

Full wwPDB NMR Structure Validation Report (i)

Apr 20, 2024 – 11:15 AM EDT

PDB ID : 1ESX

Title : 1H, 15N AND 13C STRUCTURE OF THE HIV-1 REGULATORY PRO-

TEIN VPR: COMPARISON WITH THE N-AND C-TERMINAL DOMAIN

STRUCTURE, (1-51)VPR AND (52-96)VPR

Authors: Wecker, K.; Morellet, N.; Bouaziz, S.; Roques, B.

Deposited on : 2000-04-11

This is a Full wwPDB NMR Structure Validation 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 (i) 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 (i)) were used in the production of this report:

MolProbity: 4.02b-467

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

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)

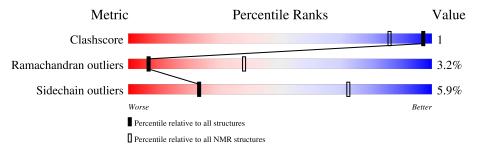
Validation Pipeline (wwPDB-VP) : 2.36.2

1 Overall quality at a glance (i)

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	NMR archive	
Metric	$(\# ext{Entries})$	$(\# ext{Entries})$	
Clashscore	158937	12864	
Ramachandran outliers	154571	11451	
Sidechain outliers	154315	11428	

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 <=5%

Mol	Chain	Length	Quality of chain				
1	A	96	78%	21%	•		



2 Ensemble composition and analysis (i)

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



3 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 1608 atoms, of which 802 are hydrogens and 0 are deuteriums.

 \bullet Molecule 1 is a protein called VPR PROTEIN.

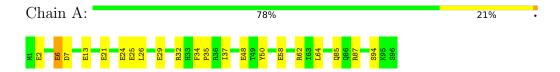
Mol	Chain	Residues	Atoms					Trace	
1	Λ	06	Total	С	Н	N	О	S	0
	A	96	1608	508	802	152	144	2	0



4 Residue-property plots (i)

These plots are provided for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic is the same as shown in the summary in section 1 of this report. The second graphic shows the sequence where residues are colour-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. Stretches of 2 or more consecutive residues without any outliers are shown as green connectors. Residues which are classified as ill-defined in the NMR ensemble, are shown in cyan with an underline colour-coded according to the previous scheme. Residues which were present in the experimental sample, but not modelled in the final structure are shown in grey.

• Molecule 1: VPR PROTEIN





5 Refinement protocol and experimental data overview (i)



The models were refined using the following method: simulated annealing, molecular dynamics, minimisations.

Of the 20 calculated structures, 1 were deposited, based on the following criterion: structures with acceptable covalent geometry, structures with favorable non-bond energy, structures with the least restraint violations, structures with the lowest energy.

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

Software name	Classification	Version	
Discover	structure solution	98.0	
Discover	refinement	98.0	

No chemical shift data was provided.



6 Model quality (i)

6.1 Standard geometry (i)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with |Z| > 5 is considered an outlier worth inspection. RMSZ is the (average) root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bo	nd lengths	Bond angles		
	Chain	RMSZ	#Z>5	RMSZ	#Z>5	
1	A	1.28	10/826 (1.2%)	1.03	1/1115 (0.1%)	
All	All	1.28	10/826 (1.2%)	1.03	1/1115 (0.1%)	

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	Chirality	Planarity
1	A	0	1
All	All	0	1

All bond outliers are listed below. They are sorted according to the Z-score.

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
1	A	29	GLU	CD-OE2	10.19	1.36	1.25
1	A	6	GLU	CD-OE2	10.17	1.36	1.25
1	A	21	GLU	CD-OE2	10.16	1.36	1.25
1	A	24	GLU	CD-OE2	10.15	1.36	1.25
1	A	58	GLU	CD-OE2	10.15	1.36	1.25
1	A	48	GLU	CD-OE2	10.14	1.36	1.25
1	A	13	GLU	CD-OE2	10.14	1.36	1.25
1	A	2	GLU	CD-OE2	10.13	1.36	1.25
1	A	25	GLU	CD-OE2	10.12	1.36	1.25
1	A	7	ASP	CG-OD2	5.04	1.36	1.25

All angle outliers are listed below.

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(^{o})$	$\operatorname{Ideal}(^{o})$
1	A	87	ARG	CB-CA-C	6.89	124.19	110.40

There are no chirality outliers.



All planar outliers are listed below.

Mol	Chain	Res	Type	Group
1	A	50	TYR	Sidechain

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	806	802	790	1
All	All	806	802	790	1

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 1.

All clashes are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\operatorname{Clash}(ext{\AA})$	$\operatorname{Distance}(\text{\AA})$
1:A:34:PHE:N	1:A:35:PRO:CD	0.49	2.75

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	94/96 (98%)	74 (79%)	17 (18%)	3 (3%)	7 38
All	All	94/96 (98%)	74 (79%)	17 (18%)	3 (3%)	7 38

All 3 Ramachandran outliers are listed below. They are sorted by the frequency of occurrence in the ensemble.

Mol	Chain	Res	Type
1	A	6	GLU
1	A	37	ILE

Continued on next page...



Continued from previous page...

Mol	Chain	Res	Type
1	A	94	SER

6.3.2 Protein sidechains (i)

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	Perce	ntiles
1	A	85/85 (100%)	80 (94%)	5 (6%)	23	72
All	All	85/85 (100%)	80 (94%)	5 (6%)	23	72

All 5 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	26	LEU
1	A	32	ARG
1	A	62	ARG
1	A	64	LEU
1	A	85	GLN

6.3.3 RNA (i)

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 monosaccharides 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 (i)

No chemical shift data were provided

