



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 02:23 PM UTC

PDB ID : 7EK0 / pdb_00007ek0
Title : Complex Structure of antibody BD-503 and RBD-N501Y of COVID-19
Authors : Xu, H.; Wang, B.; Zhao, T.N.; Su, X.D.
Deposited on : 2021-04-03
Resolution : 2.70 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 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
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

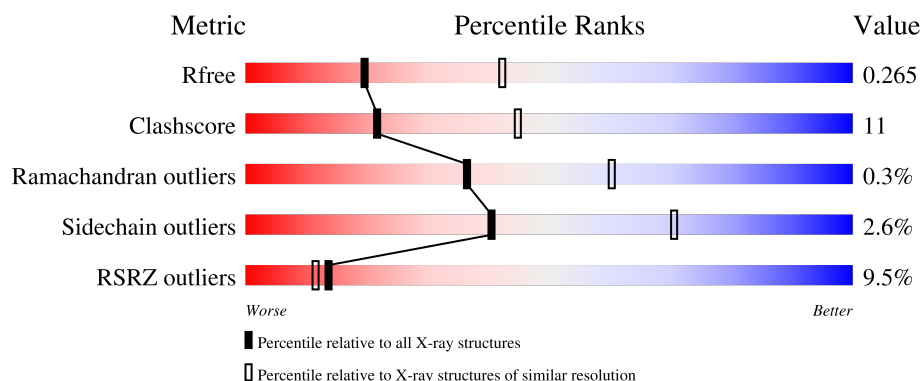
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	R	223	19% 63% 21% • 13%
2	H	222	3% 77% 18% • •
3	L	215	5% 80% 17% •

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4747 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 Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	R	193	Total	C	N	O	S	0	0	0
			1533	985	254	286	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
R	501	TYR	ASN	engineered mutation	UNP P0DTC2

- Molecule 2 is a protein called Heavy Chain of BD-503.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	214	Total	C	N	O	S	0	0	0
			1574	990	265	313	6			

- Molecule 3 is a protein called Light Chain of BD-503.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	214	Total	C	N	O	S	0	0	0
			1639	1027	272	335	5			

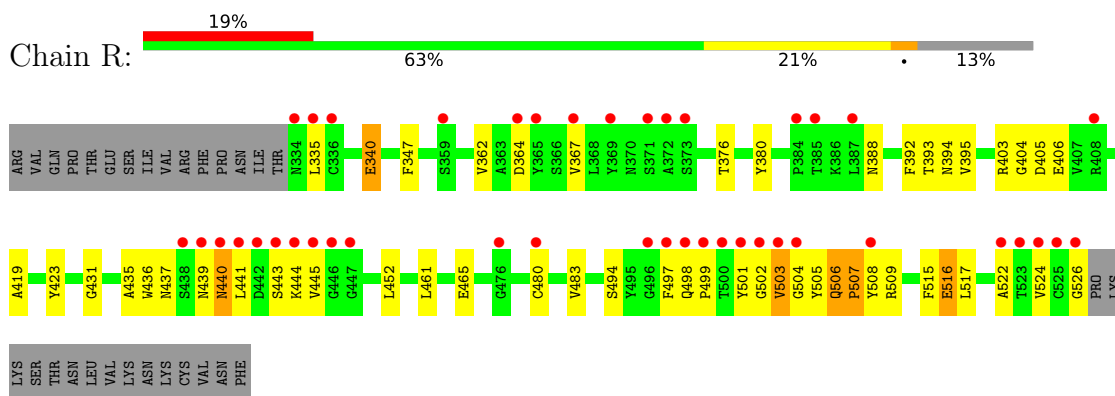
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	H	1	Total	O	0	0
			1	1		

