

In Silico Exploration of Bisphosphonate Scaffolds as Potential Inhibitors of SARS-CoV-2 RdRp for COVID-19 and PASC

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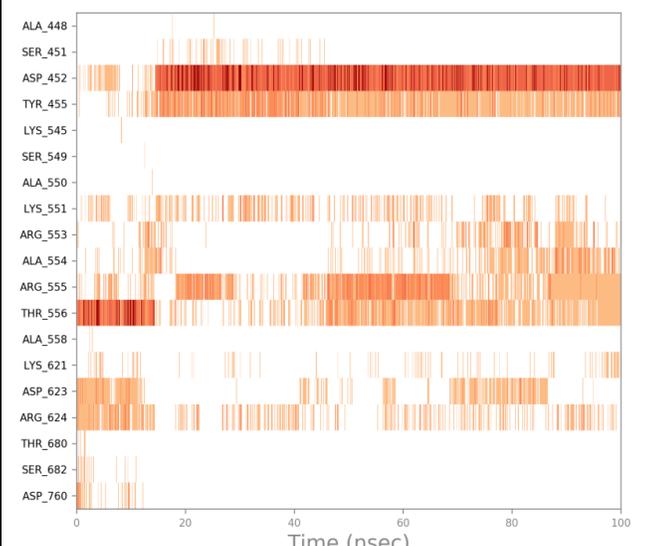
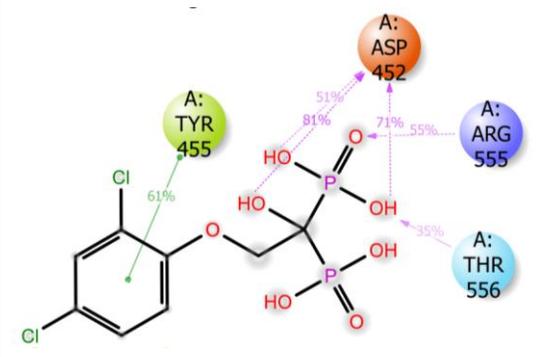
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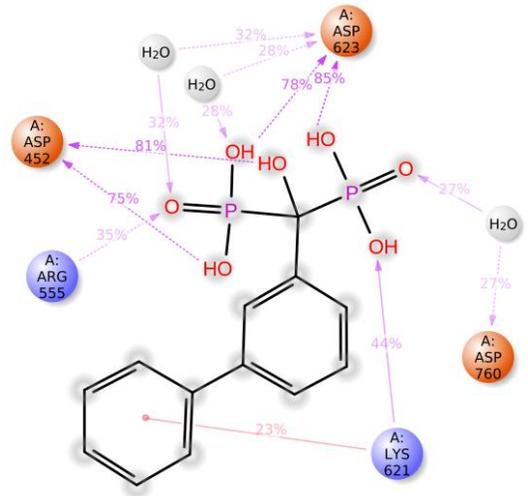
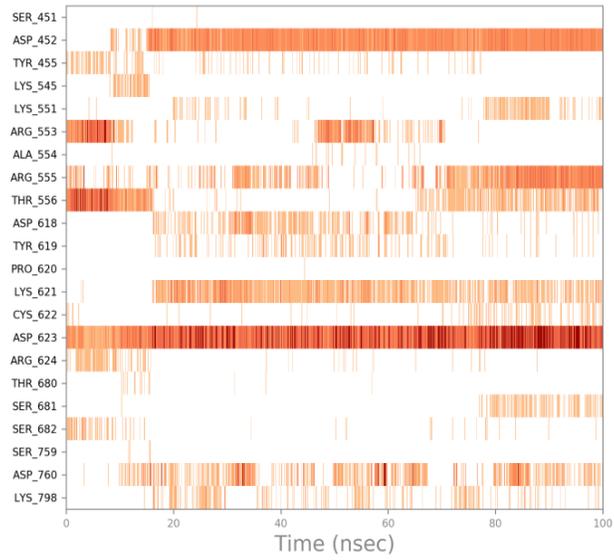
† These authors contributed equally to this work.

Table 1: Timeline plots and 2D interaction diagrams of the selected seven hit compounds

