



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:56 PM UTC

PDB ID : 4MMV / pdb_00004mmv
Title : Crystal Structure of Prefusion-stabilized RSV F Variant DS-Cav1-TriC at pH 9.5
Authors : Stewart-Jones, G.B.E.; McLellan, J.S.; Joyce, M.G.; Sastry, M.; Yang, Y.; Graham, B.S.; Kwong, P.D.
Deposited on : 2013-09-09
Resolution : 2.81 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

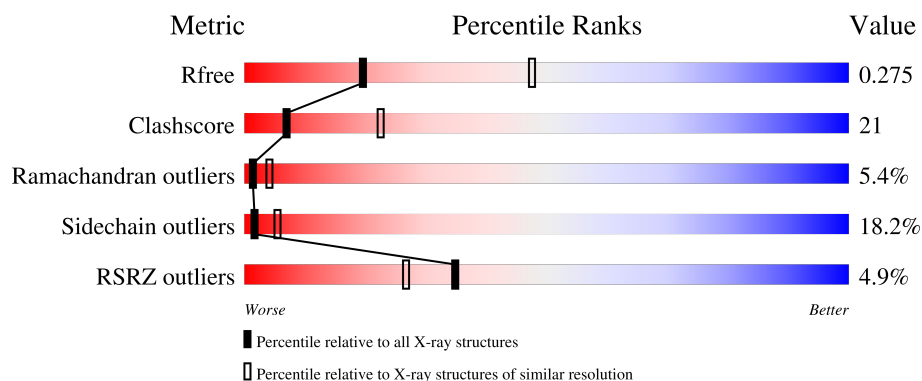
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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	82	9% 48% 37% 9% 5%
2	B	414	4% 55% 30% 11% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3844 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 Fusion glycoprotein F2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	78	Total	C	N	O	S	0	0	0
			612	385	100	124	3			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	102	ALA	PRO	engineered mutation	UNP P03420

- Molecule 2 is a protein called Fusion glycoprotein F1 fused with Fibritin trimerization domain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	408	Total	C	N	O	S	0	0	0
			3159	2003	526	610	20			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	155	CYS	SER	engineered mutation	UNP P03420
B	190	PHE	SER	engineered mutation	UNP P03420
B	207	LEU	VAL	engineered mutation	UNP P03420
B	290	CYS	SER	engineered mutation	UNP P03420
B	379	VAL	ILE	engineered mutation	UNP P03420
B	447	VAL	MET	engineered mutation	UNP P03420
B	486	HIS	ASP	engineered mutation	UNP P03420
B	487	GLN	GLU	engineered mutation	UNP P03420
B	488	TRP	PHE	engineered mutation	UNP P03420
B	489	HIS	ASP	engineered mutation	UNP P03420
B	514	SER	-	linker	UNP P03420
B	515	ALA	-	linker	UNP P03420
B	516	ILE	-	linker	UNP P03420
B	517	GLY	-	linker	UNP P03420

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Chain	Residue	Modelled	Actual	Comment	Reference
B	539	LEU	PHE	variant	UNP P10104
B	545	GLY	-	expression tag	UNP P10104
B	546	GLY	-	expression tag	UNP P10104
B	547	LEU	-	expression tag	UNP P10104
B	548	VAL	-	expression tag	UNP P10104
B	549	PRO	-	expression tag	UNP P10104
B	550	ARG	-	expression tag	UNP P10104

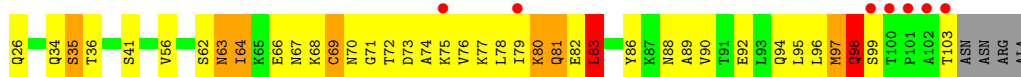
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	11	Total	O	0	0
			11	11		
3	B	62	Total	O	0	0
			62	62		

