



# Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 5YAX  
Title : Crystal structure of a human neutralizing antibody bound to a HBV preS1 peptide  
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Deposited on : 2017-09-02  
Resolution : 2.50 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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 ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

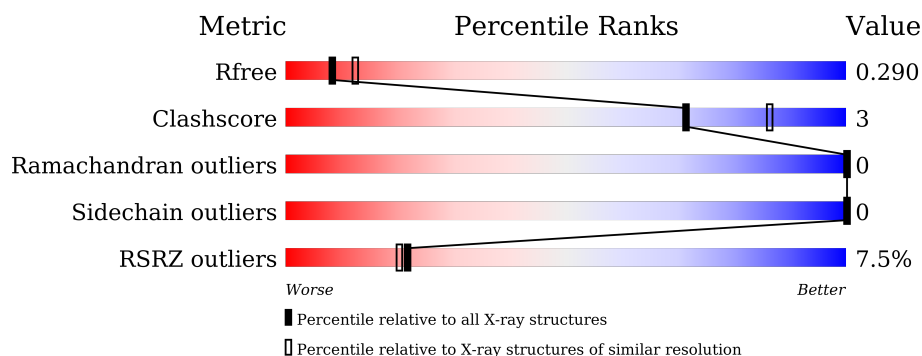
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	164625	5504 (2.50-2.50)
Clashscore	180529	6282 (2.50-2.50)
Ramachandran outliers	177936	6191 (2.50-2.50)
Sidechain outliers	177891	6193 (2.50-2.50)
RSRZ outliers	164620	5504 (2.50-2.50)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	246	<div> <div>9%</div> <div>85%</div> <div>7%</div> <div>7%</div> </div>
1	B	246	<div> <div>4%</div> <div>86%</div> <div>8%</div> <div>7%</div> </div>
2	C	60	<div> <div>5%</div> <div>7%</div> <div>7%</div> <div>87%</div> </div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3741 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 scFv1 antibody.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	0	0
			1724	1077	296	346	5			
1	B	230	Total	C	N	O	S	0	0	0
			1735	1083	298	349	5			

- Molecule 2 is a protein called Large envelope protein.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	8	Total	C	N	O	0	0	0
			54	33	9	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-11	GLY	-	expression tag	UNP B2CXZ0
C	-10	PRO	-	expression tag	UNP B2CXZ0

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	1	Total	Na	0	0
			1	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	109	Total	O	0	0
			109	109		
4	B	113	Total	O	0	0
			113	113		
4	C	5	Total	O	0	0
			5	5		



- Molecule 1: scFv1 antibody



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	142.67Å 55.66Å 67.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.92 – 2.50 19.92 – 2.50	Depositor EDS
% Data completeness (in resolution range)	94.8 (19.92-2.50) 94.5 (19.92-2.50)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.77 (at 2.50Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.242 , 0.288 0.245 , 0.290	Depositor DCC
$R_{free}$ test set	989 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	11.6	Xtriage
Anisotropy	0.452	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 39.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.86	EDS
Total number of atoms	3741	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.50% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

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

## 5 Model quality

### 5.1 Standard geometry

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

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).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.21	0/1764	0.37	0/2402
1	B	0.21	0/1775	0.39	0/2417
2	C	0.25	0/55	0.48	0/74
All	All	0.21	0/3594	0.38	0/4893

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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	1724	0	1665	10	0
1	B	1735	0	1672	12	0
2	C	54	0	43	3	0
3	C	1	0	0	0	0
4	A	109	0	0	0	0
4	B	113	0	0	0	0
4	C	5	0	0	0	0
All	All	3741	0	3380	22	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 3.

