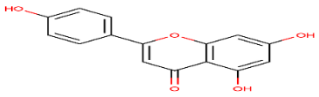
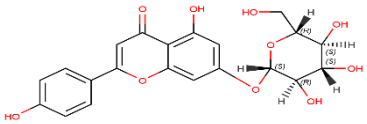
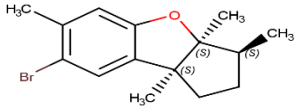
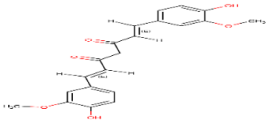
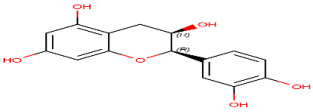
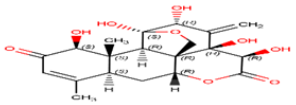
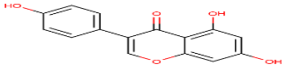
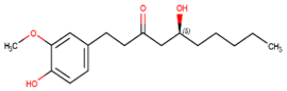
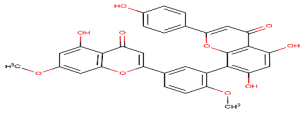
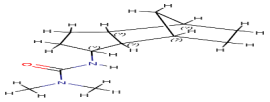
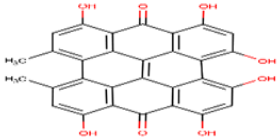
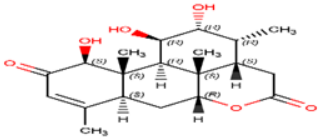
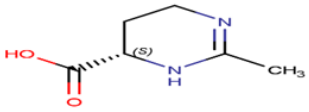
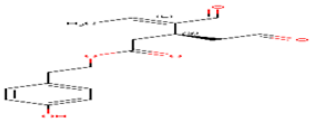
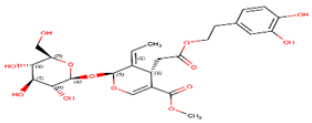
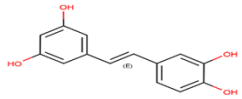
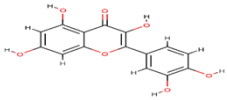
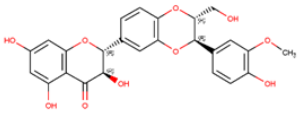
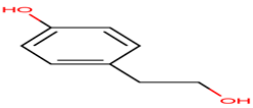
Table S1: The list of selected plant-derived compounds for molecular docking analysis

Row	Compound name	Molecular Weight (g/mol)	Molecular Formula	PubChem CID	Chemical Structure
1	Apigenin	270.24 g/mol	$C_{15}H_{10}O_5$	5280443	
2	Apigetrin	432.381 g/mol	$C_{21}H_{20}O_{10}$	5280704	
3	Aplysin	295.22 g/mol	$C_{15}H_{19}BrO$	11066347	
4	Curcumin	368.385 g/mol	$C_{21}H_{20}O_6$	969516	
5	Epicatechin	290.271 g/mol	$C_{15}H_{14}O_6$	72276	
6	Eurycomanone	408.403 g/mol	$C_{20}H_{24}O_9$	13936691	
7	Genistein	270.24 g/mol	$C_{15}H_{10}O_5$	5280961	
8	Gingerol	294.391 g/mol	$C_{17}H_{26}O_4$	442793	
9	Ginkgetin	566.518 g/mol	$C_{32}H_{22}O_{10}$	5271805	
10	Isonoruon	222.332 g/mol	$C_{13}H_{22}N_2O$	3035253	
11	Hypericin	504.45 g/mol	$C_{30}H_{16}O_8$	5281051	
12	Klaineanone	364.438 g/mol	$C_{20}H_{28}O_6$	12304890	
13	L-Ectone	142.158 g/mol	$C_6H_{10}N_2O_2$	126041	

14	Oleocanthal	304.342 g/mol	$C_{17}H_{20}O_5$	11652416	
15	Oleuropein	540.518 g/mol	$C_{25}H_{32}O_{13}$	5281544	
16	Piceatannol	244.246 g/mol	$C_{14}H_{12}O_4$	667639	
17	Pterostilbene	256.301 g/mol	$C_{16}H_{16}O_3$	5281727	
18	Quercetin	302.238 g/mol	$C_{15}H_{10}O_7$	5280343	
19	Silibinin	482.441 g/mol	$C_{25}H_{22}O_{10}$	31553	
20	Tyrosol	138.166 g/mol	$C_8H_{10}O_2$	10393	
+	Paclitaxel (positive control)	853.918 g/mol	$C_{47}H_{51}NO_{14}$	36314