

Figure 1.

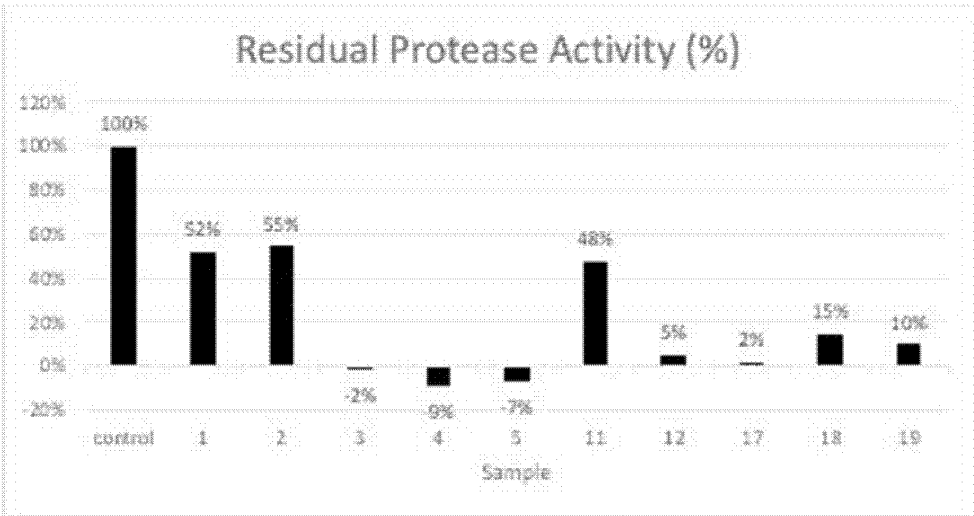


Figure 2

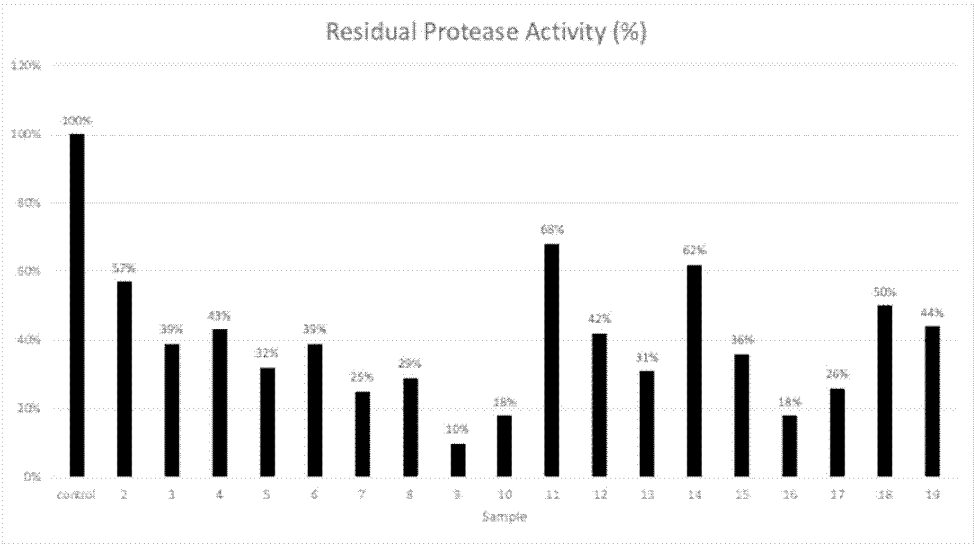


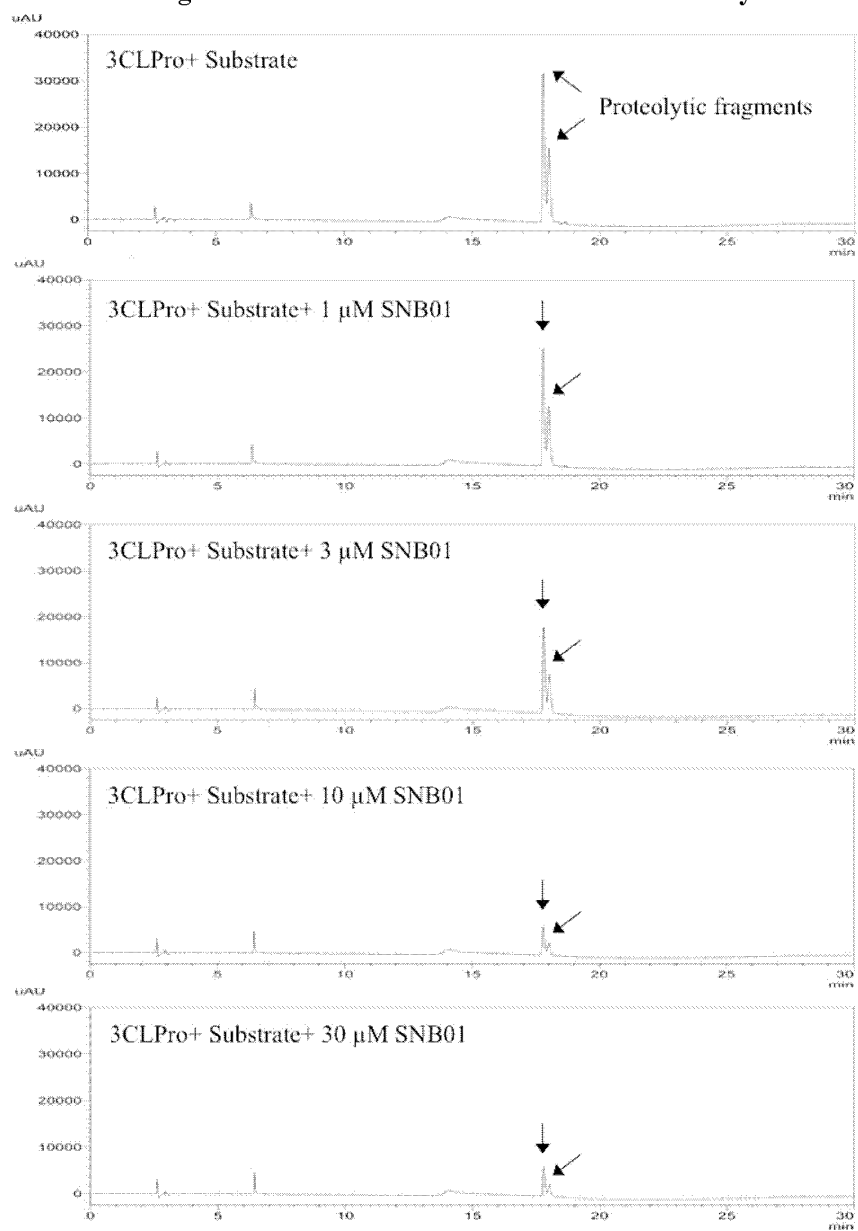
Figure 3. Inhibition of 3CLPro of SARS-CoV-2 by SNB01

Figure 4.