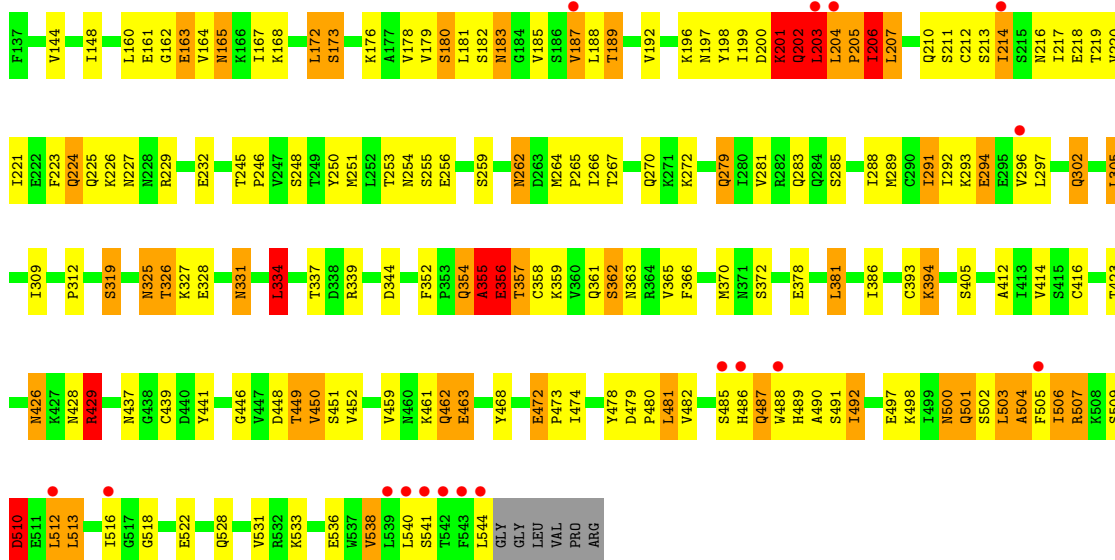
3 Residue-property plots

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: Fusion glycoprotein F2



• Molecule 2: Fusion glycoprotein F1 fused with Fibrin trimerization domain



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 3 2	Depositor
Cell constants a, b, c, α , β , γ	170.24Å 170.24Å 170.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.13 – 2.81 40.13 – 2.81	Depositor EDS
% Data completeness (in resolution range)	99.3 (40.13-2.81) 99.3 (40.13-2.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.12 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.213 , 0.279 0.208 , 0.275	Depositor DCC
R_{free} test set	1090 reflections (5.14%)	wwPDB-VP
Wilson B-factor (Å ²)	80.6	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 70.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3844	wwPDB-VP
Average B, all atoms (Å ²)	107.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.40% 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	1.15	1/619 (0.2%)	1.19	1/836 (0.1%)
2	B	1.18	7/3213 (0.2%)	1.28	20/4360 (0.5%)
All	All	1.18	8/3832 (0.2%)	1.27	21/5196 (0.4%)

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 outliers	#Planarity outliers
2	B	0	4

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	468	TYR	C-O	-6.52	1.16	1.24
2	B	334	LEU	CA-C	5.97	1.60	1.52
2	B	370	MET	C-O	-5.89	1.16	1.24
2	B	355	ALA	N-CA	5.88	1.53	1.46
1	A	34	GLN	C-O	-5.34	1.17	1.24
2	B	451	SER	CA-C	-5.14	1.46	1.52
2	B	491	SER	N-CA	-5.08	1.39	1.45
2	B	334	LEU	N-CA	-5.02	1.40	1.46

