Nano-lipids based on ginger oil and lecithin as a potential drug delivery system

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Table S1: Compounds identified by GC-MS using NIST-14 Mass Spectral library

Peak	R.T.	Name of compound	Pct Total
1	23.584	Geraniol	1.949
2	37.032	n.i	0.474
3	39.479	n.i	1.378
4	40.138	n.i	3.113
5	40.41	(6)-Shogaol	63.137
6	41.549	[8]-Shogaol	6.29
7	41.644	n.i	5.09
8	42.449	n.i	1.438
9	42.679	n.i	5.671
10	43.871	Geranyl oleate	3.735
		[6]-Gingerdiol (2E)-geranial	
11	45.544	acetal	7.726

Figure list

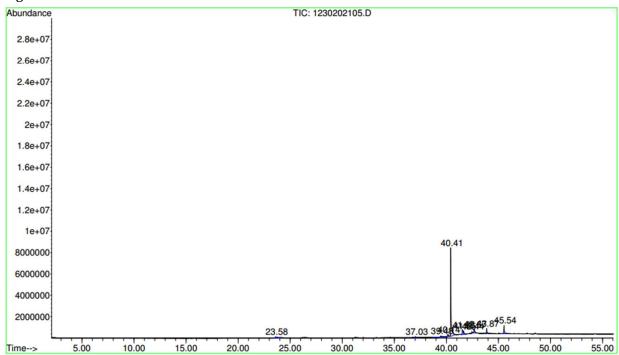
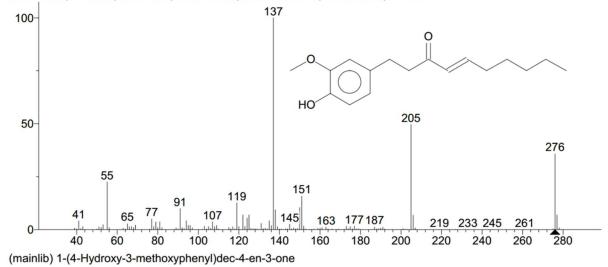


Figure S1: GC chromatograms of ginger oil

A) Hit 1 : 1-(4-Hydroxy-3-methoxyphenyl)dec-4-en-3-one C17H24O3; MF: 932; RMF: 950; Prob 96.5%; CAS: 555-66-8; Lib: mainlib; ID: 124142.



m/z 205 M⁺ m/z 276

Figure S2: Mass spectrum extracted from a peak 40.41 (A) and its fragmentation (B)

A) Hit 1 : 1-(4-Hydroxy-3-methoxyphenyl)dodec-4-en-3-one C19H28O3; MF: 654; RMF: 824; Prob 39.0%; CAS: 36700-45-5; Lib: mainlib; ID: 124143.

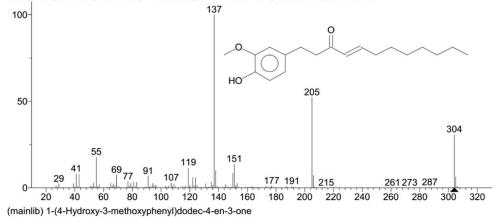
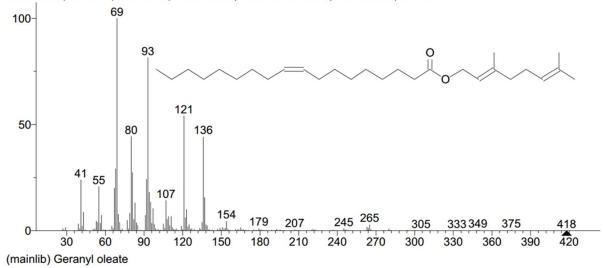


Figure S3: Mass spectrum extracted from a peak 41.45(A) and its fragmentation (B)

Hit 5 : Geranyl oleate C28H50O2; MF: 604; RMF: 678; Prob 4.85%; CAS: 81601-03-8; Lib: mainlib; ID: 35714.



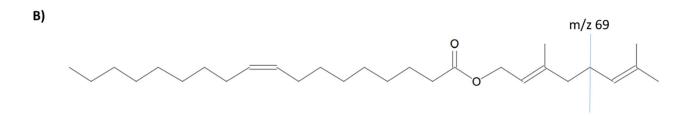
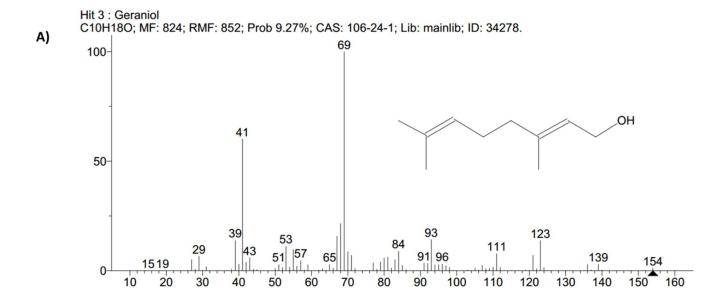
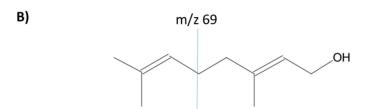


