Supplementary

In Silico Identification and Evaluation of Natural Products as Potential Tumor Necrosis Factor Function Inhibitors Using Advanced Enalos Asclepios KNIME Nodes

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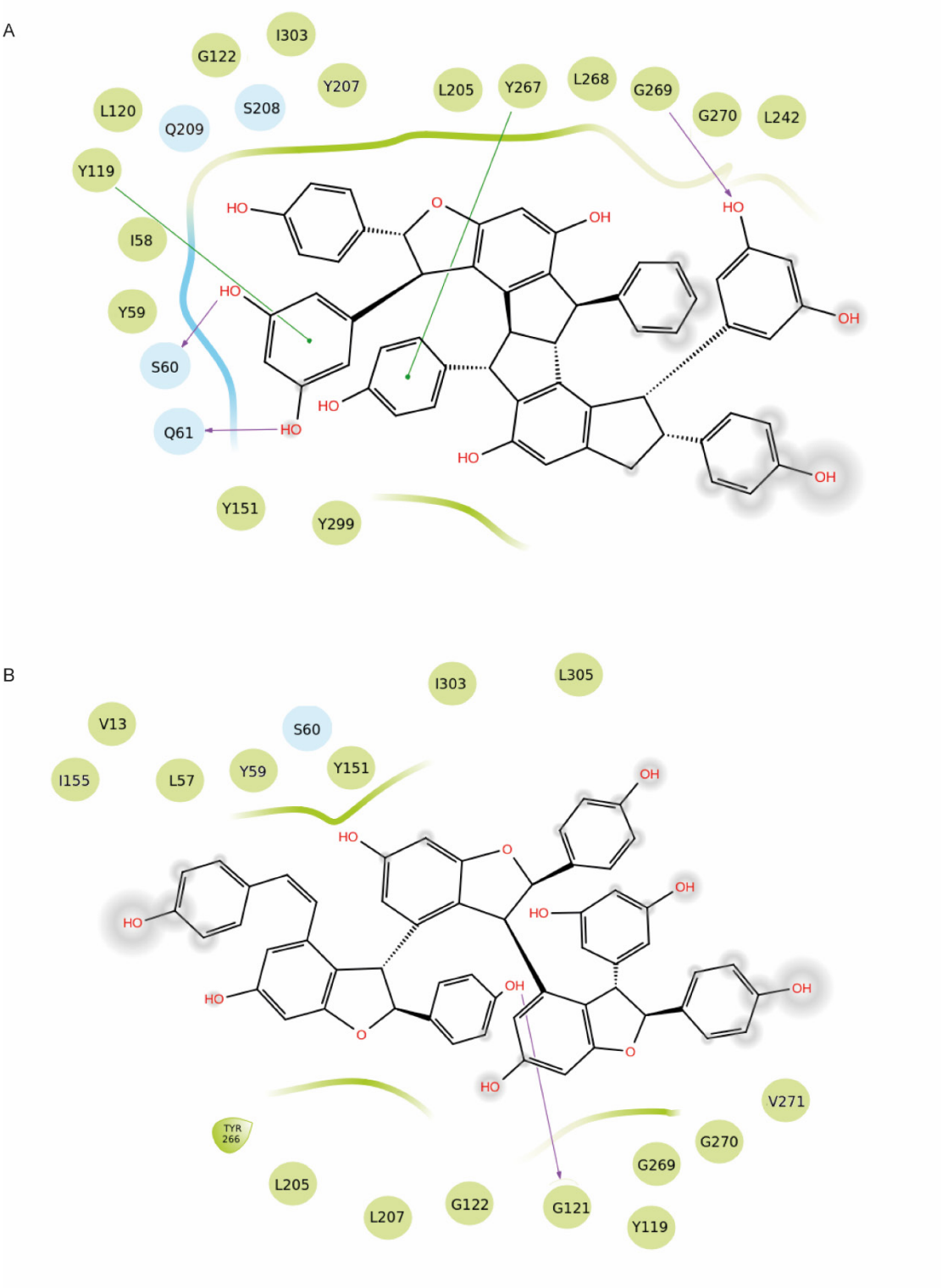
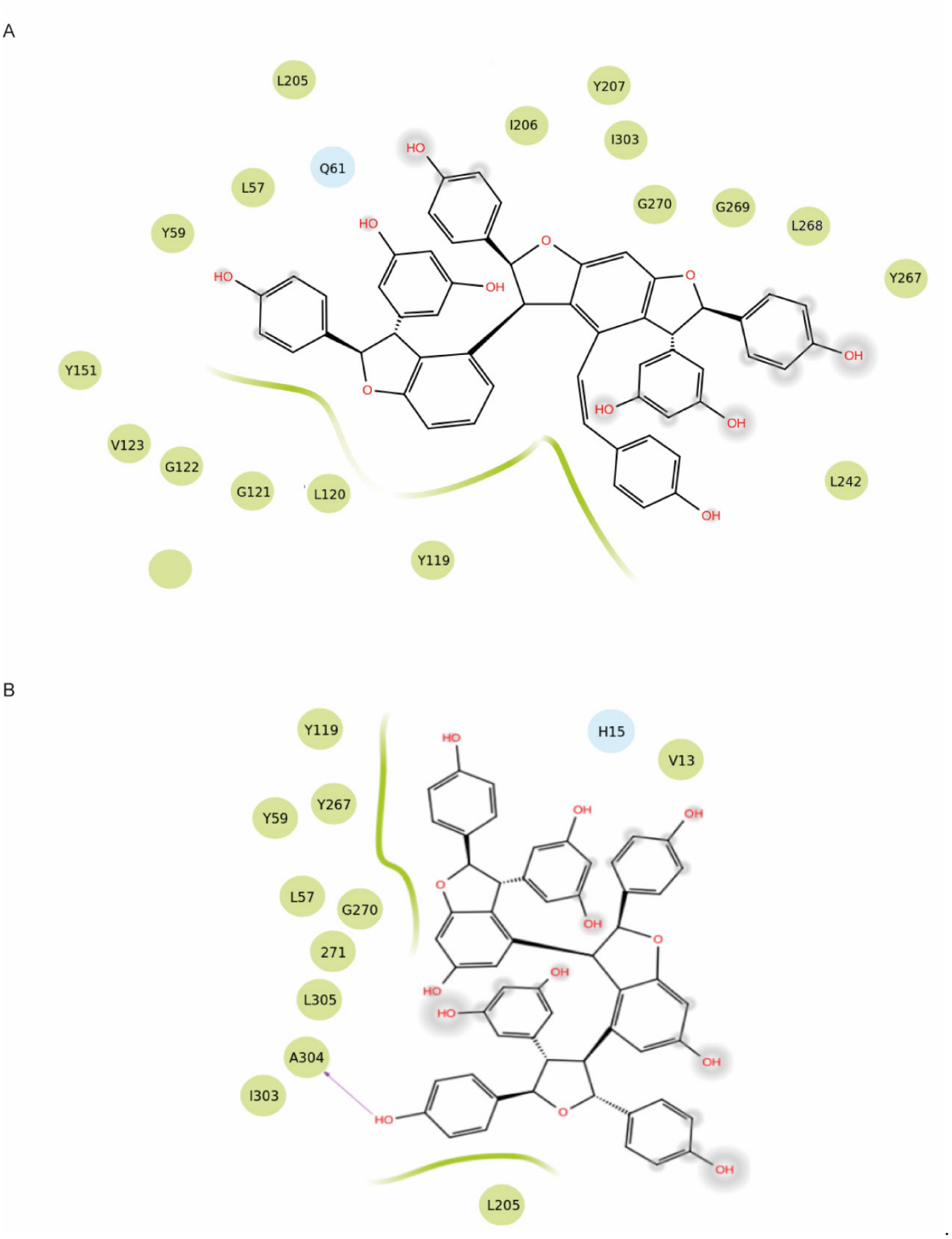


Supplementary Figure S1. SAR in 2D A. Nepalensinol B, B. Miyabenol A.



Supplementary Figure S2. SAR in 2D A. Flexuosol A, B. Kobophenol A

Ligands Name	Docking Scoore (rDock)
Ampelopsin H	-45.14
SPD304	-54.269
Nepalensinol B	-35.722
Miyabenol A	-23.104
Flexuosol A	-32.603
Kobophenol A	-43.617
Vitisin B	-37.549
trans-Diptoindonesin B	-44.674

Supplementary Figure S3. Docking scores of compounds