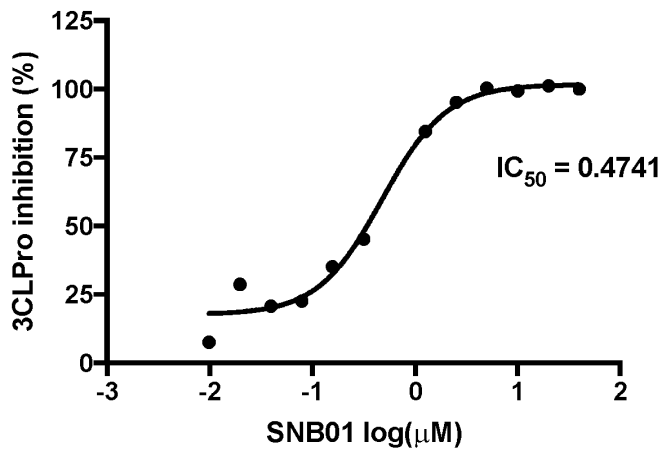


Figure 5.

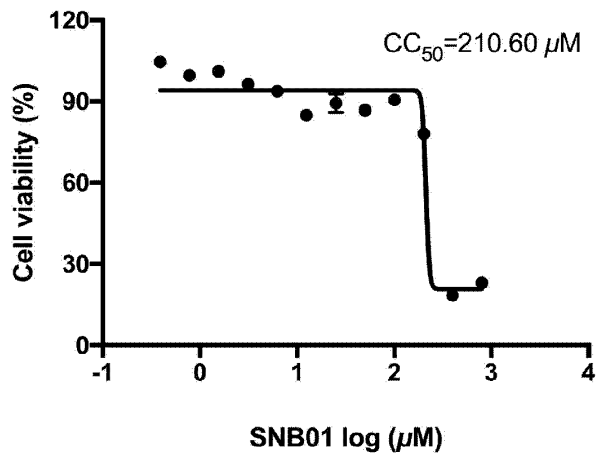


Figure 6.

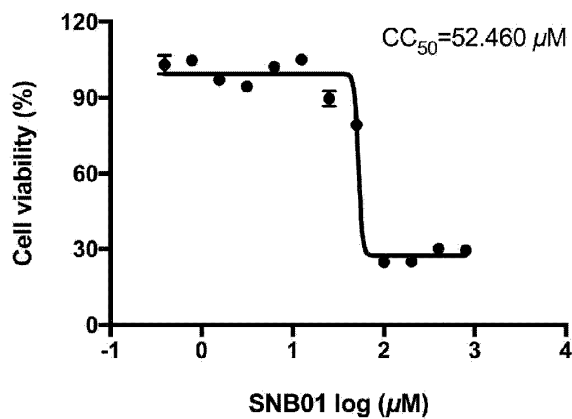


Figure 7A.

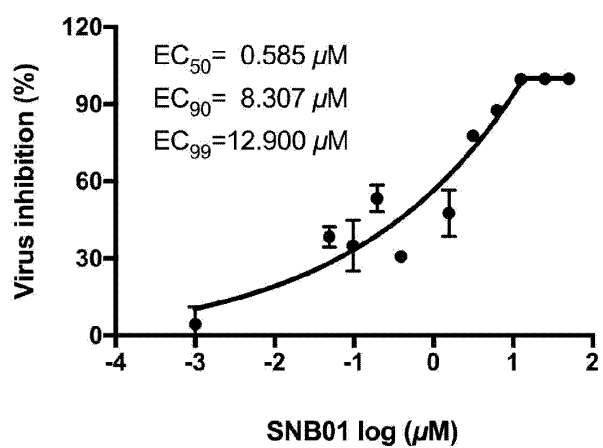


Figure 7B

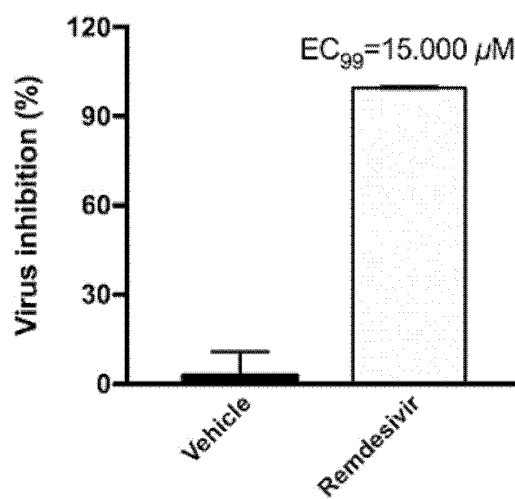


Figure 8A.

