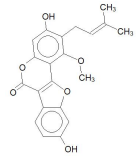
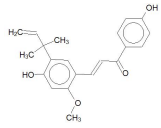
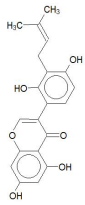
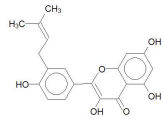
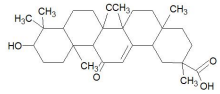
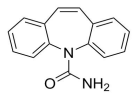
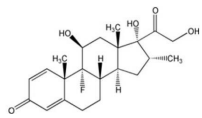
Supplementary Table 4. Molecular docking score for ingredients have strongly energy binding to CYP3A4.

Molecule Name	Structure	Score (docking of CYP3A4)
GLL		4.7912
LCA		6.0907
LIA		5.6487
ILF		6.0287
18β-GA		4.4983
carbamazepine		4.3030
Dexamethasone		4.3246