

APPENDICES

Appendix A

Appendix B

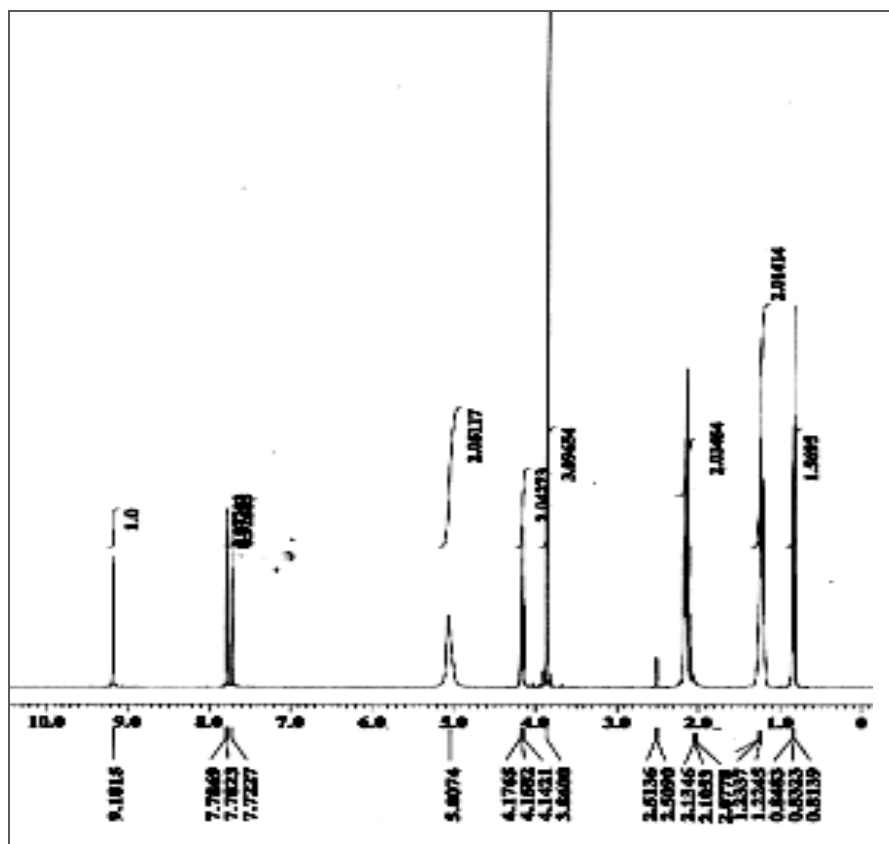


Figure B.1 Proton NMR spectrum for MSBIMHSO₄. The peaks 1-9 indicated, -N-C-CH₂-C-C-S-, -N-C-C-CH₂-C-S-, -N-C-C-C-CH₂-S-, CH₃-N-, -N-CH₂-C-C-C-S-, OH, -N-CH-C-N-, -N-C-CH-N-, and -N-CH-N-, respectively.