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	203	LEU	N-CA-C	-8.38	101.18	111.40
2	B	357	THR	N-CA-C	-7.71	100.58	112.99
1	A	98	GLN	N-CA-C	6.83	119.14	110.61
2	B	187	VAL	N-CA-C	6.19	115.52	106.55
2	B	288	ILE	N-CA-C	6.17	117.61	107.24
2	B	393	CYS	N-CA-C	6.12	119.21	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	291	ILE	CB-CA-C	-6.10	100.02	111.30
2	B	319	SER	CA-C-N	-6.02	113.49	119.99
2	B	319	SER	C-N-CA	-6.02	113.49	119.99
2	B	262	ASN	N-CA-C	5.91	119.68	112.23
2	B	331	ASN	N-CA-C	5.86	119.15	110.48
2	B	251	MET	N-CA-C	-5.84	104.88	112.23
2	B	326	THR	N-CA-C	5.77	118.78	111.69
2	B	429	ARG	CG-CD-NE	-5.70	99.47	112.00
2	B	196	LYS	N-CA-C	-5.70	104.07	111.02
2	B	248	SER	N-CA-C	5.63	118.15	110.55
2	B	356	GLU	CA-C-N	5.51	129.43	120.23
2	B	356	GLU	C-N-CA	5.51	129.43	120.23
2	B	429	ARG	N-CA-C	5.50	119.78	112.30
2	B	450	VAL	N-CA-C	-5.35	101.28	108.35
2	B	334	LEU	CB-CA-C	5.17	119.92	109.68

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	173	SER	Peptide
2	B	200	ASP	Peptide
2	B	201	LYS	Peptide
2	B	325	ASN	Peptide

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	612	0	620	27	0
2	B	3159	0	3189	142	1
3	A	11	0	0	1	0
3	B	62	0	0	20	0
All	All	3844	0	3809	160	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 21.

