BindForge Docking Analysis Report

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Protein Information

Parameter	Value
Name	Target Protein
PDB ID	Unknown
Number of Cysteines	0

Identified Cysteines

No cysteine residues identified

Molecule Information

Parameter	Value
Name	Phenylacetic acid
SMILES	C1=CC=C(C=C1)CC(=O)O

Docking Results

Summary

Metric	Value
Best Binding Affinity	-10 kcal/mol
Covalent Binding Potential	N/A
Warhead Distance	None Å

Docking Poses

Pose	Affinity (kcal/mol)	RMSD LB	RMSD UB
1	-10	0	0

Analysis and Recommendations

No specific recommendations available.