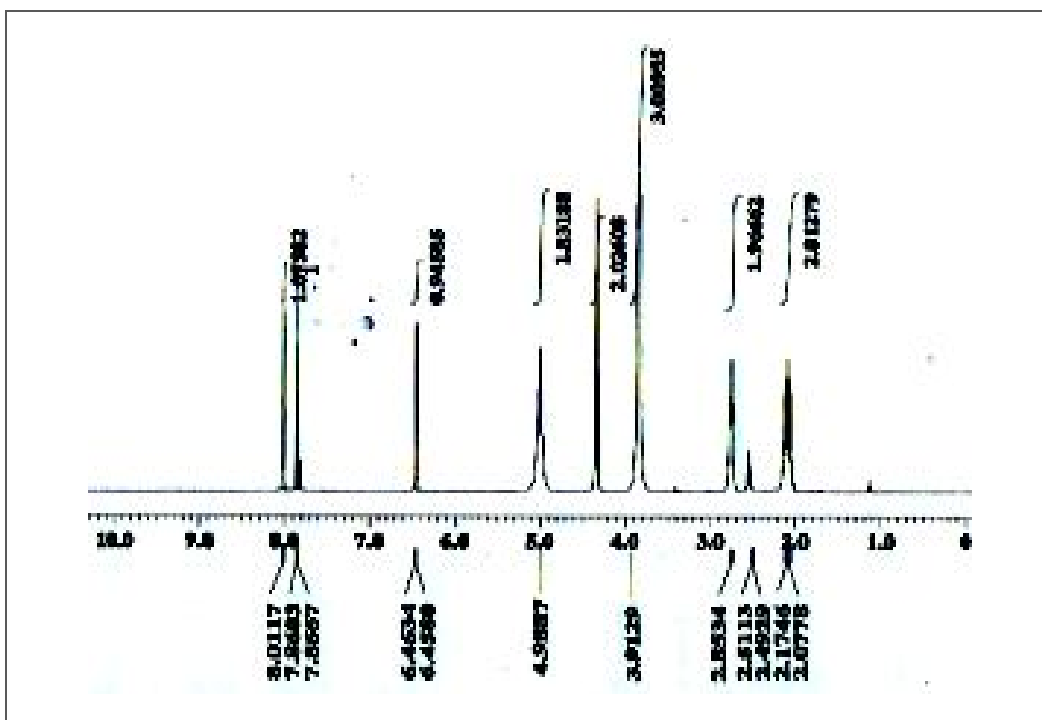


Figure B.2 Proton NMR spectrum for MSPPHSO₄. The peaks 1-8 indicated, -N-C-CH₂-C-S-, -N- CH₂-C-C-S-, CH₃-C-, -N-C-C-CH₂-S-, -OH, -N-C-CH-C-N-, -N-CH-C-N-, and -N-C-CH-N-, respectively.

