



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 26, 2021 – 06:07 PM JST

PDB ID : 7VMU
Title : Crystal Structure of SARS-CoV Spike Receptor-Binding Domain Complexed with Neutralizing Antibody
Authors : Guo, Y.; Wang, W.; Jiao, P.; Yang, H.; Rao, Z.; Cheng, G.
Deposited on : 2021-10-09
Resolution : 2.89 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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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	:	2.23.2
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)	:	2.23.2

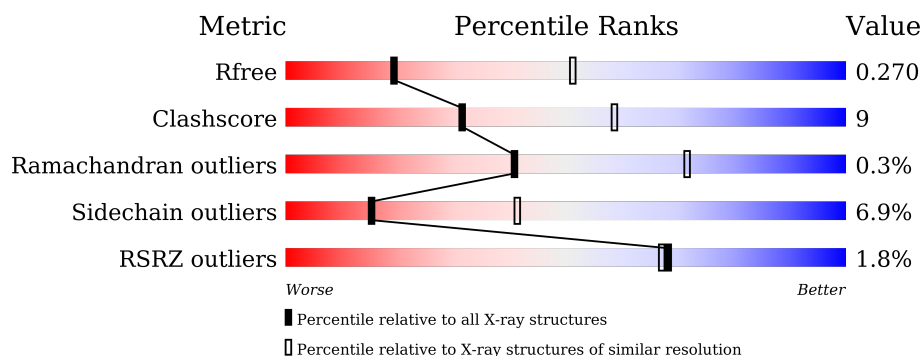
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.89 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 $\leq 5\%$. 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	245	
2	B	181	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3064 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.

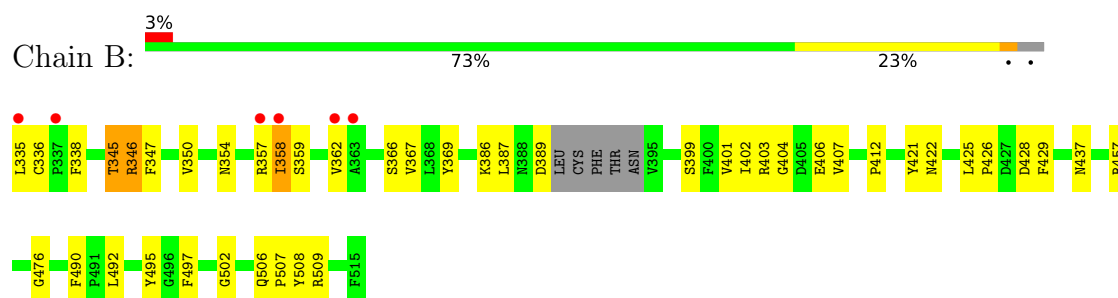
- Molecule 1 is a protein called scFv E4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	220	Total	C	N	O	S	0	0	0
			1659	1039	285	327	8			

- Molecule 2 is a protein called Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	176	Total	C	N	O	S	0	0	0
			1405	902	234	263	6			

- Molecule 1: scFv E4



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 2 2	Depositor
Cell constants a, b, c, α , β , γ	97.02Å 97.02Å 93.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.39 – 2.89 43.39 – 2.89	Depositor EDS
% Data completeness (in resolution range)	90.7 (43.39-2.89) 90.7 (43.39-2.89)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.241 , 0.273 0.244 , 0.270	Depositor DCC
R_{free} test set	497 reflections (5.24%)	wwPDB-VP
Wilson B-factor (Å ²)	56.3	Xtriage
Anisotropy	0.143	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 31.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.019 for -h,-l,-k 0.013 for l,-k,h	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	3064	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.49% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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

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	0.47	0/1693	0.70	0/2298
2	B	0.66	0/1444	0.73	0/1961
All	All	0.57	0/3137	0.71	0/4259

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	1659	0	1622	36	0
2	B	1405	0	1327	19	0
All	All	3064	0	2949	54	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.

All (54) 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:A:35:SER:HB3	1:A:50:LEU:HD23	1.56	0.88
1:A:82:MET:HB3	1:A:85:LEU:HD11	1.58	0.85