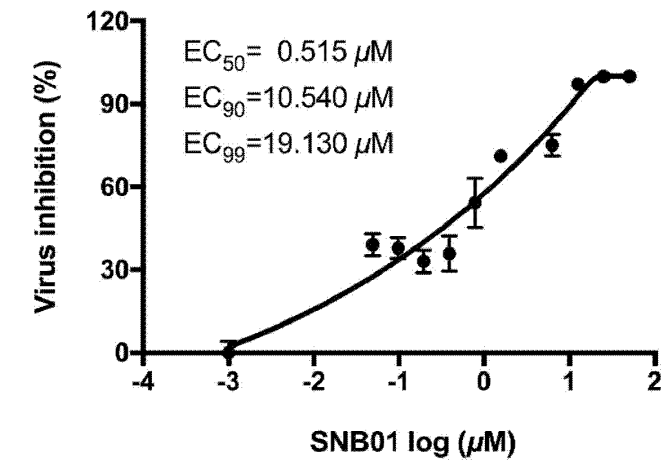


Figure 8B

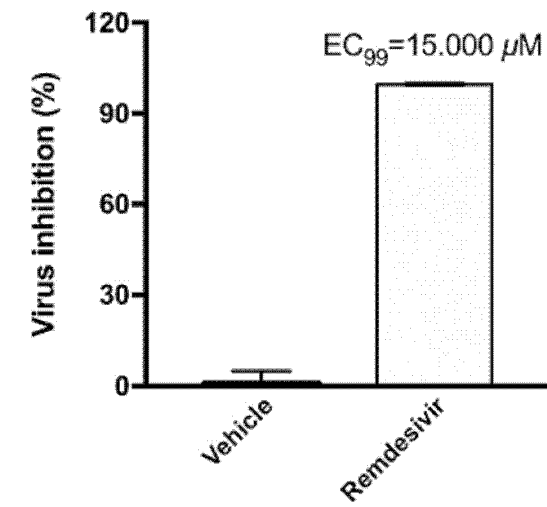
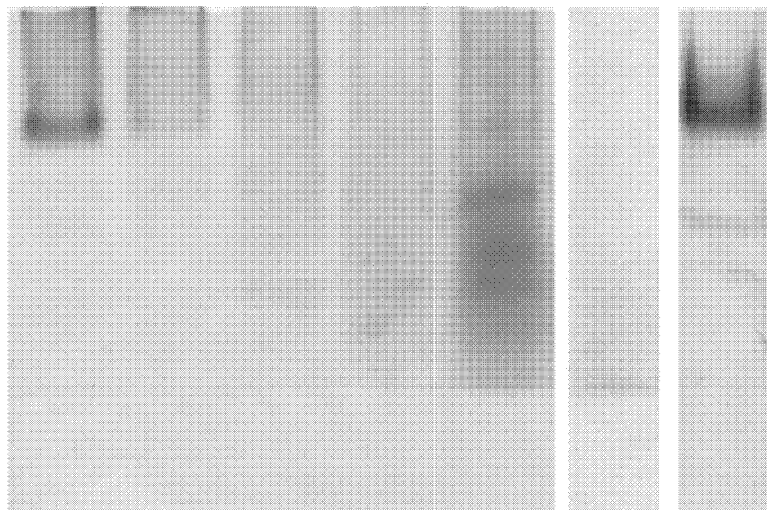


Figure 9.

SNB01 (μM)	31.25	62.5	125	250	500	500	0
3CLPro (μM)	20	20	20	20	20	0	20

**Figure 10.**

β -10G (μM)	31.25	62.5	125	250	500	500	0
3CLPro (μM)	20	20	20	20	20	0	20

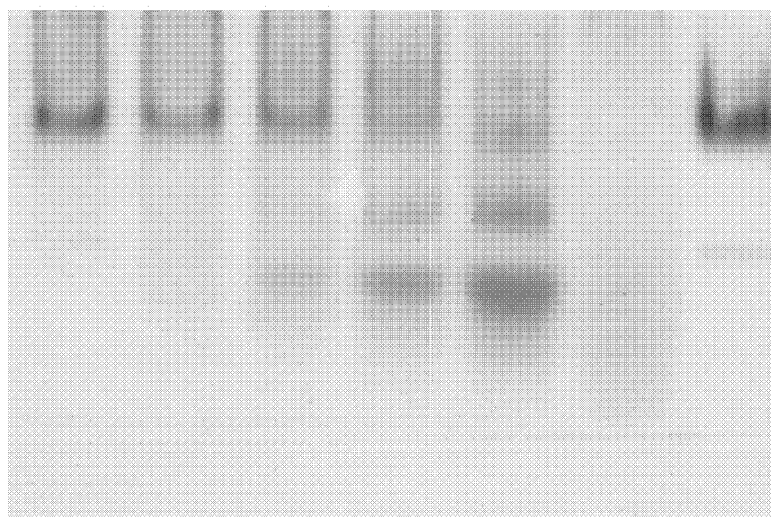
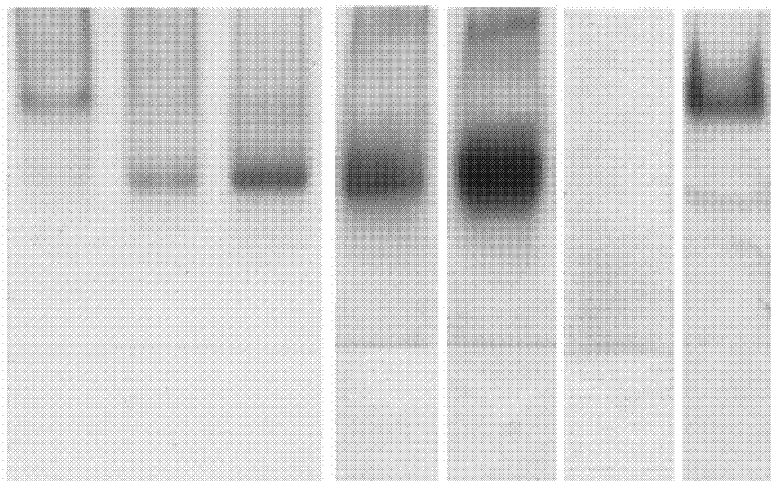


Figure 11.

β -15G (μ M)	31.25	62.5	125	250	500	500	0
3CLPro (μ M)	20	20	20	20	20	0	20

**Figure 12.**

GA (μ M)	31.25	62.5	125	250	500	500	0
3CLPro (μ M)	20	20	20	20	20	0	20

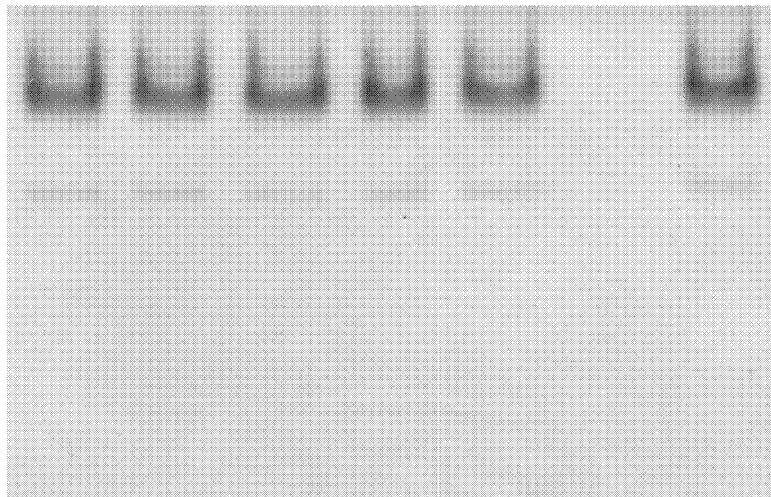


Figure 13.