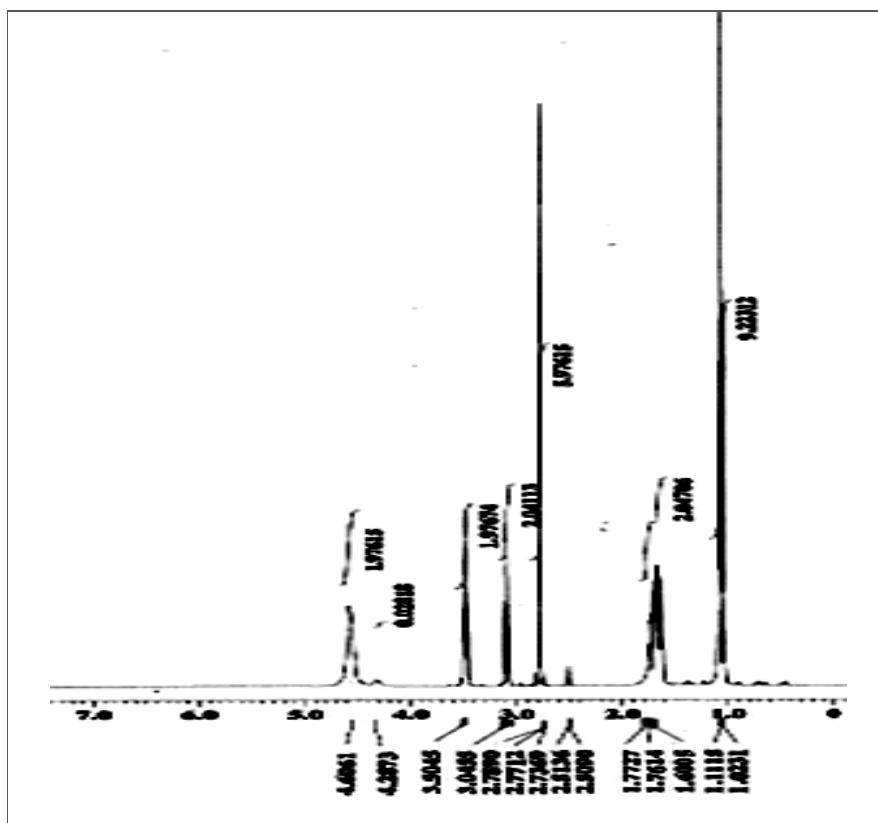


Figure B.3 Proton NMR spectrum for TESPAMHSO₄. The peaks 1-6 indicated, $\text{CH}_2\text{-C-}$, $\text{-N-C-CH}_2\text{-C-S-}$, $\text{-N-C-C-CH}_2\text{-S-}$, $\text{C-CH}_2\text{-N-}$, $\text{-N-CH}_2\text{-C-C-S-}$, and -S-OH , respectively.

Appendix C

Table C.1 Carbon (C, 12.0107 g/mol), Hydrogen (H, 1.0079 g/mol), Nitrogen (N, 14.0067 g/mol) and Sulfur (S, 32.066 g/mol) percentages in the investigated ILs

IL	C	H	N	S	Predicted
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	(%)	(%)	(%)	(%)	formula
MPSIM	43.23 ± 0.02	6.07 ± 0.018	13.37 ± 0.044	16.58 ± 0.015	C ₇ H ₁₂ N ₂ O ₃ S
MSPIMHSO ₄	27.85 ± 0.047	4.58 ± 0.053	9.29 ± 0.081	21.26 ± 0.046	C ₇ H ₁₄ N ₂ O ₇ S ₂
MBSIM	45.03 ± 0.026	6.556 ± 0.031	12.23 ± 0.029	15.44 ± 0.030	C ₈ H ₁₄ N ₂ O ₃ S
MSBIMHSO ₄	30.32 ± 0.095	5.094 ± 0.024	8.99 ± 0.0164	20.26 ± 0.073	C ₈ H ₁₆ N ₂ O ₇ S ₂
BPSIM	49.05 ± 0.03	7.551 ± 0.005	11.38 ± 0.045	13.34 ± 0.038	C ₁₀ H ₁₈ N ₂ O ₃ S
BBSIM	50.93 ± 0.038	7.772 ± 0.010	10.80 ± 0.075	12.47 ± 0.045	C ₁₁ H ₂₀ N ₂ O ₃ S
BSBIMHSO ₄	36.82 ± 0.01	6.17 ± 0.058	7.91 ± 0.086	18.11 ± 0.03	C ₁₁ H ₂₂ N ₂ O ₇ S ₂
PSP	48.71 ± 0.049	9.66 ± 0.058	5.92 ± 0.053	14.73 ± 0.026	C ₆ H ₁₀ N ₂ O ₃ S
SPPHSO ₄	24.78 ± 0.015	4.31 ± 0.004	10.03 ± 0.011	22.95 ± 0.02	C ₆ H ₁₂ N ₂ O ₇ S ₂
SBPHSO ₄	27.77 ± 0.032	4.599 ± 0.016	9.29 ± 0.033	21.24 ± 0.020	C ₇ H ₁₄ N ₂ O ₇ S ₂
MSPPHSO ₄	27.77 ± 0.064	4.62 ± 0.008	9.28 ± 0.010	21.25 ± 0.06	C ₇ H ₁₄ N ₂ O ₇ S ₂
MSBPHSO ₄	30.23 ± 0.040	5.095 ± 0.025	8.985 ± 0.022	20.26 ± 0.035	C ₈ H ₁₆ N ₂ O ₇ S ₂
DEMSPAM	48.93 ± 0.035	9.57 ± 0.020	6.24 ± 0.01	14.51 ± 0.015	C ₉ H ₂₁ NO ₃ S
DEMSBAMHSO ₄	33.90 ± 0.020	7.312 ± 0.010	4.42 ± 0.004	20.24 ± 0.072	C ₉ H ₂₃ NO ₇ S ₂
TEPSAM	49.26 ± 0.036	9.64 ± 0.020	6.30 ± 0.006	14.62 ± 0.035	C ₉ H ₂₁ NO ₃ S
TESPAMHSO ₄	33.73 ± 0.038	7.31 ± 0.005	4.42 ± 0.003	20.24 ± 0.059	C ₉ H ₂₃ NO ₇ S ₂
TEBSAM	50.97 ± 0.011	9.96 ± 0.025	6.004 ± 0.034	13.98 ± 0.086	C ₁₀ H ₂₃ NO ₃ S
TESBAMHSO ₄	36.05 ± 0.026	7.58 ± 0.005	4.24 ± 0.009	19.07 ± 0.050	C ₁₀ H ₂₅ NO ₇ S ₂