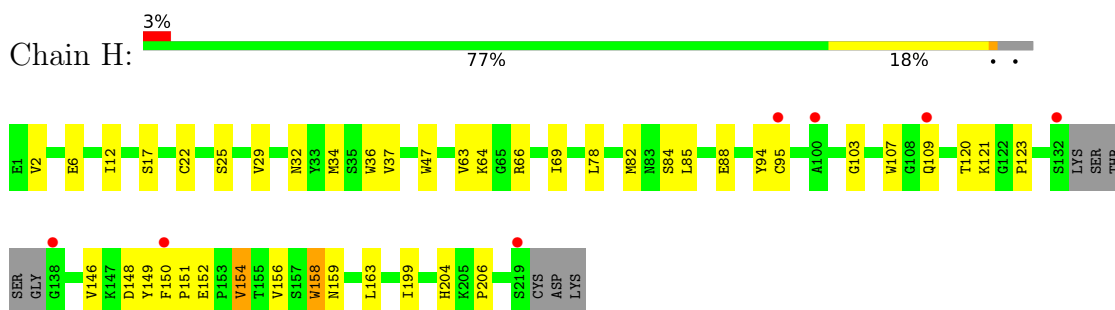
3 Residue-property plots [i](#)

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.

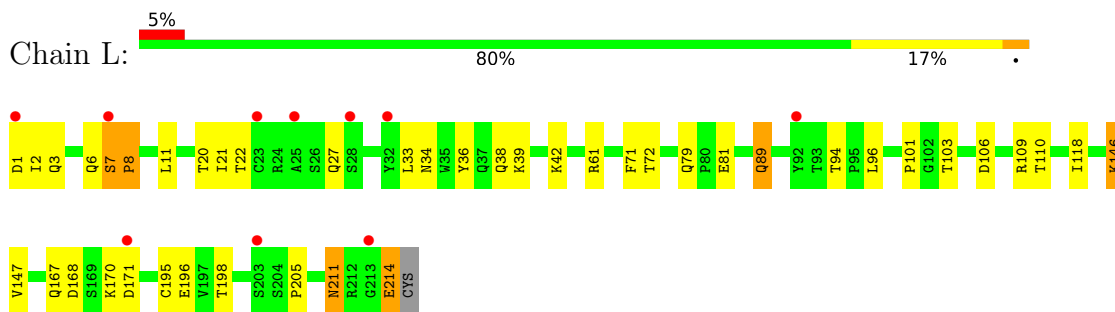
- Molecule 1: Spike protein S1



- Molecule 2: Heavy Chain of BD-503



- Molecule 3: Light Chain of BD-503



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	81.43Å 149.66Å 146.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.54 – 2.70 42.54 – 2.70	Depositor EDS
% Data completeness (in resolution range)	94.7 (42.54-2.70) 94.7 (42.54-2.70)	Depositor EDS
R_{merge}	0.23	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.80 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.214 , 0.266 0.216 , 0.265	Depositor DCC
R_{free} test set	1987 reflections (7.98%)	wwPDB-VP
Wilson B-factor (Å ²)	41.8	Xtriage
Anisotropy	0.180	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.016 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.027 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	4747	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.35% 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 [i](#)

5.1 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 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	R	0.70	1/1577 (0.1%)	0.94	3/2146 (0.1%)
2	H	0.60	1/1607 (0.1%)	0.80	4/2190 (0.2%)
3	L	0.79	6/1675 (0.4%)	0.88	1/2276 (0.0%)
All	All	0.70	8/4859 (0.2%)	0.87	8/6612 (0.1%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	R	507	PRO	C-O	-7.94	1.15	1.23
3	L	147	VAL	C-O	-6.74	1.16	1.24
3	L	211	ASN	C-O	-6.18	1.16	1.24
3	L	8	PRO	C-O	-5.80	1.16	1.24
3	L	196	GLU	C-O	-5.66	1.17	1.24
3	L	89	GLN	C-O	-5.54	1.17	1.23
2	H	156	VAL	C-O	-5.17	1.18	1.24
3	L	146	LYS	C-O	-5.02	1.18	1.24

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	506	GLN	CB-CA-C	6.83	119.40	109.09
2	H	148	ASP	N-CA-C	6.17	120.07	111.74
1	R	406	GLU	N-CA-C	5.50	119.83	113.18
2	H	64	LYS	CA-C-N	-5.39	117.58	123.30
2	H	64	LYS	C-N-CA	-5.39	117.58	123.30
1	R	505	TYR	N-CA-C	-5.18	105.00	113.19
2	H	158	TRP	N-CA-C	5.17	116.82	108.34
3	L	170	LYS	CB-CG-CD	5.06	122.94	111.30

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	R	1533	0	1449	44	0
2	H	1574	0	1555	32	0
3	L	1639	0	1595	28	0
4	H	1	0	0	0	0
All	All	4747	0	4599	101	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 11.