All (160) 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)
2:B:429:ARG:HB2	2:B:429:ARG:CZ	1.81	1.10
2:B:463:GLU:HB3	3:B:618:HOH:O	1.66	0.95
2:B:285:SER:HB3	3:B:651:HOH:O	1.70	0.91
2:B:416:CYS:HB3	3:B:655:HOH:O	1.76	0.83
2:B:188:LEU:C	2:B:188:LEU:HD23	2.07	0.80
2:B:173:SER:N	3:B:649:HOH:O	2.15	0.79
1:A:103:THR:HG22	2:B:148:ILE:HG12	1.66	0.76
2:B:178:VAL:HG12	2:B:178:VAL:O	1.86	0.75
2:B:302:GLN:O	3:B:651:HOH:O	2.04	0.75
1:A:97:MET:SD	1:A:98:GLN:NE2	2.60	0.75
2:B:416:CYS:SG	3:B:655:HOH:O	2.44	0.75
2:B:429:ARG:CZ	2:B:429:ARG:CB	2.63	0.70
2:B:188:LEU:HD23	2:B:189:THR:N	2.07	0.70
2:B:167:ILE:HG23	2:B:189:THR:HG21	1.74	0.70
2:B:462:GLN:HB2	3:B:609:HOH:O	1.92	0.69
2:B:279:GLN:H	2:B:279:GLN:CD	2.01	0.68
2:B:426:ASN:OD1	2:B:429:ARG:NH1	2.27	0.67
2:B:354:GLN:O	2:B:356:GLU:N	2.23	0.67
2:B:354:GLN:O	2:B:356:GLU:OE1	2.13	0.67
2:B:202:GLN:HE21	2:B:202:GLN:C	2.03	0.66
2:B:504:ALA:O	2:B:506:ILE:N	2.29	0.66
2:B:426:ASN:HB2	2:B:446:GLY:O	1.96	0.65
2:B:500:ASN:O	2:B:501:GLN:C	2.40	0.65
1:A:63:ASN:OD1	1:A:63:ASN:N	2.30	0.65
2:B:488:TRP:O	2:B:489:HIS:ND1	2.30	0.65
2:B:183:ASN:HD21	2:B:185:VAL:HG13	1.61	0.64
2:B:245:THR:HA	2:B:246:PRO:C	2.22	0.64
2:B:160:LEU:O	2:B:162:GLY:N	2.31	0.63
1:A:82:GLU:O	1:A:83:LEU:HB2	1.99	0.63
2:B:223:PHE:O	2:B:224:GLN:C	2.41	0.63
2:B:246:PRO:HB3	2:B:283:GLN:HA	1.80	0.62
2:B:188:LEU:C	2:B:188:LEU:CD2	2.73	0.61
2:B:266:ILE:HD12	2:B:270:GLN:HB3	1.83	0.60
2:B:198:TYR:HD1	3:B:628:HOH:O	1.84	0.60
2:B:223:PHE:O	2:B:226:LYS:N	2.36	0.58
2:B:497:GLU:HB3	3:B:653:HOH:O	2.03	0.58
2:B:192:VAL:HB	2:B:229:ARG:HH12	1.68	0.58
2:B:334:LEU:C	2:B:334:LEU:HD23	2.29	0.57
2:B:232:GLU:HG3	2:B:250:TYR:CE1	2.40	0.57
2:B:512:LEU:C	2:B:512:LEU:HD23	2.29	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:223:PHE:O	2:B:225:GLN:N	2.39	0.56
2:B:428:ASN:HB2	2:B:429:ARG:HH11	1.69	0.56
2:B:355:ALA:C	2:B:356:GLU:HG3	2.30	0.56
2:B:199:ILE:HA	3:B:628:HOH:O	2.05	0.56
2:B:270:GLN:HG2	2:B:309:ILE:HD12	1.88	0.56
2:B:416:CYS:CB	3:B:655:HOH:O	2.39	0.55
2:B:506:ILE:HG22	2:B:507:ARG:HG3	1.87	0.55
2:B:163:GLU:OE2	2:B:182:SER:N	2.38	0.55
2:B:253:THR:HG1	2:B:256:GLU:CD	2.15	0.54
2:B:504:ALA:C	2:B:506:ILE:H	2.15	0.54
1:A:82:GLU:O	2:B:223:PHE:HE2	1.88	0.54
1:A:35:SER:HB2	1:A:36:THR:HG23	1.89	0.54
2:B:179:VAL:HG12	2:B:180:SER:N	2.21	0.54
2:B:202:GLN:O	2:B:202:GLN:NE2	2.41	0.54
2:B:378:GLU:O	2:B:381:LEU:HB2	2.08	0.53
2:B:358:CYS:HA	2:B:366:PHE:O	2.09	0.53
2:B:293:LYS:O	2:B:294:GLU:C	2.51	0.52
2:B:497:GLU:CB	3:B:653:HOH:O	2.56	0.52
2:B:507:ARG:HA	2:B:510:ASP:HB3	1.91	0.52
2:B:264:MET:HE2	2:B:266:ILE:HD13	1.92	0.52
2:B:428:ASN:C	2:B:429:ARG:HG3	2.35	0.51
2:B:354:GLN:O	2:B:356:GLU:CD	2.54	0.51
2:B:472:GLU:O	2:B:473:PRO:C	2.53	0.51