Appendix D

Table D.1 The thermal decomposition of investigated zwitterionic ILs and RTILs

IL	T _{dec} ± 2.5 (°C)
MPSIM	370
MSPIMHSO ₄	315

MBSIM	367.9
MSBIMHSO ₄	306
BPSIM	372.8
BSPIMHSO ₄	323
BBSIM	360.2
BSBIMHSO ₄	311.5
PSP	347.4
SPPHSO ₄	300.4
BSP	352.7
SBPHSO ₄	308.3
MPSP	359.8
MSPPHSO ₄	320
MBSP	350.8
MSBPHSO ₄	308
DEMPSAM	328
DEMSPAMHSO ₄	332.7
DEMBSAM	318
DEMSBAMHSO ₄	330
TEPSAM	296.8
TESPAMHSO ₄	295.4
TEBSAM	288.8
TESBAMHSO ₄	280

Appendix E

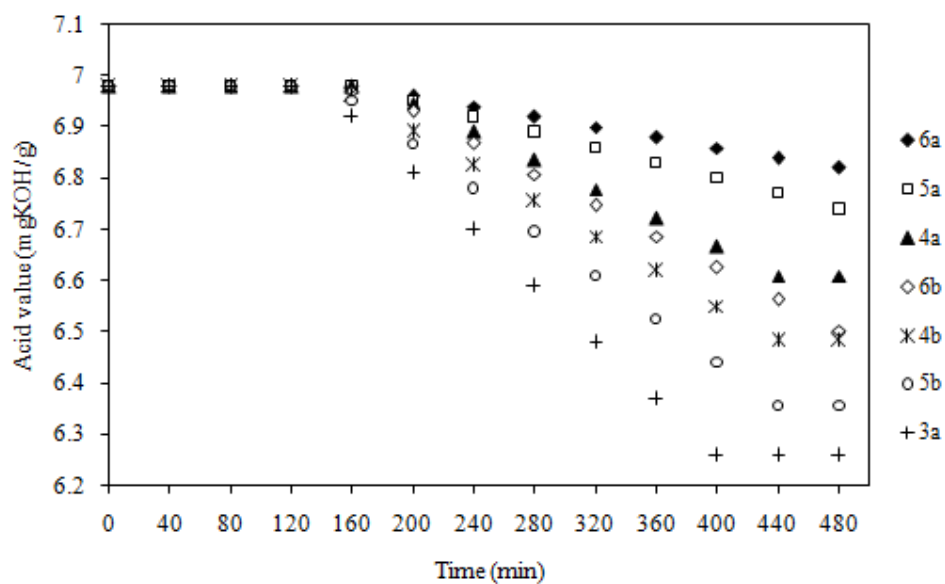


Figure E.1 The effect of different catalysts on the CPO acidity reduction during the transesterification reactions. 3a = SPPHSO₄; 4a = MSPPHSO₄; 4b = MSBPHSO₄, 5a = DEMSPAMHSO₄, 5b = DEMSBAMHSO₄, 6a = TESPAMHSO₄, 6b = TESBAMHSO₄.