S. no	Time-line plot	2D interaction diagram
1.	<p>CHEMBL164344</p> 	

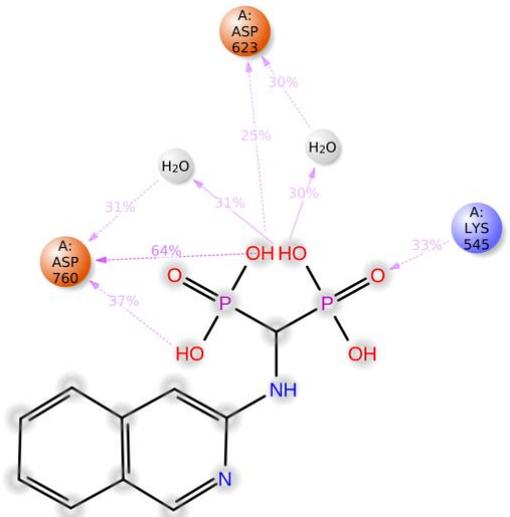
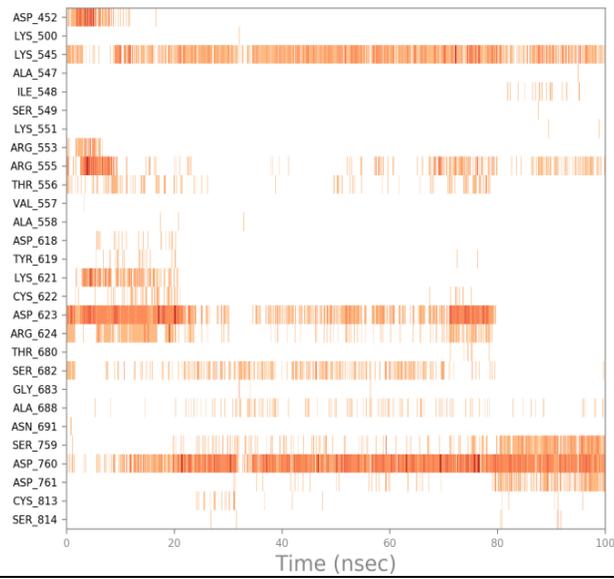
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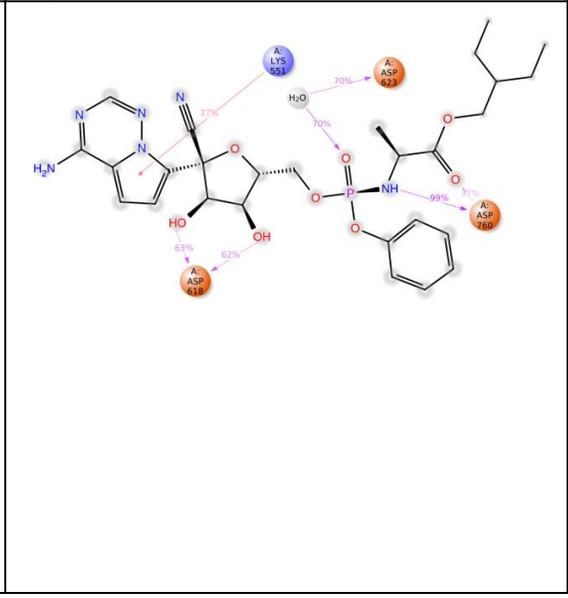
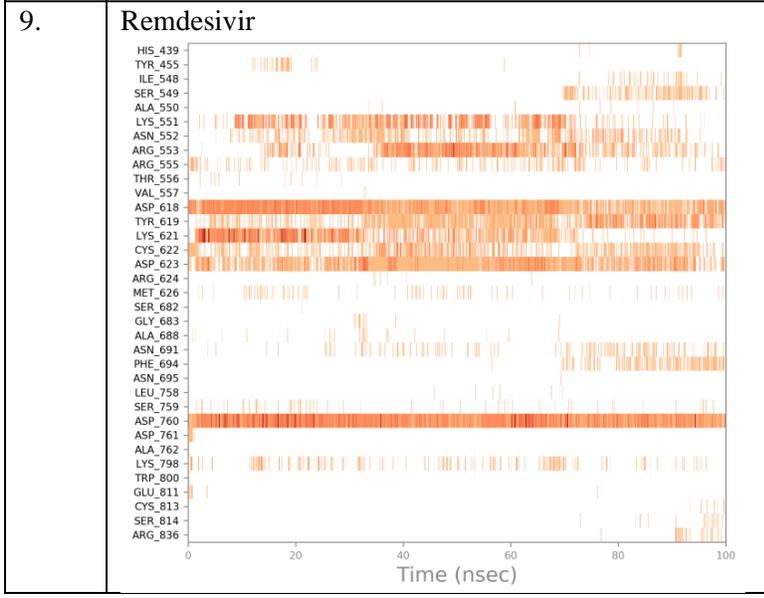
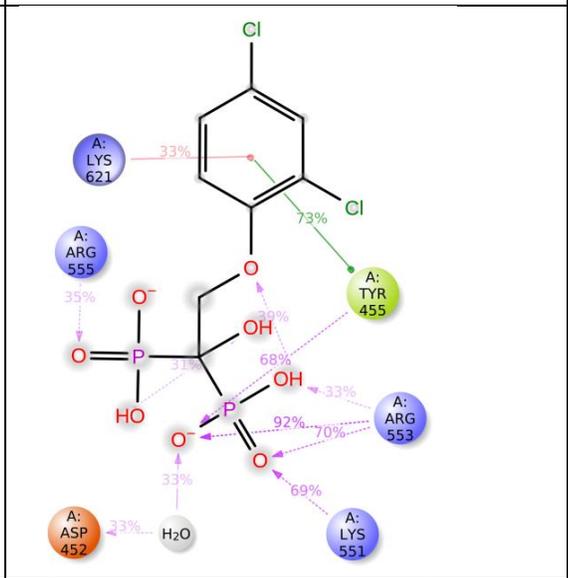