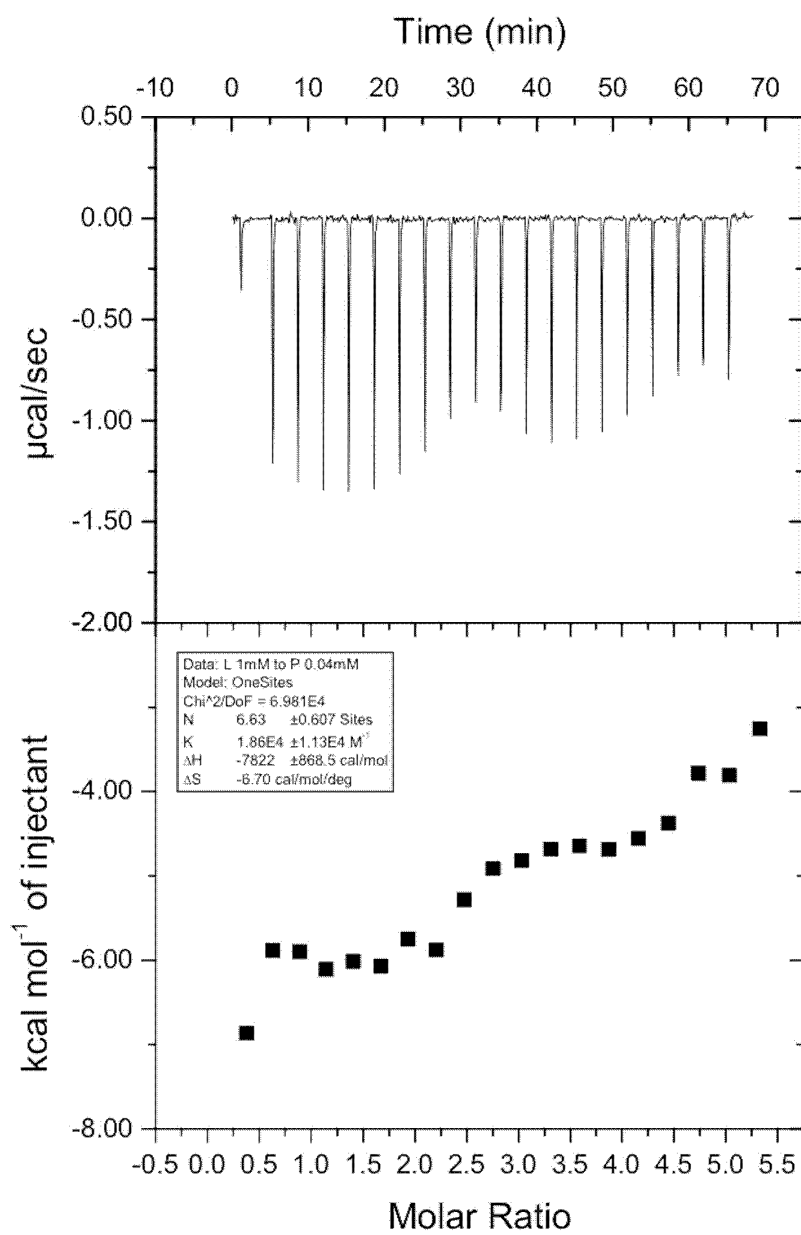


Figure 14.

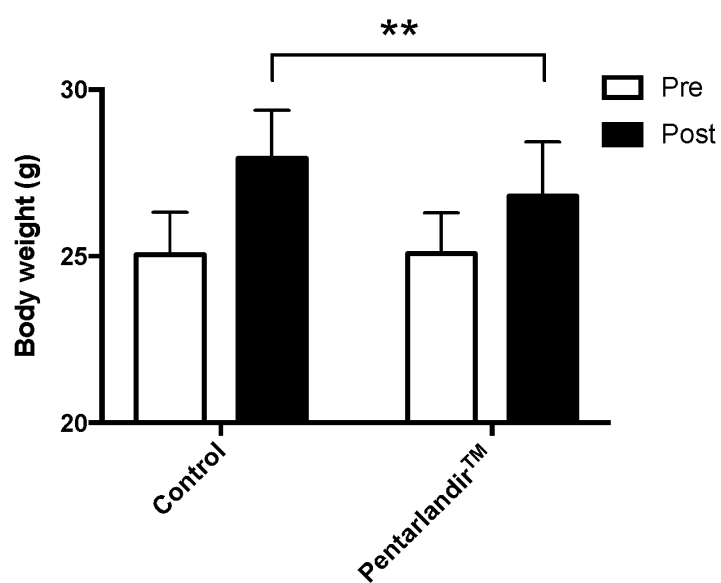


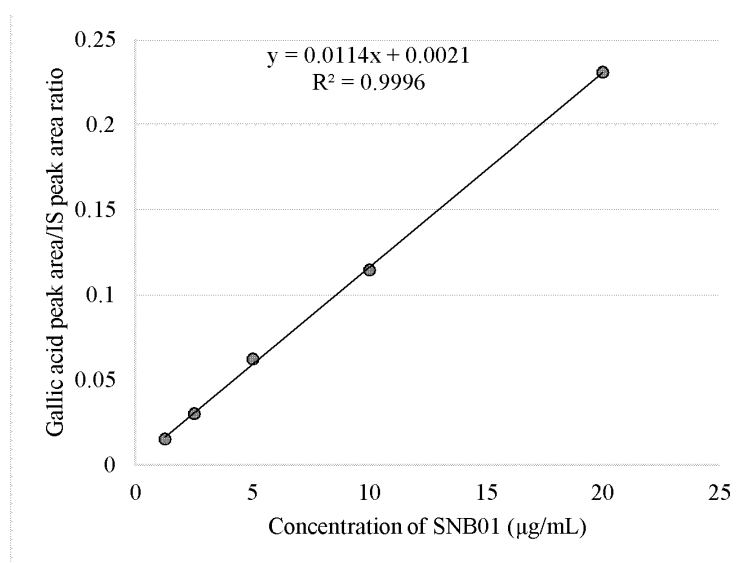
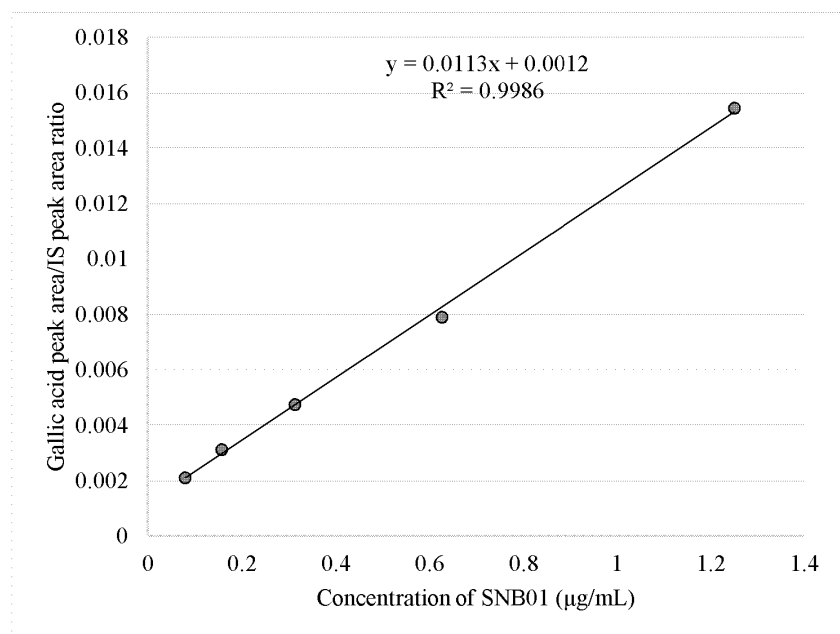
Figure 15A.**Figure 15B**