All (101) 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)
3:L:7:SER:HB3	3:L:8:PRO:CD	1.79	1.13
3:L:7:SER:HB2	3:L:22:THR:OG1	1.51	1.09
3:L:39:LYS:H	3:L:42:LYS:HE3	1.20	1.06
3:L:7:SER:HB3	3:L:8:PRO:HD3	1.06	1.04
1:R:443:SER:HB2	1:R:498:GLN:O	1.61	1.01
3:L:7:SER:CB	3:L:8:PRO:HD3	1.99	0.86
3:L:7:SER:CB	3:L:22:THR:OG1	2.31	0.78
1:R:393:THR:HA	1:R:522:ALA:HA	1.66	0.77
1:R:395:VAL:HG23	1:R:524:VAL:HG21	1.70	0.73
1:R:394:ASN:HB2	1:R:516:GLU:OE2	1.90	0.70
2:H:120:THR:HA	2:H:150:PHE:O	1.92	0.69
1:R:443:SER:CB	1:R:498:GLN:O	2.38	0.69
3:L:39:LYS:N	3:L:42:LYS:HE3	2.03	0.69
1:R:440:ASN:OD1	1:R:440:ASN:N	2.21	0.68
2:H:29:VAL:HG13	2:H:34:MET:HG3	1.75	0.67
1:R:440:ASN:HB2	1:R:441:LEU:HD12	1.81	0.63
2:H:159:ASN:HD21	2:H:199:ILE:H	1.48	0.62
1:R:340:GLU:O	1:R:340:GLU:HG3	1.98	0.61
1:R:392:PHE:CE1	1:R:515:PHE:HB3	2.36	0.61
1:R:498:GLN:HB3	1:R:499:PRO:CD	2.31	0.60
1:R:461:LEU:HD22	1:R:465:GLU:HB3	1.83	0.60
3:L:168:ASP:HB3	3:L:171:ASP:HB2	1.84	0.60
2:H:12:ILE:HG21	2:H:85:LEU:HD13	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:364:ASP:O	1:R:367:VAL:HG22	2.03	0.58
2:H:2:VAL:HA	2:H:25:SER:O	2.03	0.58
2:H:154:VAL:CG2	2:H:204:HIS:HB2	2.35	0.57
3:L:198:THR:HG22	3:L:205:PRO:HB3	1.86	0.57
3:L:6:GLN:HG3	3:L:101:PRO:HD2	1.87	0.57
2:H:204:HIS:CD2	2:H:206:PRO:HD2	2.40	0.56
3:L:39:LYS:NZ	3:L:81:GLU:O	2.38	0.56
1:R:443:SER:HA	1:R:497:PHE:O	2.06	0.55
2:H:159:ASN:ND2	2:H:199:ILE:H	2.03	0.55
2:H:123:PRO:HB3	2:H:149:TYR:HB3	1.90	0.54
3:L:109:ARG:NH1	3:L:110:THR:O	2.39	0.54
1:R:439:ASN:OD1	1:R:499:PRO:HA	2.07	0.54
2:H:32:ASN:HB3	2:H:34:MET:HE3	1.89	0.54
1:R:501:TYR:C	1:R:506:GLN:HE21	2.17	0.53
1:R:431:GLY:HA2	1:R:515:PHE:CD2	2.42	0.53
1:R:395:VAL:CG2	1:R:524:VAL:HG21	2.38	0.53
1:R:498:GLN:H	1:R:501:TYR:HE2	1.55	0.53
3:L:106:ASP:HB2	3:L:167:GLN:OE1	2.09	0.53
2:H:121:LYS:N	2:H:150:PHE:O	2.39	0.52
3:L:39:LYS:H	3:L:42:LYS:CE	2.07	0.52
3:L:33:LEU:HD13	3:L:71:PHE:CD2	2.45	0.52
2:H:2:VAL:O	2:H:2:VAL:HG13	2.11	0.51
1:R:404:GLY:HA3	1:R:504:GLY:HA2	1.92	0.50
3:L:61:ARG:CZ	3:L:79:GLN:HG3	2.41	0.50
1:R:436:TRP:C	1:R:508:TYR:HD1	2.20	0.50
1:R:437:ASN:HA	1:R:508:TYR:CD1	2.47	0.50
1:R:393:THR:HG23	1:R:517:LEU:HD12	1.93	0.50
1:R:403:ARG:HH21	1:R:405:ASP:CG	2.18	0.50
1:R:376:THR:HB	1:R:435:ALA:HB3	1.93	0.50
2:H:120:THR:HG22	2:H:151:PRO:HD3	1.92	0.50