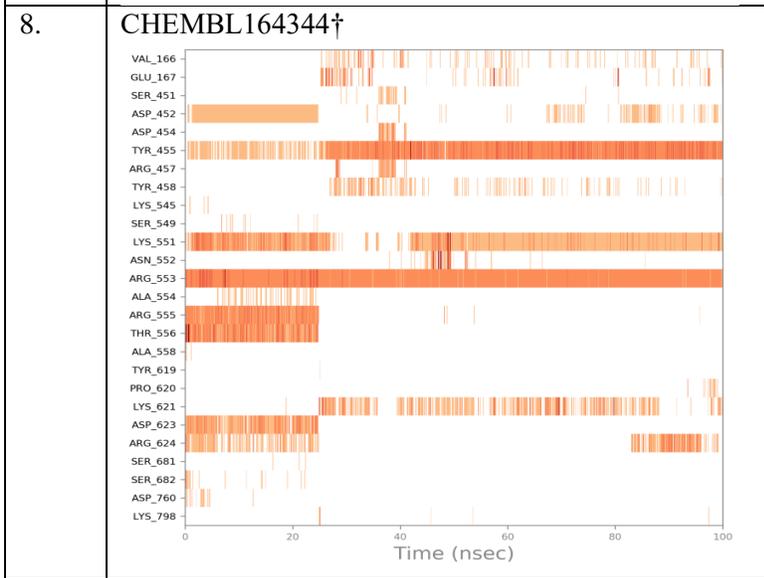
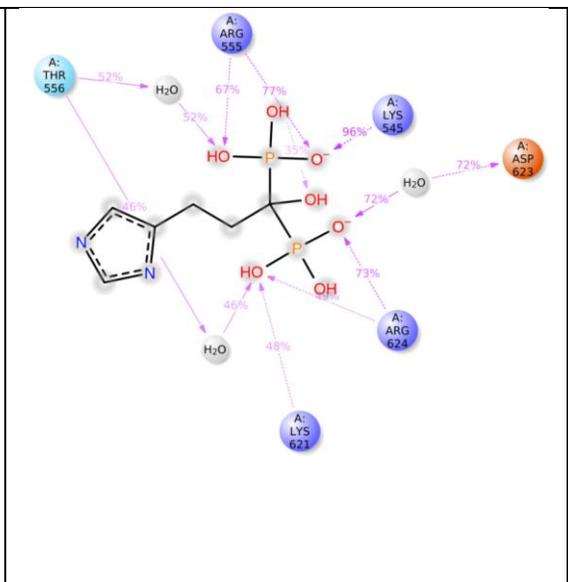
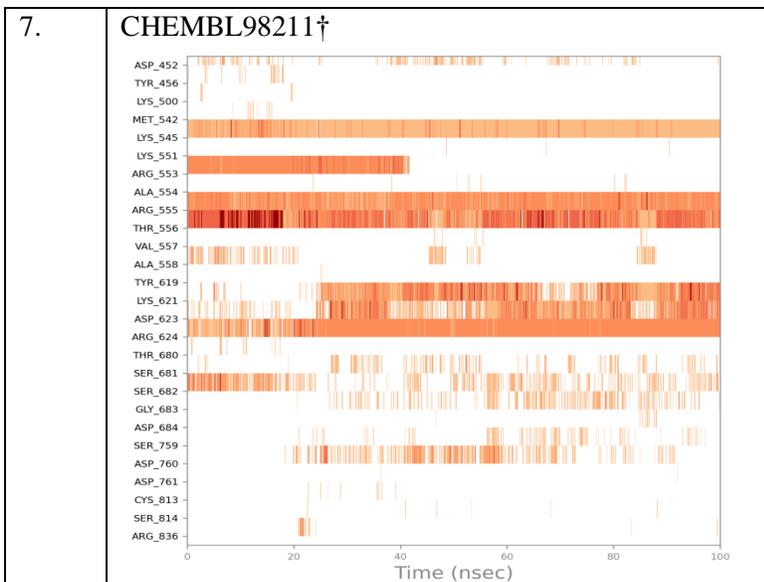
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3.

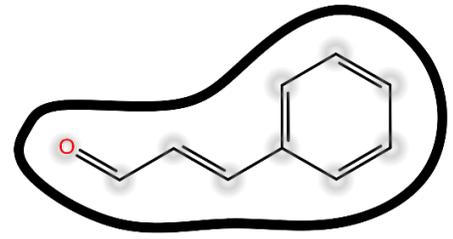
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10.

Cinnamaldehyde



† indicates protomers