Figure 16.

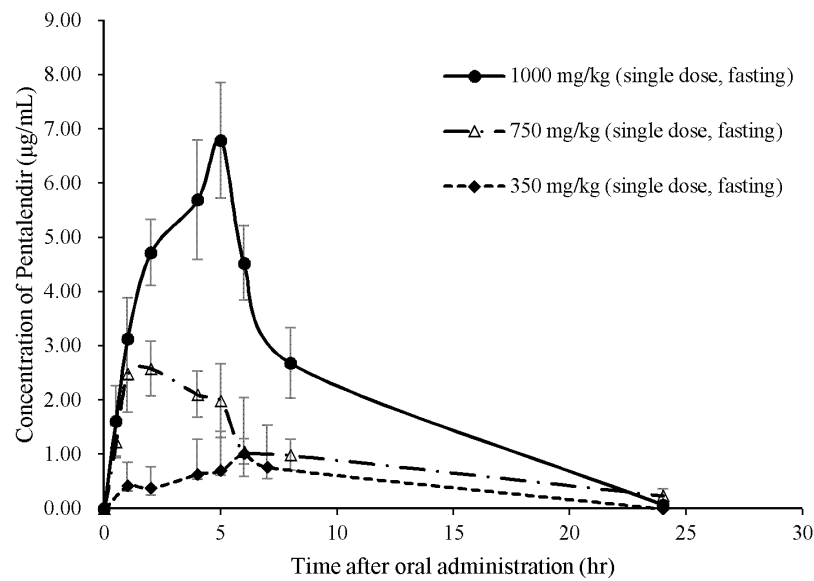


Figure 17.

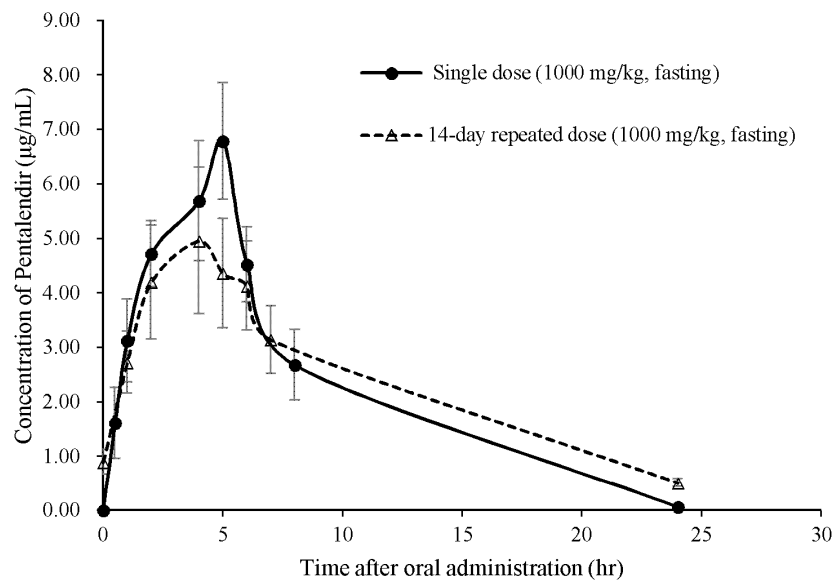


Figure 18.

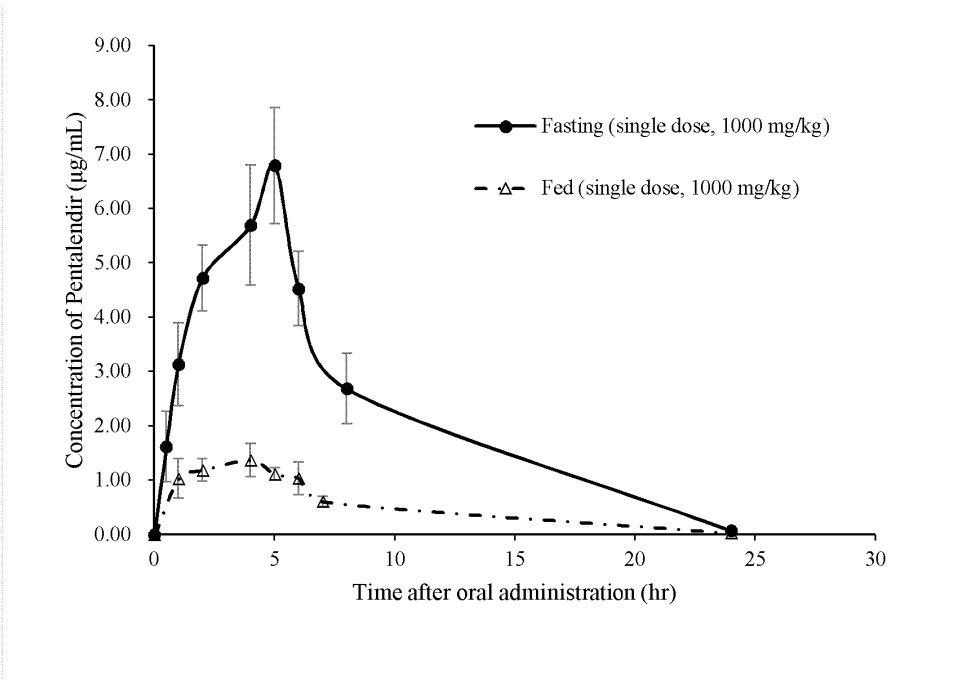


Figure 19.