Appendix F

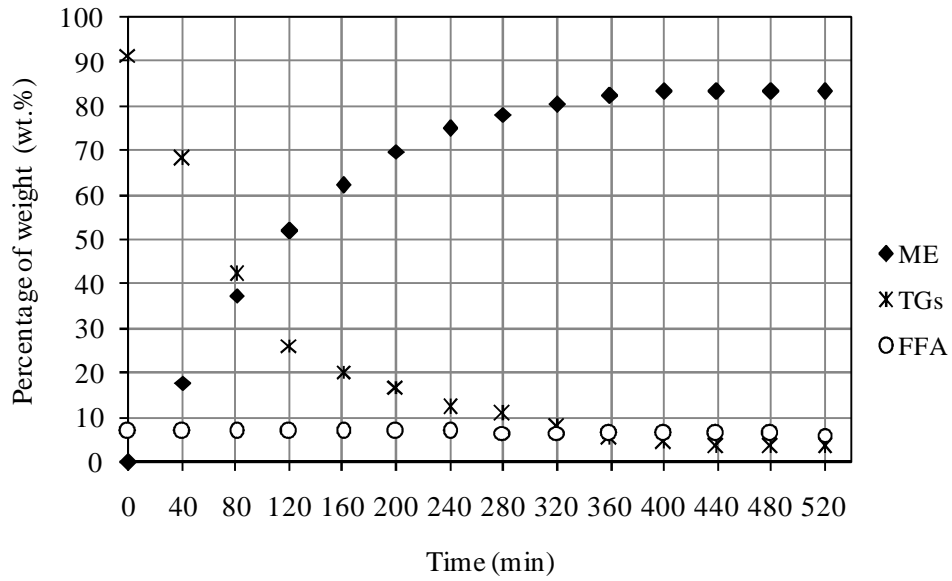


Figure F.1 The effect of MSPIMHSO₄ catalyst on the TGs and FFA conversion to ME.

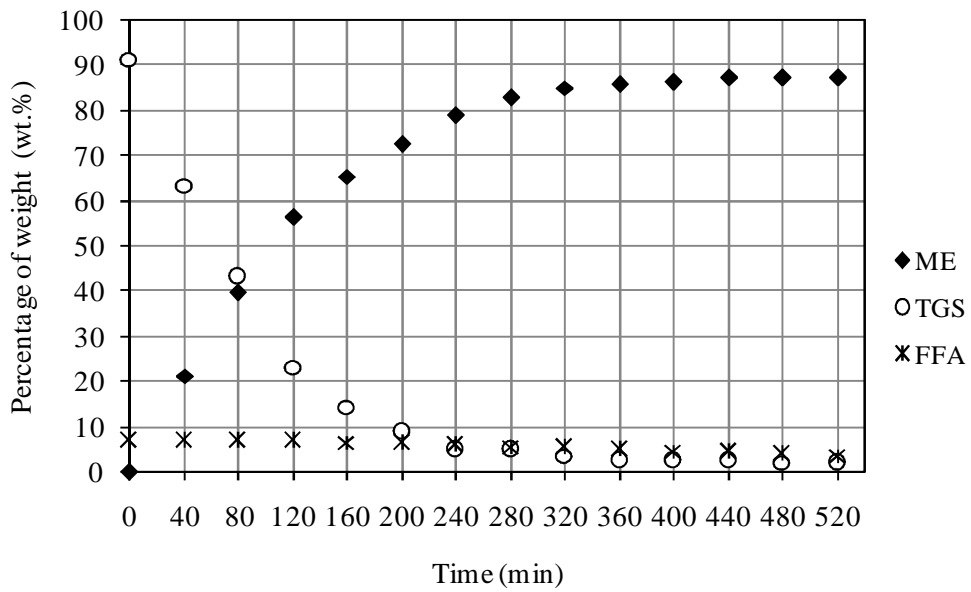


Figure F.2 The effect of MSBIMHSO₄ catalyst on the TGs and FFA conversion to ME.

