



wwPDB NMR Structure Validation Summary Report ⓘ

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PDB ID : 2JYL / pdb_00002jyl
Title : Solution Structure of A Double Mutant of The Carboxy-terminal Dimerization Domain of The HIV-1 Capsid Protein
Authors : Wong, H.C.; Shin, R.; Krishna, N.R.
Deposited on : 2007-12-14

This is a wwPDB NMR Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/NMRValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
wwPDB-RCI	:	v_1n_11_5_13_A (Berjanski et al., 2005)
PANAV	:	Wang et al. (2010)
wwPDB-ShiftChecker	:	v1.2
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

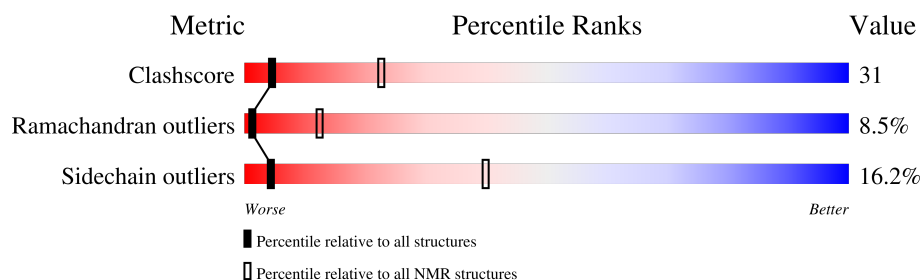
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

SOLUTION NMR

The overall completeness of chemical shifts assignment was not calculated.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	NMR archive (#Entries)
Clashscore	229148	14424
Ramachandran outliers	224038	12848
Sidechain outliers	223484	12823

The table below summarises the geometric issues observed across the polymeric chains and their fit to the experimental data. The red, orange, yellow and green segments indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A cyan segment indicates the fraction of residues that are not part of the well-defined cores, and a grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	105	

2 Ensemble composition and analysis ⓘ

This entry contains 1 models. Identification of well-defined residues and clustering analysis are not possible.

3 Entry composition

There is only 1 type of molecule in this entry. The entry contains 1301 atoms, of which 661 are hydrogens and 0 are deuteriums.

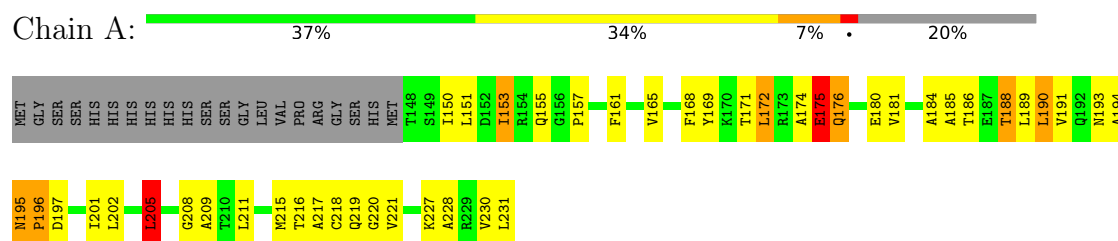
- Molecule 1 is a protein called Capsid protein p24 (CA).

Mol	Chain	Residues	Atoms						Trace
1	A	84	Total	C	H	N	O	S	0
			1301	398	661	115	123	4	

There are 23 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	127	MET	-	expression tag	UNP P12497
A	128	GLY	-	expression tag	UNP P12497
A	129	SER	-	expression tag	UNP P12497
A	130	SER	-	expression tag	UNP P12497
A	131	HIS	-	expression tag	UNP P12497
A	132	HIS	-	expression tag	UNP P12497
A	133	HIS	-	expression tag	UNP P12497
A	134	HIS	-	expression tag	UNP P12497
A	135	HIS	-	expression tag	UNP P12497
A	136	HIS	-	expression tag	UNP P12497
A	137	SER	-	expression tag	UNP P12497
A	138	SER	-	expression tag	UNP P12497
A	139	GLY	-	expression tag	UNP P12497
A	140	LEU	-	expression tag	UNP P12497
A	141	VAL	-	expression tag	UNP P12497
A	142	PRO	-	expression tag	UNP P12497
A	143	ARG	-	expression tag	UNP P12497
A	144	GLY	-	expression tag	UNP P12497
A	145	SER	-	expression tag	UNP P12497
A	146	HIS	-	expression tag	UNP P12497
A	147	MET	-	expression tag	UNP P12497
A	184	ALA	TRP	engineered mutation	UNP P12497
A	185	ALA	MET	engineered mutation	UNP P12497

- Molecule 1: Capsid protein p24 (CA)



5 Refinement protocol and experimental data overview ⓘ

The models were refined using the following method: *torsion angle dynamics*.

Of the 1 calculated structures, 1 were deposited, based on the following criterion: ?.

The following table shows the software used for structure solution, optimisation and refinement.

Software name	Classification	Version
CYANA	refinement	

No chemical shift data was provided.

6 Model quality [i](#)

6.1 Standard geometry [i](#)

There are no covalent bond-length or bond-angle outliers.

There are no bond-length outliers.

There are no bond-angle outliers.

There are no chirality outliers.

There are no planarity outliers.

6.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in each chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes averaged over the ensemble.

Mol	Chain	Non-H	H(model)	H(added)	Clashes
1	A	640	661	649	40
All	All	640	661	649	40

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 31.

5 of 40 clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Clash(Å)	Distance(Å)
1:A:201:ILE:CG2	1:A:217:ALA:HB1	0.86	2.00
1:A:201:ILE:HG22	1:A:217:ALA:HB1	0.84	1.48
1:A:172:LEU:C	1:A:172:LEU:HD12	0.79	2.03
1:A:205:LEU:HD23	1:A:209:ALA:CB	0.78	2.08
1:A:191:VAL:HG22	1:A:202:LEU:HD13	0.70	1.63

6.3 Torsion angles [i](#)

6.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the backbone conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	82/105 (78%)	65 (79%)	10 (12%)	7 (9%)	1	12
All	All	82/105 (78%)	65 (79%)	10 (12%)	7 (9%)	1	12

5 of 7 Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	157	PRO
1	A	175	GLU
1	A	176	GLN
1	A	195	ASN
1	A	196	PRO

6.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all NMR entries. The Analysed column shows the number of residues for which the sidechain conformation was analysed and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	68/86 (79%)	57 (84%)	11 (16%)	4	40
All	All	68/86 (79%)	57 (84%)	11 (16%)	4	40

5 of 11 residues with a non-rotameric sidechain are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	151	LEU
1	A	153	ILE
1	A	172	LEU
1	A	175	GLU
1	A	188	THR

6.3.3 RNA ⓘ

There are no RNA molecules in this entry.

6.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.6 Ligand geometry [i](#)

There are no ligands in this entry.

6.7 Other polymers [i](#)

There are no such molecules in this entry.

6.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

7 Chemical shift validation

No chemical shift data were provided