

wwPDB X-ray Structure Validation Summary Report (i)

Apr 29, 2024 – 05:57 pm BST

PDB ID : 2VX8

Title: Vamp7 longin domain Hrb peptide complex

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Deposited on : 2008-07-01

Resolution : 2.20 Å(reported)

This is a wwPDB X-ray 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/XrayValidationReportHelp
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 (1)) were used in the production of this report:

MolProbity: 4.02b-467 Xtriage (Phenix): 1.13

EDS : 2.36.2

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

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

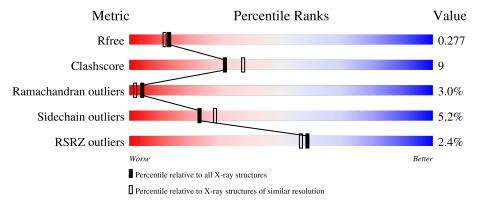
Validation Pipeline (wwPDB-VP) : 2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \ resolution} \\ (\#{\rm Entries, \ resolution \ range(\AA)}) \end{array}$
R_{free}	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain					
1	A	169	64%	12% • • 20%				
1	В	169	68%	13% • 18%				
1	С	169	5% 23%	• 22%				
1	D	169	66%	12% 21%				



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4418 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

• Molecule 1 is a protein called NUCLEOPORIN-LIKE PROTEIN RIP, VESICLE-ASSOCIATED MEMBRANE PROTEIN 7.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Λ	136	Total	С	N	О	S	0	0	0
1	A	130	1065	682	182	196	5	0	0	0
1	В	138	Total	С	N	О	S	0	0	0
1	Ъ	130	1087	696	185	201	5	0		
1	C	132	Total	С	N	О	S	0	0	0
1		132	1046	671	177	193	5	U	0	0
1	D	194	Total	С	N	О	S	0	0	0
1		D 134		674	176	192	5	U	0	

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP P70280
A	2	SER	-	expression tag	UNP P70280
A	43	GLY	-	linker	UNP P70280
A	164	HIS	-	expression tag	UNP P70280
A	165	HIS	-	expression tag	UNP P70280
A	166	HIS	-	expression tag	UNP P70280
A	167	HIS	-	expression tag	UNP P70280
A	168	HIS	-	expression tag	UNP P70280
A	169	HIS	-	expression tag	UNP P70280
В	1	GLY	-	expression tag	UNP P70280
В	2	SER	-	expression tag	UNP P70280
В	43	GLY	-	linker	UNP P70280
В	164	HIS	_	expression tag	UNP P70280
В	165	HIS	-	expression tag	UNP P70280
В	166	HIS	_	expression tag	UNP P70280
В	167	HIS	-	expression tag	UNP P70280
В	168	HIS	-	expression tag	UNP P70280
В	169	HIS		expression tag	UNP P70280
С	1	GLY	-	expression tag	UNP P70280
С	2	SER	-	expression tag	UNP P70280

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Chain	Residue	Modelled	Actual	Comment	Reference
С	43	GLY	=	linker	UNP P70280
С	164	HIS	=	expression tag	UNP P70280
С	165	HIS	-	expression tag	UNP P70280
С	166	HIS	-	expression tag	UNP P70280
С	167	HIS	-	expression tag	UNP P70280
С	168	HIS	-	expression tag	UNP P70280
С	169	HIS	-	expression tag	UNP P70280
D	1	GLY	-	expression tag	UNP P70280
D	2	SER	-	expression tag	UNP P70280
D	43	GLY	-	linker	UNP P70280
D	164	HIS	-	expression tag	UNP P70280
D	165	HIS	-	expression tag	UNP P70280
D	166	HIS	-	expression tag	UNP P70280
D	167	HIS	-	expression tag	UNP P70280
D	168	HIS	=	expression tag	UNP P70280
D	169	HIS	-	expression tag	UNP P70280

• Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	В	2	Total Cl 2 2	0	0
2	D	1	Total Cl 1 1	0	0

• Molecule 3 is water.

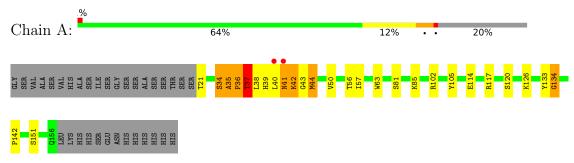
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	55	Total O 55 55	0	0
3	В	49	Total O 49 49	0	0
3	С	22	Total O 22 22	0	0
3	D	44	Total O 44 44	0	0



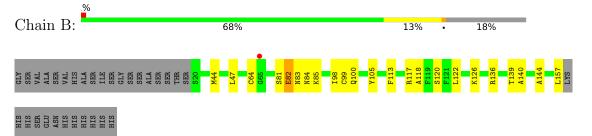
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-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. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

 \bullet Molecule 1: NUCLEOPORIN-LIKE PROTEIN RIP, VESICLE-ASSOCIATED MEMBRANE PROTEIN 7



• Molecule 1: NUCLEOPORIN-LIKE PROTEIN RIP, VESICLE-ASSOCIATED MEMBRANE PROTEIN 7



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 \bullet Molecule 1: NUCLEOPORIN-LIKE PROTEIN RIP, VESICLE-ASSOCIATED MEMBRANE PROTEIN 7









4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants	105.51Å 115.10Å 55.44Å	Donositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	57.71 - 2.20	Depositor
rtesolution (A)	57.55 - 2.20	EDS
% Data completeness	99.0 (57.71-2.20)	Depositor
(in resolution range)	99.3 (57.55-2.20)	EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	5.38 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0035	Depositor
D D.	0.220 , 0.280	Depositor
R, R_{free}	0.217 , 0.277	DCC
R_{free} test set	1747 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	37.4	Xtriage
Anisotropy	0.586	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35, 43.6	EDS
L-test for twinning ²	$ < L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4418	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.54% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities estimated from amplitudes.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: CL

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 root-mean-square of all Z scores of the bond lengths (or angles).

