.2476	1.381897	0.000197	0.2525	1.380552	0.000133
0.3385	1.383748	0.000236	0.3416	1.381883	0.000173
0.4354	1.385704	0.000259	0.4622	1.383648	0.000191
0.5373	1.387746	0.000267	0.5402	1.384783	0.000194
0.6418	1.389826	0.000262	0.6455	1.386300	0.000187
0.7546	1.392053	0.000236	0.7560	1.387883	0.000168
0.8742	1.394364	0.000160	0.8747	1.389550	0.000116
0.9360	1.395549	0.000110	0.9369	1.390407	0.000072
1-Chloro-2-methylpropane(1) + butyl ethyl ether (2)			2-Chloro-2-methylpropane(1) + butyl ethyl ether (2)		
0.0386	1.377391	0.000024	0.0417	1.376921	0.000058
0.0795	1.378066	0.000058	0.0851	1.377094	0.000124
0.1623	1.379428	0.000121	0.1687	1.377363	0.000188
0.2502	1.380852	0.000165	0.2563	1.377600	0.000210
0.3424	1.382327	0.000194	0.3505	1.377812	0.000191
0.4389	1.383860	0.000213	0.4440	1.378009	0.000158

(Table S1). Continued.

ϕ_1	n_D	Δn_D	ϕ_1	n_D	Δn_D
0.5381	1.385418	0.000214	0.5466	1.378219	0.000117
0.6430	1.387048	0.000198	0.6517	1.378435	0.000076
0.7553	1.388775	0.000163	0.7628	1.378674	0.000043
0.8737	1.390568	0.000099	0.8783	1.378934	0.000020
0.9352	1.391488	0.000054	0.9376	1.379070	0.000011
T = 308.15 K					
1-Chlorobutane (1) + butyl ethyl ether (2)			2-Chlorobutane (1) + butyl ethyl ether (2)		
0.0400	1.374883	0.000027	0.0413	1.374674	0.000017
0.0798	1.375708	0.000060	0.0775	1.375213	0.000035
0.1616	1.377407	0.000134	0.1629	1.376494	0.000087
0.2476	1.379175	0.000192	0.2525	1.377820	0.000126
0.3385	1.381020	0.000230	0.3416	1.379142	0.000166
0.4354	1.382970	0.000253	0.4622	1.380895	0.000185
0.5373	1.385004	0.000260	0.5402	1.382021	0.000188
0.6418	1.387073	0.000251	0.6455	1.383527	0.000181
0.7546	1.389284	0.000219	0.7560	1.385096	0.000160
0.8742	1.391590	0.000145	0.8747	1.386752	0.000110
0.9360	1.392770	0.000094	0.9369	1.387603	0.000066
1-Chloro-2-methylpropane(1) + butyl ethyl ether (2)			2-Chloro-2-methylpropane(1) + butyl ethyl ether (2)		
0.0386	1.374688	0.000024	0.0417	1.374222	0.000070
0.0796	1.375359	0.000057	0.0851	1.374393	0.000148
0.1623	1.376708	0.000119	0.1687	1.374637	0.000214
0.2503	1.378121	0.000163	0.2563	1.374838	0.000228
0.3425	1.379584	0.000191	0.3505	1.375020	0.000210
0.4389	1.381104	0.000210	0.4440	1.375187	0.000178
0.5381	1.382648	0.000211	0.5466	1.375370	0.000143
0.6430	1.384268	0.000198	0.6517	1.375550	0.000100
0.7553	1.385980	0.000162	0.7628	1.375749	0.000063
0.8737	1.387758	0.000098	0.8783	1.375960	0.000029
0.9352	1.388672	0.000054	0.9376	1.376071	0.000014
T = 313.15 K					
1-Chlorobutane (1) + butyl ethyl ether (2)			2-Chlorobutane (1) + butyl ethyl ether (2)		
0.0400	1.372226	0.000023	0.0413	1.372008	0.000012
0.0798	1.373046	0.000051	0.0775	1.372540	0.000030
0.1616	1.374743	0.000122	0.1629	1.373804	0.000079
0.2476	1.376509	0.000177	0.2525	1.375116	0.000119
0.3385	1.378361	0.000222	0.3416	1.376423	0.000160
0.4354	1.380314	0.000247	0.4622	1.378154	0.000178
0.5373	1.382352	0.000255	0.5402	1.379270	0.000184
0.6418	1.384421	0.000244	0.6455	1.380756	0.000175
0.7546	1.386632	0.000210	0.7560	1.382306	0.000154

(Table S1). Continued.

ϕ_1	n_D	Δn_D	ϕ_1	n_D	Δn_D
0.8742	1.388934	0.000130	0.8747	1.383940	0.000102
0.9360	1.390112	0.000077	0.9369	1.384780	0.000059
1-Chloro-2-methylpropane(1) + butyl ethyl ether (2)			2-Chloro-2-methylpropane(1) + butyl ethyl ether (2)		
0.0386	1.372027	0.000023	0.0417	1.371559	0.000079
0.0796	1.372692	0.000056	0.0851	1.371720	0.000165
0.1624	1.374029	0.000116	0.1687	1.371928	0.000230
0.2504	1.375430	0.000160	0.2563	1.372088	0.000240
0.3426	1.376878	0.000187	0.3505	1.372236	0.000227
0.4391	1.378385	0.000206	0.4440	1.372368	0.000199
0.5383	1.379916	0.000207	0.5466	1.372511	0.000168
0.6431	1.381518	0.000194	0.6517	1.372651	0.000128
0.7554	1.383213	0.000157	0.7628	1.372786	0.000074
0.8738	1.384976	0.000095	0.8783	1.372947	0.000039
0.9353	1.385880	0.000051	0.9376	1.373027	0.000018