**Table 1. Information of Molecular Docking**

|  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- |
| Ingredient | Ingredient Structure | Target | PDB ID | Target Structure | Affinity (kcal/mol) |
| Tanshinone IIA |  | ATM | 6HKA |  | -6.4 |
| GADD45 | 2KG4 |  | -7.2 |
| ORC | 5UJ8 |  | -8.3 |