Mal	Chain	Boı	nd lengths	Bond angles		
IVIOI	Mol Chain		# Z > 5	RMSZ	# Z > 5	
1	A	1.17	0/1089	1.02	3/1477 (0.2%)	
1	В	1.09	1/1111 (0.1%)	0.96	0/1505	
1	С	1.00	2/1069~(0.2%)	0.97	1/1448 (0.1%)	
1	D	1.04	2/1069 (0.2%)	0.98	1/1447 (0.1%)	
All	All	1.08	5/4338 (0.1%)	0.98	5/5877 (0.1%)	

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(ext{\AA})$
1	С	99	CYS	CB-SG	6.88	1.94	1.82
1	С	50	VAL	CB-CG1	-6.23	1.39	1.52
1	В	113	PHE	CE2-CZ	5.83	1.48	1.37
1	D	148	GLU	CG-CD	5.32	1.59	1.51
1	D	148	GLU	CB-CG	5.00	1.61	1.52

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	D	71	VAL	CB-CA-C	-7.57	97.01	111.40
1	A	102	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	A	38	LEU	CA-CB-CG	5.25	127.37	115.30
1	A	102	ARG	NE-CZ-NH1	5.09	122.84	120.30
1	С	102	ARG	NE-CZ-NH2	-5.02	117.79	120.30

There are no chirality outliers.

There are no planarity outliers.



5.2 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 the 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 within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1065	0	1053	25	0
1	В	1087	0	1084	15	0
1	С	1046	0	1038	29	0
1	D	1047	0	1036	10	0
2	В	2	0	0	0	0
2	D	1	0	0	0	0
3	A	55	0	0	4	0
3	В	49	0	0	3	0
3	С	22	0	0	4	0
3	D	44	0	0	3	0
All	All	4418	0	4211	76	0

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 9.

The worst 5 of 76 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$	
1:A:35:ALA:HB1	1:A:36:PRO:HD3	1.43	0.99	
1:A:41:ASN:HA	1:A:63:TRP:CD1	1.98	0.98	
1:B:44:MET:SD	1:B:157:LEU:HD21	2.06	0.96	
1:A:37:THR:HG21	3:A:2006:HOH:O	1.69	0.92	
1:C:44:MET:HB2	1:C:47:LEU:HD21	1.59	0.85	

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 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 X-ray entries followed by that with respect to entries of similar resolution.

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	Perce	ntiles
1	A	$134/169 \ (79\%)$	119 (89%)	8 (6%)	7 (5%)	2	0
1	В	136/169 (80%)	134 (98%)	2 (2%)	0	100	100
1	С	128/169 (76%)	108 (84%)	12 (9%)	8 (6%)	1	0
1	D	128/169 (76%)	123 (96%)	4 (3%)	1 (1%)	19	19
All	All	526/676 (78%)	484 (92%)	26 (5%)	16 (3%)	4	2

5 of 16 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	35	ALA
1	A	42	LYS
1	A	44	MET
1	С	148	GLU
1	С	149	PHE

5.3.2 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 X-ray entries followed by that with respect to entries of similar resolution.

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 Outlier		Percentiles		
1	A	114/144 (79%)	108 (95%)	6 (5%)	22	27	
1	В	118/144 (82%)	112 (95%)	6 (5%)	24	29	
1	С	114/144 (79%)	108 (95%)	6 (5%)	22	27	
1	D	112/144 (78%)	106 (95%)	6 (5%)	22	26	
All	All	458/576~(80%)	434 (95%)	24 (5%)	23	28	

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	С	116	SER
1	С	135	SER
1	С	126	LYS
1	D	31	LEU

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Mol	Chain	Res	Type	
1	В	82	GLU	

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	138	GLN
1	В	39	HIS
1	В	74	GLN
1	С	130	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains (i)

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.



5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.



6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ>2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95^{th} percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<rsrz></rsrz>	$\#\mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	136/169 (80%)	0.03	2 (1%) 73 72	21, 34, 67, 76	0
1	В	138/169 (81%)	-0.11	1 (0%) 87 86	19, 33, 56, 67	0
1	С	132/169 (78%)	0.50	8 (6%) 21 20	29, 54, 76, 80	0
1	D	134/169 (79%)	-0.03	2 (1%) 73 72	27, 42, 66, 76	0
All	All	540/676 (79%)	0.09	13 (2%) 59 56	19, 40, 72, 80	0

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res Type		RSRZ
1	С	117	ARG	5.4
1	С	113	PHE	3.2
1	D	21	THR	3.1
1	A	41	ASN	2.8
1	D	116	SER	2.7

6.2 Non-standard residues in protein, DNA, RNA chains (i)

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.



Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	CL	D	1159	1/1	0.92	0.12	58,58,58,58	0
2	CL	В	1159	1/1	0.98	0.09	29,29,29,29	0
2	CL	В	1158	1/1	0.99	0.11	28,28,28,28	0

6.5 Other polymers (i)

There are no such residues in this entry.