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:59:TYR:HB2	1:A:64:LYS:HG2	1.71	0.72
1:A:178:LYS:HG2	1:A:223:ALA:HB2	1.79	0.64
1:A:161:THR:HG22	1:A:211:THR:HG22	1.81	0.63
1:A:4:LEU:HD13	1:A:107:VAL:HG12	1.81	0.62
2:B:490:PHE:CZ	2:B:492:LEU:HB2	2.35	0.62
2:B:425:LEU:HD12	2:B:426:PRO:HD2	1.83	0.60
1:A:61:ASP:HA	1:A:64:LYS:HE2	1.85	0.59
2:B:403:ARG:HB3	2:B:497:PHE:HE1	1.67	0.59
2:B:338:PHE:CE1	2:B:358:ILE:HD12	2.38	0.57
1:A:2:VAL:HG22	1:A:26:GLY:HA3	1.86	0.56
1:A:35:SER:HB3	1:A:50:LEU:CD2	2.33	0.56
1:A:228:GLN:HE21	1:A:234:PRO:HB3	1.71	0.56
2:B:345:THR:HG22	2:B:346:ARG:HH21	1.70	0.56
1:A:34:MET:HB3	1:A:78:LEU:HD22	1.88	0.54
1:A:36:TRP:CE2	1:A:80:LEU:HB2	2.44	0.52
2:B:412:PRO:HG3	2:B:429:PHE:HB3	1.90	0.52
1:A:39:GLN:HB2	1:A:45:LEU:HD23	1.93	0.51
2:B:402:ILE:HD12	2:B:406:GLU:HG3	1.93	0.51
2:B:338:PHE:HE1	2:B:358:ILE:HD12	1.74	0.51
2:B:403:ARG:HB3	2:B:497:PHE:CE1	2.45	0.51
1:A:218:GLN:HG3	1:A:219:PRO:HD2	1.93	0.49
2:B:404:GLY:O	2:B:407:VAL:HG12	2.13	0.49
2:B:497:PHE:CD2	2:B:507:PRO:HB3	2.48	0.48
1:A:172:LEU:HD13	1:A:210:PHE:CD1	2.49	0.48
2:B:401:VAL:HG22	2:B:509:ARG:HG2	1.95	0.47
1:A:106:ASP:OD1	1:A:107:VAL:HG23	2.15	0.46
2:B:421:TYR:CD1	2:B:457:ARG:HB3	2.50	0.46
1:A:234:PRO:O	1:A:235:ARG:HD3	2.16	0.46
1:A:106:ASP:CG	1:A:107:VAL:HG23	2.36	0.45
2:B:350:VAL:HG22	2:B:422:ASN:HB3	1.97	0.45
1:A:9:GLY:O	1:A:18:LEU:HD11	2.17	0.45
1:A:157:ARG:HH11	1:A:157:ARG:HG2	1.82	0.45
2:B:502:GLY:O	2:B:506:GLN:HG3	2.17	0.44
1:A:6:GLN:HB2	1:A:21:SER:O	2.17	0.44
1:A:221:ASP:O	1:A:225:TYR:OH	2.29	0.44
1:A:6:GLN:HA	1:A:22:CYS:HA	1.99	0.44
1:A:69:ILE:HD11	1:A:78:LEU:HD11	2.00	0.44
1:A:33:TYR:HB2	1:A:98:ASP:O	2.18	0.42
1:A:82:MET:CB	1:A:85:LEU:HD11	2.40	0.42
2:B:347:PHE:CE2	2:B:399:SER:HB2	2.55	0.41
1:A:222:PHE:CE1	1:A:244:ILE:HG23	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:VAL:HA	2:B:476:GLY:HA2	2.03	0.41
1:A:32:ASN:ND2	1:A:97:ARG:HH11	2.18	0.41
2:B:437:ASN:HA	2:B:508:TYR:CD1	2.56	0.41
1:A:200:ARG:NE	1:A:221:ASP:OD2	2.53	0.41
1:A:60:ALA:O	1:A:64:LYS:HG3	2.21	0.40
1:A:164:ALA:HB3	1:A:208:THR:HA	2.04	0.40
1:A:155:GLY:N	1:A:217:LEU:O	2.36	0.40
1:A:175:TYR:O	1:A:225:TYR:HA	2.22	0.40
1:A:80:LEU:HD12	1:A:80:LEU:HA	1.95	0.40
1:A:86:ARG:C	1:A:116:VAL:HG11	2.42	0.40
2:B:425:LEU:HD11	2:B:429:PHE:CE1	2.57	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	216/245 (88%)	197 (91%)	18 (8%)	1 (0%)	29	61
2	B	172/181 (95%)	156 (91%)	16 (9%)	0	100	100
All	All	388/426 (91%)	353 (91%)	34 (9%)	1 (0%)	41	71

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	103	GLY

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	Outliers	Percentiles	
1	A	182/190 (96%)	176 (97%)	6 (3%)	38	72
2	B	152/157 (97%)	135 (89%)	17 (11%)	6	18
All	All	334/347 (96%)	311 (93%)	23 (7%)	15	41

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	101	GLU
1	A	102	LYS
1	A	105	MET
1	A	148	SER
1	A	149	SER
1	A	150	LEU
2	B	335	LEU
2	B	336	CYS
2	B	345	THR
2	B	346	ARG
2	B	354	ASN
2	B	357	ARG
2	B	358	ILE
2	B	359	SER
2	B	362	VAL
2	B	366	SER
2	B	367	VAL
2	B	369	TYR
2	B	386	LYS
2	B	387	LEU
2	B	389	ASP
2	B	428	ASP
2	B	495	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	32	ASN
1	A	166	GLN
1	A	177	GLN
1	A	229	GLN

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Mol	Chain	Res	Type
2	B	388	ASN
2	B	498	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](#)

There are no ligands in this entry.

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, 95th 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 ²)	Q<0.9
1	A	220/245 (89%)	0.07	1 (0%) 91 91	31, 45, 64, 75	0
2	B	176/181 (97%)	0.03	6 (3%) 45 40	30, 47, 82, 92	0
All	All	396/426 (92%)	0.05	7 (1%) 68 67	30, 46, 75, 92	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	362	VAL	5.0
2	B	335	LEU	2.9
2	B	357	ARG	2.7
2	B	363	ALA	2.7
2	B	358	ILE	2.4
2	B	337	PRO	2.1
1	A	18	LEU	2.0

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.