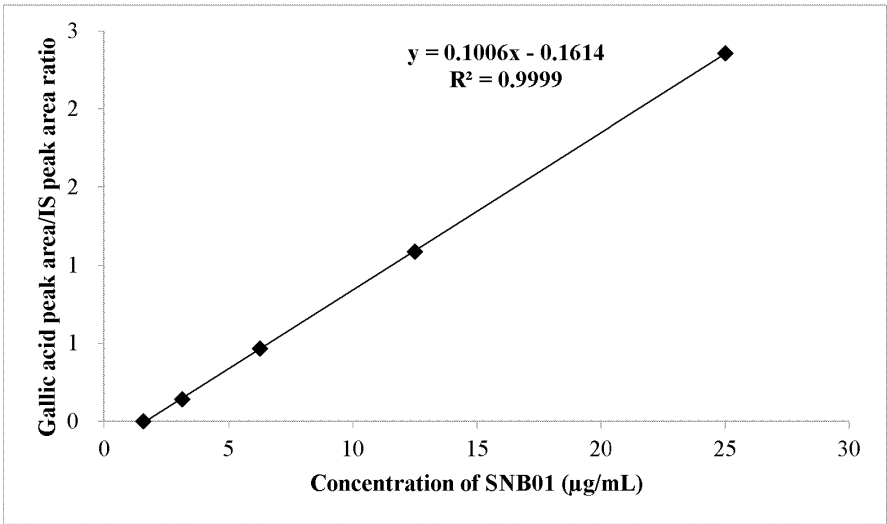
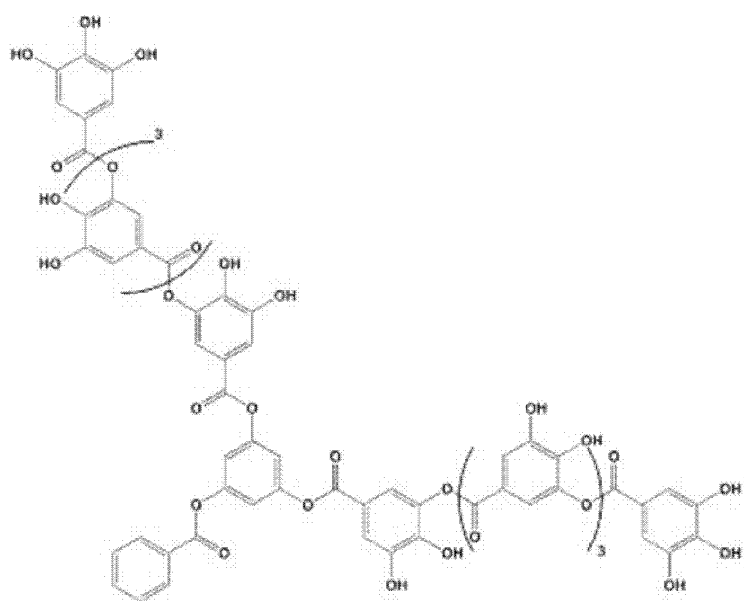
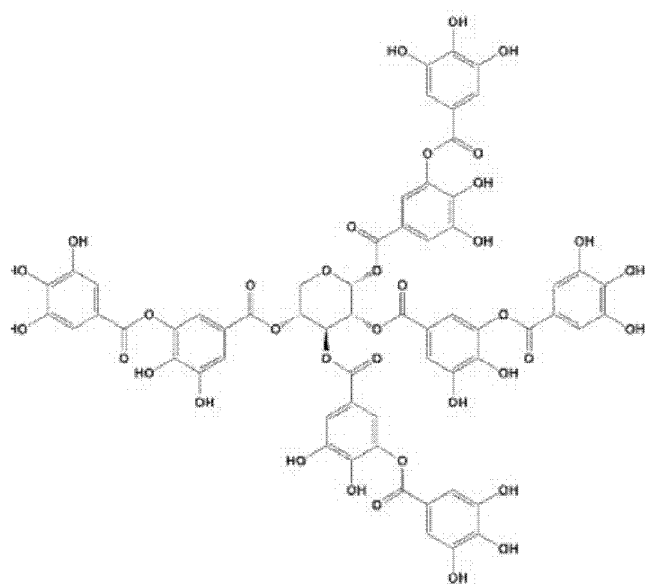


Figure 20

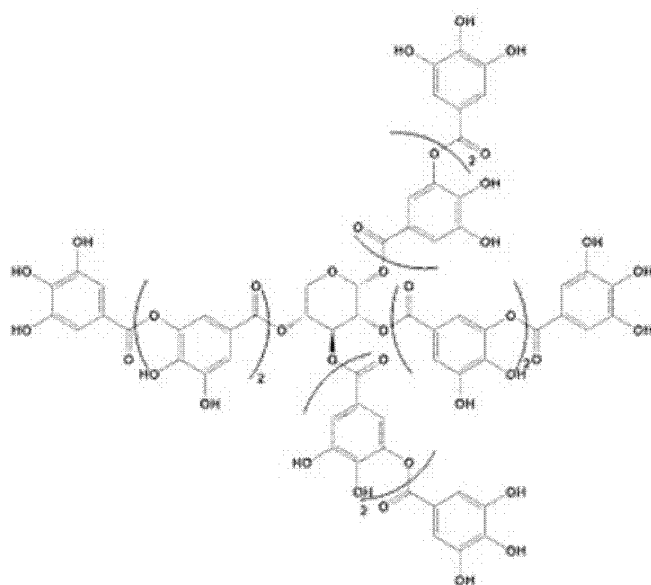


Compound 103

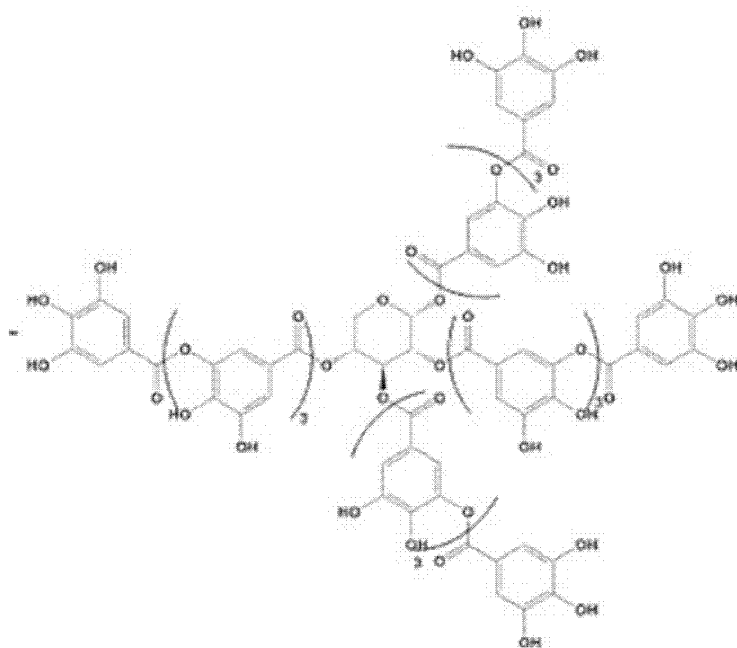


Compound 117

Figure 20 (cont'd)