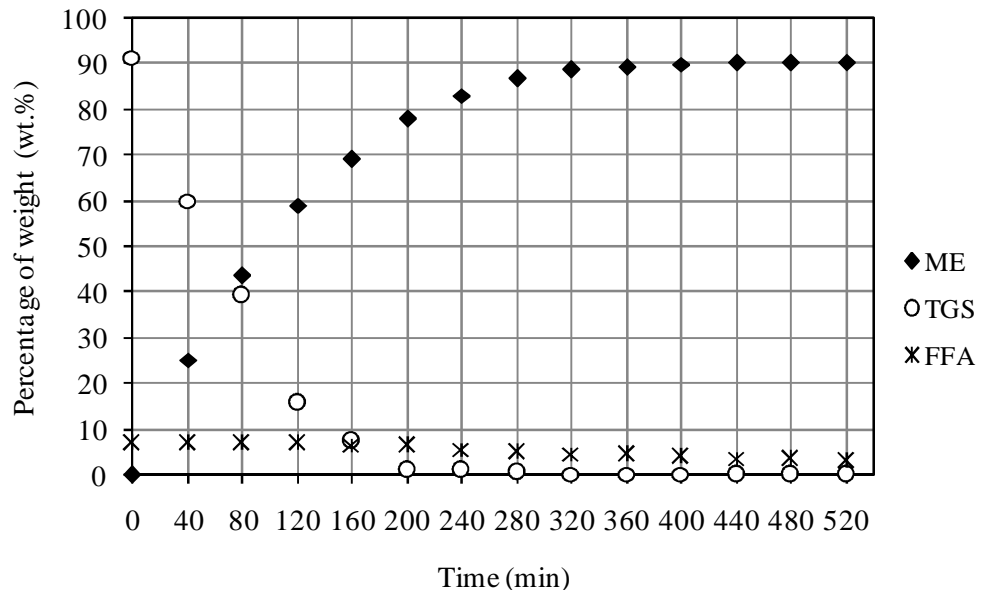


Figure F.3 The effect of BSPIMHSO₄ catalyst on the TGs and FFA conversion to ME.

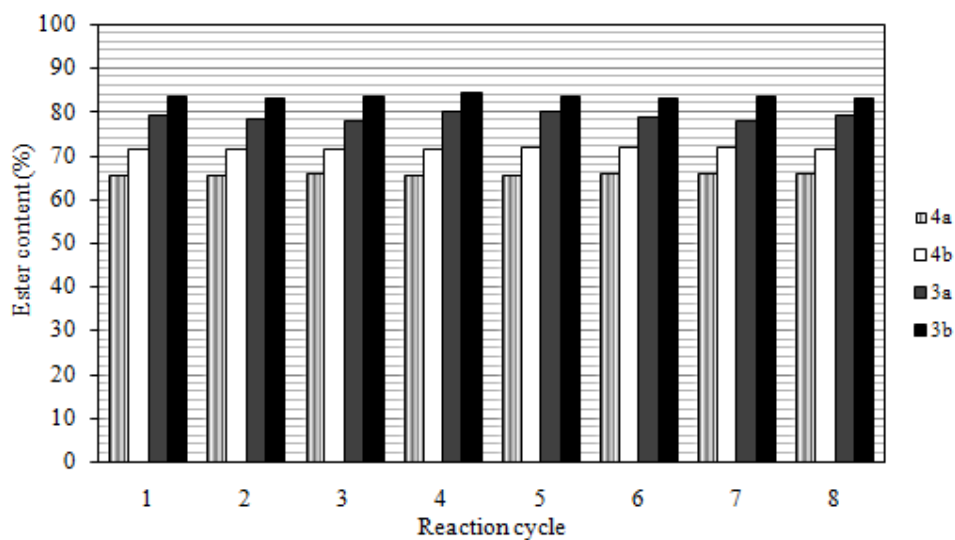


Figure G.1 Functionalized pyrazolium acidic ILs recyclability for catalyzing the transesterification of CPO with methanol under their optimal conditions. 3a = SPPHSO₄; 4a = MSPPHSO₄; 3b = SBPHSO₄; 4b = MSBPHSO₄.