All (22) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:52:TRP:HZ2	1:B:55:ARG:HG2	1.61	0.66
1:B:21:LEU:HD12	1:B:89:LEU:HD23	1.85	0.56
1:A:45:SER:HB3	1:A:48:ARG:HB2	1.89	0.55
1:B:1037:TYR:HA	2:C:27:SER:HA	1.89	0.55
1:A:106:ARG:NH2	1:A:116:ASP:OD2	2.36	0.51
1:A:1096:ALA:HA	1:A:1126:VAL:HG11	1.91	0.51
1:A:52:TRP:CG	1:A:1116:VAL:HB	2.47	0.49
2:C:23:PHE:HB2	2:C:25:ALA:H	1.81	0.46
1:B:1019:VAL:HG21	1:B:1094:LEU:HD22	1.97	0.45
1:A:1012:SER:HB2	1:A:1127:LEU:HD11	1.99	0.45
1:B:1002:SER:HA	1:B:1113:LEU:HD11	1.99	0.45
1:A:1002:SER:N	1:A:1117:ILE:HG13	2.32	0.44
1:B:59:ARG:C	1:B:62:LYS:H	2.21	0.44
1:A:41:TRP:CE2	1:A:89:LEU:HB2	2.52	0.43
1:A:1094:LEU:HD21	1:A:1126:VAL:HG22	2.00	0.43
1:B:28:ASP:OD1	1:B:29:SER:N	2.50	0.43
1:B:55:ARG:NH2	2:C:20:ASP:OD2	2.52	0.43
1:B:1041:TRP:CD2	1:B:1089:LEU:HB2	2.54	0.41
1:B:41:TRP:CE2	1:B:89:LEU:HB2	2.55	0.41
1:A:1005:THR:HG21	1:B:127:SER:HB3	2.03	0.41
1:A:68:ALA:HB3	1:A:71:VAL:HG22	2.01	0.41
1:B:34:SER:HB3	1:B:37:ALA:HB3	2.02	0.40

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	Percentiles	
1	A	224/246 (91%)	213 (95%)	11 (5%)	0	100	100
1	B	226/246 (92%)	218 (96%)	8 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	C	6/60 (10%)	6 (100%)	0	0	100	100
All	All	456/552 (83%)	437 (96%)	19 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 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 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	Outliers	Percentiles	
1	A	192/198 (97%)	192 (100%)	0	100	100
1	B	193/198 (98%)	193 (100%)	0	100	100
2	C	5/49 (10%)	5 (100%)	0	100	100
All	All	390/445 (88%)	390 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.



## 5.6 Ligand geometry

Of 1 ligands modelled in this entry, 1 is 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

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

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<sup>th</sup> 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>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	228/246 (92%)	0.87	21 (9%) 16 15	7, 15, 30, 44	0
1	B	230/246 (93%)	0.71	11 (4%) 36 34	5, 13, 25, 47	0
2	C	8/60 (13%)	2.19	3 (37%) 1 1	16, 23, 26, 28	0
All	All	466/552 (84%)	0.81	35 (7%) 22 20	5, 14, 29, 47	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1003	VAL	5.0
1	B	1097	GLU	4.4
1	A	1037	TYR	4.2
2	C	20	ASP	4.0
1	B	128	SER	3.9
1	A	16	SER	3.9
1	A	1002	SER	3.8
1	A	1109	SER	3.3
1	B	1	GLN	3.3
1	A	120	GLN	3.1
2	C	24	GLY	2.9
1	B	5	GLN	2.8
1	A	61	SER	2.7
1	B	16	SER	2.6
1	A	1107	TYR	2.6
1	A	1026	SER	2.5
1	A	1114	SER	2.5
2	C	22	ALA	2.5
1	B	1001	GLN	2.5
1	A	1	GLN	2.4
1	A	1036	SER	2.4
1	A	1039	VAL	2.4
1	B	23	CYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	1117	ILE	2.3
1	A	1108	ASP	2.3
1	A	77	THR	2.2
1	B	3	GLN	2.2
1	A	1110	SER	2.2
1	A	23	CYS	2.2
1	A	1072	PRO	2.2
1	A	31	SER	2.2
1	B	97	GLU	2.1
1	B	109	MET	2.1
1	A	97	GLU	2.1
1	B	1095	GLN	2.1

## 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<sup>th</sup> 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	B-factors(Å <sup>2</sup> )	Q<0.9
3	NA	C	101	1/1	0.99	0.15	20,20,20,20	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.