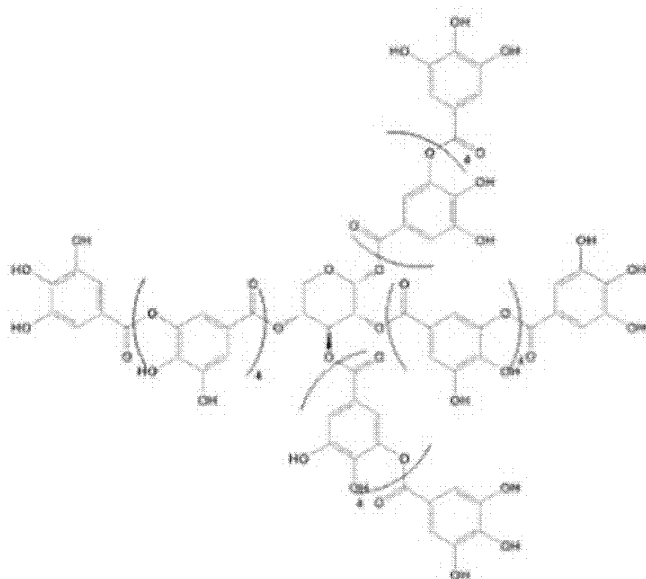


Compound 119

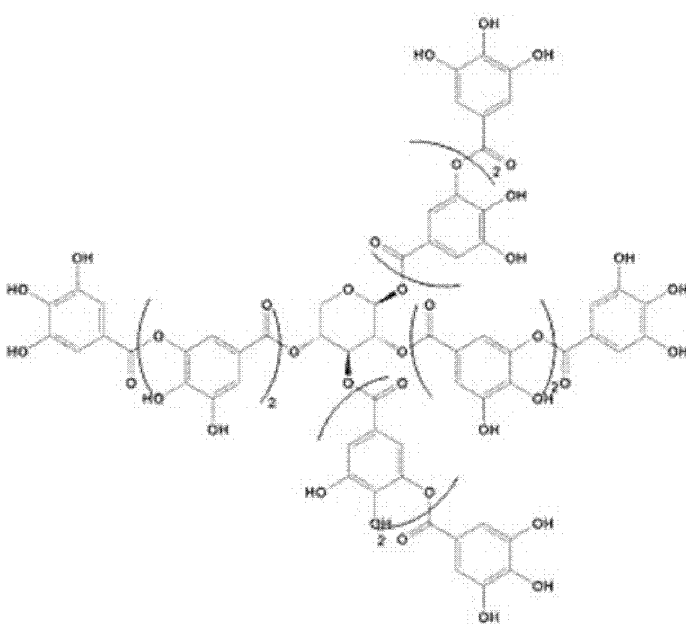


Compound 121

Figure 20 (cont'd)

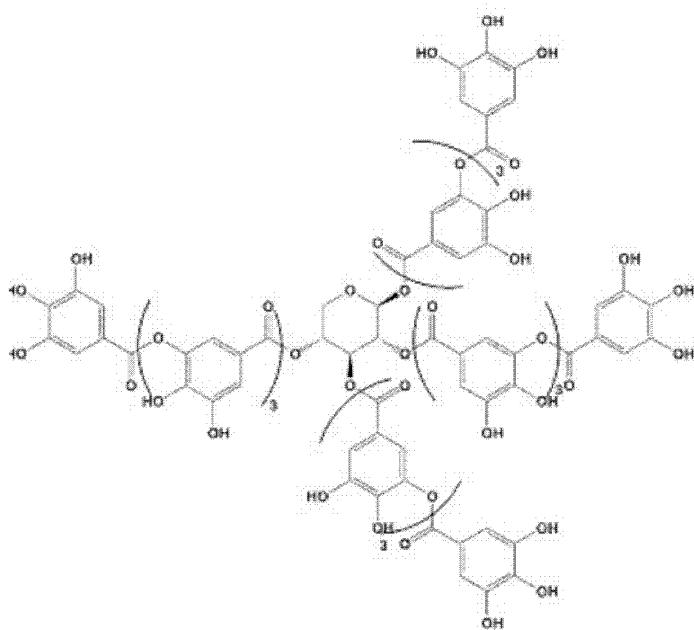


Compound 123

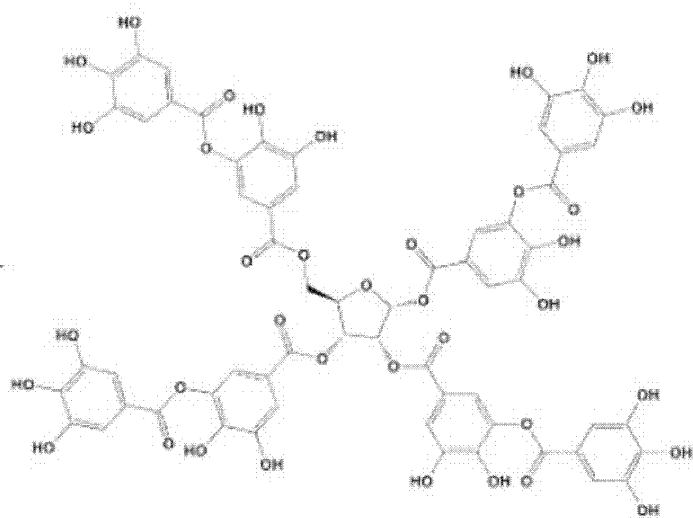


Compound 126

Figure 20 (cont'd)

