**Table S9. PDB ligands supporting the benzamide prediction for DAPK1**

|  |  |  |  |  |  |
| --- | --- | --- | --- | --- | --- |
| **Ligand ID** | **Chemical Structure** | **# PDB**  **Occurrence** | **Ligand ID** | **Chemical Structure** | **# PDB**  **Occurrence** |
| 0CE |  | 1 | SKE |  | 1 |
| 609 |  | 1 | STU |  | 46 |
| KSA |  | 3 |  |  |  |

Benzamide substructure is shown in pink.