On Bisphosphonates and COVID-19: *In silico* Model Suggests Inhibition of SARS-CoV-2 RdRp as Potential Explanation

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

|  |  |  |
| --- | --- | --- |
| **S. no** | **Time-line plot** | **2D interaction diagram** |
| 1. | CHEMBL164344 |  |
| 2. | CHEMBL196676 |  |
| 3. | CHEMBL387132 |  |
| 4. | CHEMBL4291724 |  |
| 5. | CHEMBL608526† |  |
| 6. | CHEMBL4569308† |  |
| 7. | CHEMBL98211† |  |
| 8. | CHEMBL164344† |  |
| 9. | Remdesivir |  |
| 10. | Cinnamaldehyde |  |

† indicates protomers