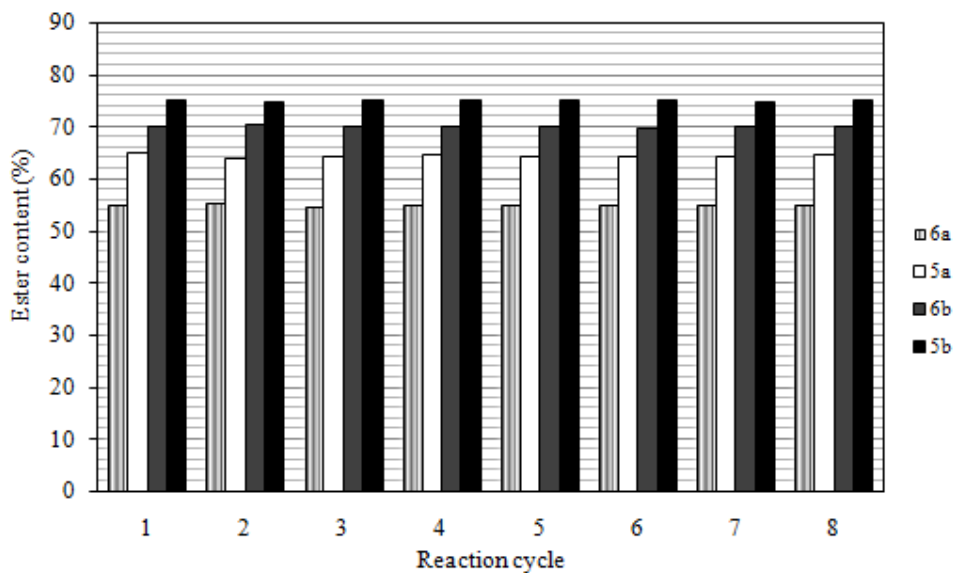
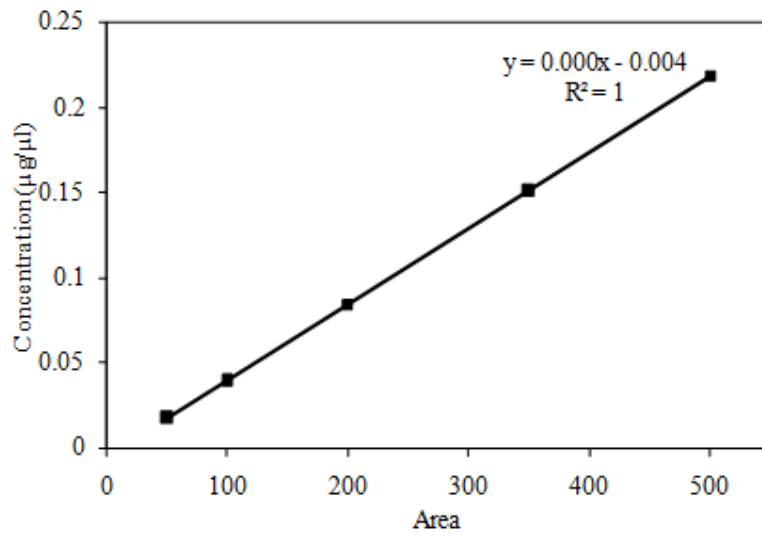
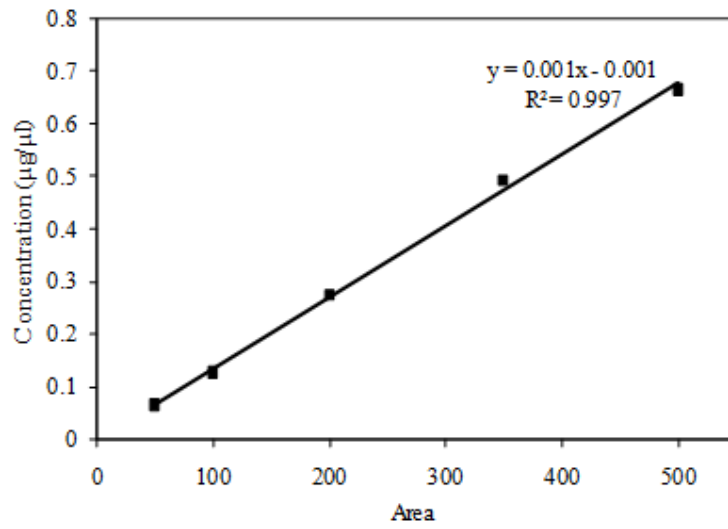