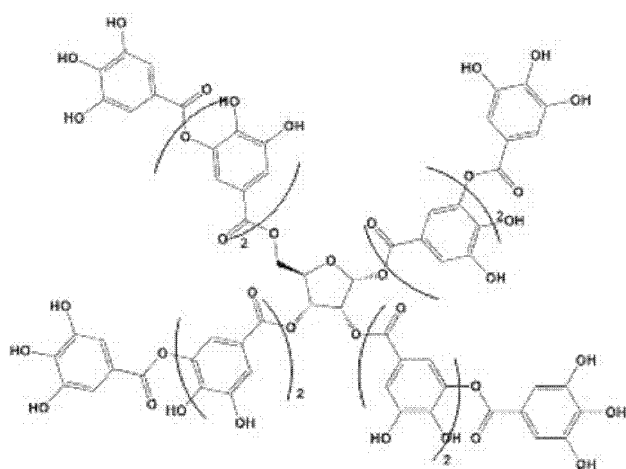


Compound 128

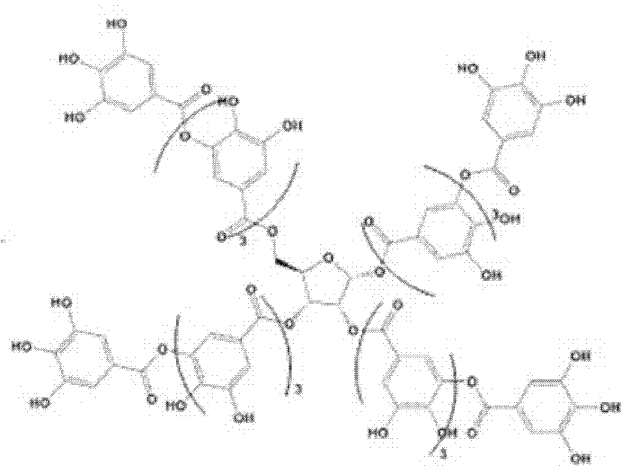


Compound 134

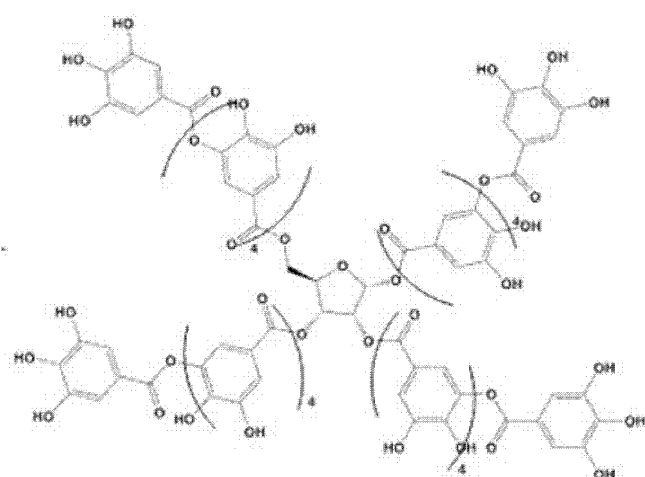
Figure 20 (cont'd)



Compound 136

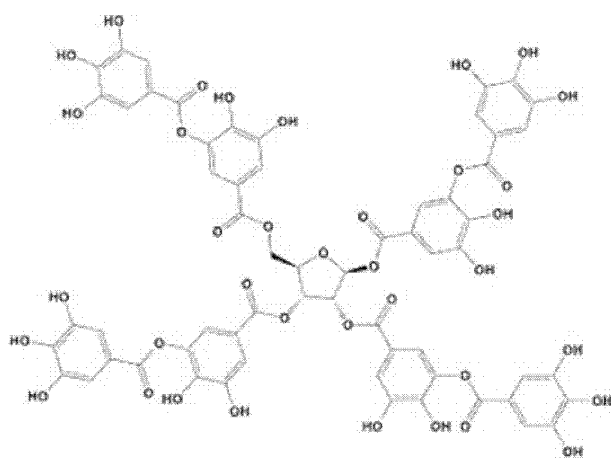


Compound 138

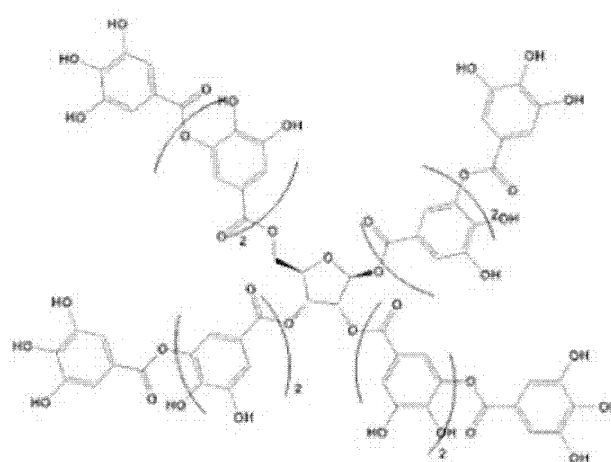


Compound 140

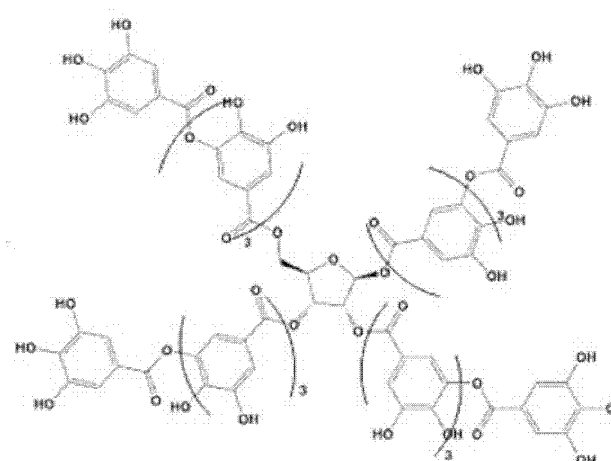
Figure 20 (cont'd)



Compound 142



Compound 144



Compound 146

The chemical structure shows a central furan ring substituted at the 2 and 5 positions with two 2,4,6-trihydroxyphenyl groups. At the 3 and 4 positions of the furan ring, there are two ether linkages connecting to two additional 2,4,6-trihydroxyphenyl groups. Each of these four phenyl groups is further substituted with a 2,4,6-trihydroxyphenyl group via an ester linkage. The structure is symmetrical and features multiple hydroxyl groups, suggesting it is a precursor for a poly(phenylene ether) or a similar polymer.

O=C(O)c1ccc(O)c(OC(=O)c2ccc(O)c(O)c2OC(=O)c3ccc(O)c(O)c3)c1

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