2:H:146:VAL:HG11	2:H:154:VAL:HG11	1.94	0.49
2:H:66:ARG:NH2	2:H:84:SER:O	2.44	0.49
3:L:20:THR:HG23	3:L:72:THR:HG23	1.94	0.49
1:R:404:GLY:HA2	1:R:508:TYR:CD2	2.46	0.49
1:R:419:ALA:HA	1:R:423:TYR:O	2.12	0.49
2:H:32:ASN:CB	2:H:34:MET:HE3	2.43	0.48
2:H:36:TRP:HD1	2:H:69:ILE:HD12	1.79	0.48
2:H:6:GLU:OE2	2:H:95:CYS:N	2.45	0.48
1:R:388:ASN:O	1:R:526:GLY:HA3	2.14	0.47
1:R:364:ASP:CG	1:R:367:VAL:HG13	2.40	0.47
2:H:47:TRP:CH2	3:L:96:LEU:CD2	2.99	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:497:PHE:CD2	1:R:507:PRO:HB3	2.50	0.46
1:R:443:SER:HB2	1:R:498:GLN:C	2.39	0.46
2:H:34:MET:HB3	2:H:78:LEU:HD22	1.98	0.46
2:H:103:GLY:HA2	3:L:34:ASN:ND2	2.31	0.46
1:R:436:TRP:O	1:R:508:TYR:HD1	1.98	0.46
1:R:443:SER:CB	1:R:497:PHE:HB3	2.46	0.46
1:R:443:SER:HB3	1:R:497:PHE:HB3	1.98	0.46
3:L:118:ILE:HD12	3:L:195:CYS:HB2	1.98	0.45
2:H:82:MET:HE2	2:H:85:LEU:HD21	1.98	0.45
1:R:480:CYS:O	1:R:483:VAL:HG12	2.16	0.45
2:H:22:CYS:HB3	2:H:78:LEU:HB3	1.97	0.45
3:L:211:ASN:HB2	3:L:214:GLU:HG3	1.98	0.45
2:H:37:VAL:HG13	2:H:94:TYR:HB2	1.99	0.45
2:H:17:SER:HA	2:H:82:MET:O	2.16	0.45
3:L:21:ILE:HD12	3:L:103:THR:HG21	1.98	0.45
2:H:6:GLU:OE1	2:H:109:GLN:N	2.50	0.44
2:H:120:THR:CA	2:H:150:PHE:O	2.64	0.44
2:H:34:MET:HA	2:H:34:MET:HE2	2.00	0.43
1:R:437:ASN:HA	1:R:508:TYR:CE1	2.54	0.43
1:R:439:ASN:O	1:R:443:SER:OG	2.36	0.43
3:L:168:ASP:CB	3:L:171:ASP:HB2	2.47	0.43
3:L:36:TYR:HE1	3:L:89:GLN:HB3	1.84	0.43
2:H:158:TRP:O	2:H:163:LEU:HB3	2.19	0.42
3:L:1:ASP:OD1	3:L:2:ILE:N	2.53	0.42
1:R:452:LEU:HD23	1:R:494:SER:HA	2.01	0.42
1:R:380:TYR:N	1:R:431:GLY:O	2.42	0.42
3:L:109:ARG:HG2	3:L:110:THR:N	2.35	0.42
1:R:507:PRO:O	1:R:507:PRO:HG2	2.20	0.42
1:R:335:LEU:HB2	1:R:362:VAL:O	2.20	0.41
1:R:524:VAL:O	1:R:524:VAL:HG23	2.19	0.41
2:H:88:GLU:OE1	2:H:88:GLU:N	2.36	0.41
2:H:120:THR:CG2	2:H:151:PRO:HD3	2.50	0.41
1:R:347:PHE:CE2	1:R:509:ARG:HB3	2.56	0.41
1:R:431:GLY:HA2	1:R:515:PHE:CE2	2.56	0.41
3:L:38:GLN:HA	3:L:42:LYS:NZ	2.36	0.40
2:H:37:VAL:HG11	2:H:107:TRP:CZ3	2.56	0.40
1:R:502:GLY:HA3	3:L:27:GLN:NE2	2.36	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	R	191/223 (86%)	170 (89%)	20 (10%)	1 (0%)	24	48
2	H	210/222 (95%)	201 (96%)	9 (4%)	0	100	100
3	L	212/215 (99%)	203 (96%)	8 (4%)	1 (0%)	24	48
All	All	613/660 (93%)	574 (94%)	37 (6%)	2 (0%)	36	60