Figure G.2 Functionalized ammonium acidic ILs recyclability for catalyzing the transesterification of CPO with methanol under their optimal conditions. 3a = SPPHSO₄; 4a = MSPPHSO₄; 3b = SBPHSO₄; 4b = MSBPHSO₄.

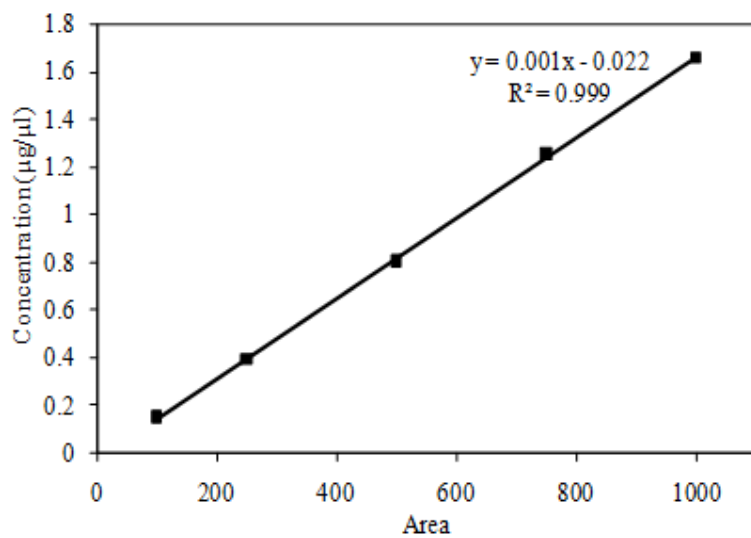
Appendix H



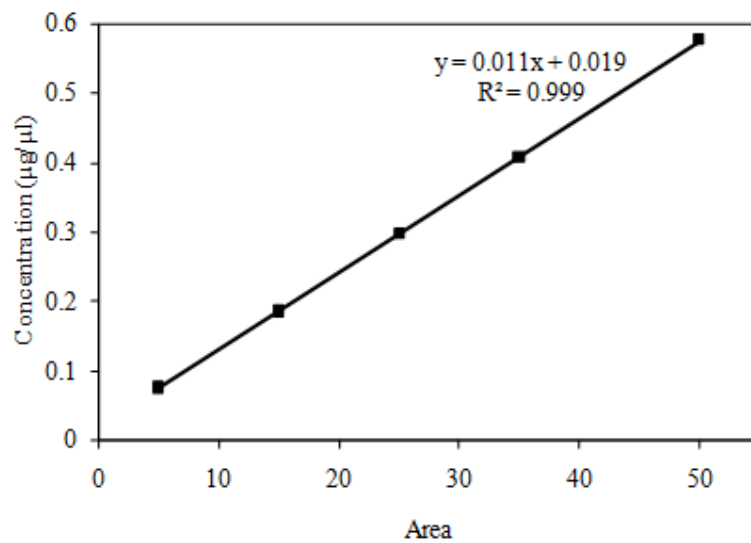
(a)



(b)



(c)



(d)

Figure I.1 GC standard curves for: (a) TG; (b) DG; (c) MG; (d) GI.