Figure S4: Mass spectrum extracted from a peak 43.88 (A) and its fragmentation (B)





(mainlib) Geraniol

Figure S5: Mass spectrum extracted from a peak 23.57 (A) and its fragmentation (B)

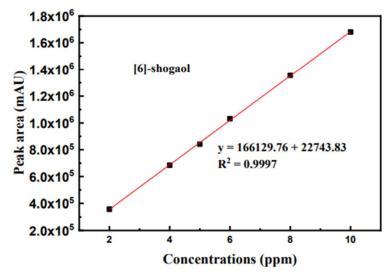


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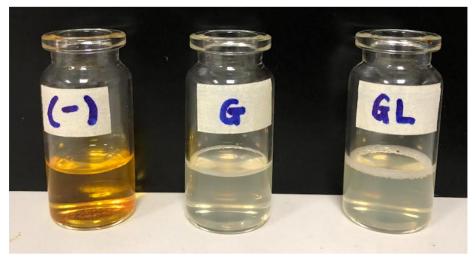


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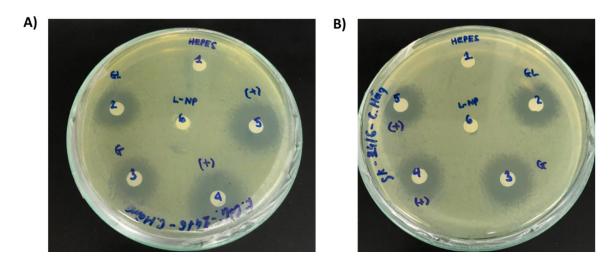


Figure S8: Zones of Inhibitions formed on nutrient agar surface by Hepes (1) GL nano-lipid (2), ginger oil (3), and two positive controls (Ampicillin (4) and Streptomycin(5)) L nano-lipid (6) against E.coli (A) and S.aureus (B).

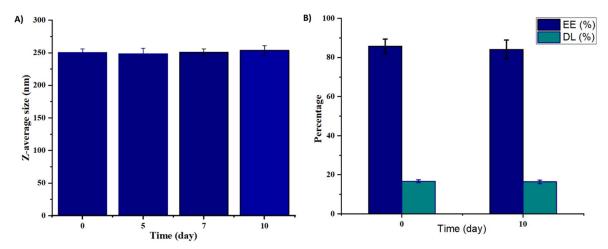


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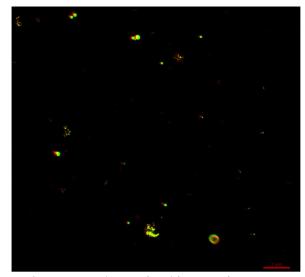


Figure S10: The confocal image of Cur@ GL

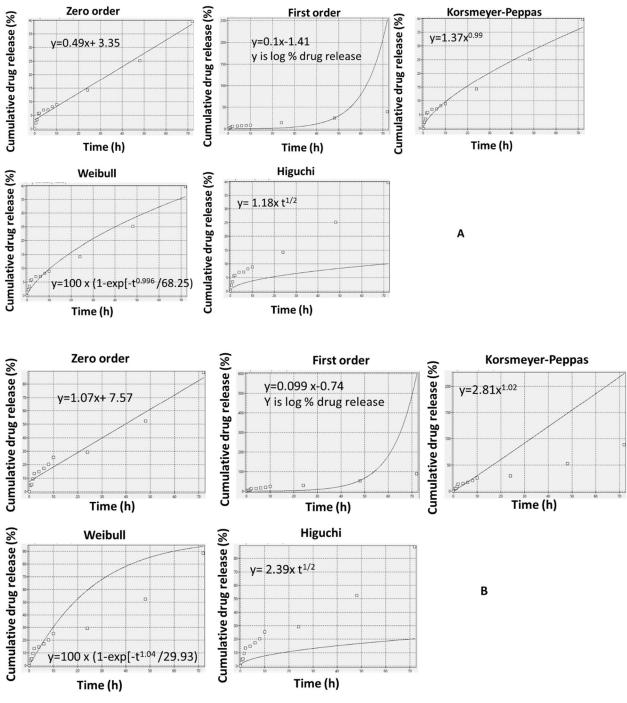


Figure S11: Release kinetics of Cur@ GL at pH 7.4 (A) and at pH 5.5 (B) fitted to 5 kinetic models contructing by KinetDS