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	R	503	VAL
3	L	7	SER

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	R	166/196 (85%)	160 (96%)	6 (4%)	31	60
2	H	176/183 (96%)	173 (98%)	3 (2%)	53	79
3	L	189/190 (100%)	184 (97%)	5 (3%)	40	70
All	All	531/569 (93%)	517 (97%)	14 (3%)	40	70

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	R	340	GLU

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Mol	Chain	Res	Type
1	R	440	ASN
1	R	444	LYS
1	R	445	VAL
1	R	503	VAL
1	R	516	GLU
2	H	63	VAL
2	H	152	GLU
2	H	154	VAL
3	L	3	GLN
3	L	11	LEU
3	L	94	THR
3	L	146	LYS
3	L	214	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	R	334	ASN
1	R	414	GLN
1	R	474	GLN
1	R	506	GLN
2	H	39	GLN
2	H	159	ASN
2	H	196	GLN
3	L	38	GLN
3	L	161	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 oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

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, 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	R	193/223 (86%)	1.11	42 (21%) 2 2	29, 53, 96, 120	0
2	H	214/222 (96%)	0.19	7 (3%) 49 45	20, 38, 60, 93	0
3	L	214/215 (99%)	0.22	10 (4%) 36 33	20, 38, 59, 87	0
All	All	621/660 (94%)	0.49	59 (9%) 14 11	20, 41, 79, 120	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	R	501	TYR	5.8
1	R	503	VAL	5.3
1	R	498	GLN	5.3
1	R	372	ALA	4.7
1	R	369	TYR	4.7
1	R	526	GLY	4.2
1	R	443	SER	4.2
1	R	441	LEU	4.2
1	R	499	PRO	4.2
1	R	445	VAL	4.0
1	R	497	PHE	3.9
3	L	171	ASP	3.9
3	L	7	SER	3.8
1	R	502	GLY	3.7
3	L	213	GLY	3.7
1	R	387	LEU	3.7
1	R	525	CYS	3.6
1	R	500	THR	3.5
1	R	442	ASP	3.5
2	H	132	SER	3.4
3	L	92	TYR	3.4
1	R	438	SER	3.2
1	R	385	THR	3.2

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Mol	Chain	Res	Type	RSRZ
1	R	439	ASN	3.1
1	R	367	VAL	3.0
1	R	371	SER	2.9
1	R	504	GLY	2.9
2	H	109	GLN	2.8
1	R	365	TYR	2.8
2	H	138	GLY	2.8
3	L	203	SER	2.7
1	R	335	LEU	2.7
1	R	524	VAL	2.7
1	R	373	SER	2.6
1	R	444	LYS	2.6
1	R	440	ASN	2.6
2	H	150	PHE	2.5
3	L	23	CYS	2.5
1	R	523	THR	2.4
1	R	334	ASN	2.4
2	H	219	SER	2.4
1	R	522	ALA	2.4
1	R	408	ARG	2.4
1	R	364	ASP	2.4
1	R	447	GLY	2.4
1	R	446	GLY	2.4
2	H	95	CYS	2.3
3	L	1	ASP	2.3
1	R	476	GLY	2.3
1	R	480	CYS	2.2
1	R	336	CYS	2.2
1	R	496	GLY	2.2
3	L	28	SER	2.2
3	L	32	TYR	2.1
1	R	359	SER	2.1
3	L	25	ALA	2.1
2	H	100	ALA	2.1
1	R	508	TYR	2.1
1	R	384	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides 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.