1:A:92:GLU:O	1:A:95:LEU:N	2.44	0.51
2:B:173:SER:C	3:B:649:HOH:O	2.53	0.51
2:B:362:SER:OG	2:B:363:ASN:N	2.44	0.51
1:A:35:SER:O	2:B:474:ILE:HG23	2.11	0.50
2:B:167:ILE:CG2	2:B:189:THR:HG21	2.40	0.50
2:B:206:ILE:O	2:B:210:GLN:O	2.29	0.50
2:B:289:MET:HE1	2:B:297:LEU:HD12	1.93	0.50
2:B:203:LEU:O	2:B:207:LEU:HB2	2.11	0.50
2:B:354:GLN:C	2:B:356:GLU:CD	2.80	0.50
1:A:98:GLN:HE21	1:A:98:GLN:H	1.60	0.50
2:B:339:ARG:HD2	3:B:642:HOH:O	2.12	0.50
2:B:334:LEU:HD23	2:B:334:LEU:O	2.12	0.50
2:B:394:LYS:HD2	3:B:614:HOH:O	2.12	0.50
2:B:449:THR:HB	3:B:630:HOH:O	2.12	0.49
1:A:75:LYS:HB2	2:B:214:ILE:HG13	1.95	0.49
2:B:354:GLN:O	2:B:357:THR:HG23	2.12	0.49
2:B:281:VAL:O	2:B:285:SER:OG	2.28	0.49
2:B:179:VAL:HB	2:B:187:VAL:HG23	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:312:PRO:HD2	2:B:344:ASP:HB2	1.95	0.49
1:A:82:GLU:OE2	2:B:227:ASN:HB3	2.13	0.48
2:B:179:VAL:CG1	2:B:180:SER:N	2.76	0.48
2:B:486:HIS:CD2	2:B:487:GLN:N	2.82	0.48
2:B:500:ASN:O	2:B:502:SER:N	2.46	0.48
2:B:416:CYS:O	2:B:437:ASN:HA	2.13	0.48
1:A:88:ASN:HA	3:A:209:HOH:O	2.14	0.48
1:A:97:MET:SD	1:A:98:GLN:N	2.87	0.48
2:B:211:SER:HA	2:B:212:CYS:SG	2.53	0.48
1:A:70:ASN:N	2:B:212:CYS:SG	2.87	0.48
2:B:207:LEU:O	2:B:207:LEU:HD13	2.14	0.48
2:B:478:TYR:O	2:B:480:PRO:HD3	2.14	0.48
2:B:173:SER:CA	3:B:649:HOH:O	2.58	0.47
2:B:183:ASN:ND2	2:B:185:VAL:HG13	2.28	0.47
2:B:183:ASN:OD1	2:B:185:VAL:HG22	2.14	0.47
2:B:217:ILE:HG23	2:B:218:GLU:N	2.29	0.47
1:A:63:ASN:O	1:A:64:ILE:O	2.32	0.47
2:B:164:VAL:O	2:B:167:ILE:N	2.46	0.47
2:B:253:THR:O	2:B:254:ASN:C	2.58	0.47
2:B:488:TRP:O	2:B:489:HIS:CG	2.67	0.47
2:B:165:ASN:OD1	2:B:165:ASN:N	2.45	0.47
1:A:89:ALA:O	1:A:90:VAL:C	2.56	0.47
2:B:266:ILE:HD12	2:B:270:GLN:CB	2.45	0.47
1:A:98:GLN:CG	1:A:98:GLN:O	2.63	0.46
2:B:192:VAL:HA	3:B:627:HOH:O	2.15	0.46
2:B:448:ASP:OD2	2:B:461:LYS:CE	2.63	0.46
1:A:71:GLY:N	2:B:212:CYS:SG	2.89	0.46
2:B:325:ASN:C	2:B:325:ASN:OD1	2.58	0.46
2:B:305:LEU:HD22	3:B:638:HOH:O	2.16	0.46
2:B:386:ILE:HG12	2:B:492:ILE:HD13	1.97	0.46
1:A:72:THR:HG21	1:A:76:VAL:HG23	1.98	0.45
2:B:352:PHE:CE2	2:B:372:SER:HB3	2.52	0.45
2:B:412:ALA:O	2:B:441:TYR:HA	2.17	0.45
2:B:217:ILE:HG23	2:B:218:GLU:HG3	1.98	0.44
2:B:264:MET:HA	2:B:265:PRO:HD3	1.72	0.44
2:B:232:GLU:HG3	2:B:250:TYR:CZ	2.52	0.44
2:B:337:THR:HG21	2:B:489:HIS:CD2	2.53	0.44
2:B:279:GLN:CD	2:B:279:GLN:N	2.74	0.44
2:B:178:VAL:O	2:B:178:VAL:CG1	2.56	0.44
2:B:448:ASP:OD2	2:B:461:LYS:HE2	2.18	0.43
1:A:81:GLN:O	1:A:83:LEU:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:197:ASN:O	2:B:198:TYR:C	2.61	0.43
1:A:75:LYS:HD3	2:B:214:ILE:HD12	1.99	0.43
2:B:289:MET:HE1	2:B:297:LEU:CD1	2.49	0.43
2:B:428:ASN:HB2	2:B:429:ARG:NH1	2.34	0.43
2:B:386:ILE:HG12	2:B:492:ILE:CD1	2.49	0.43
2:B:426:ASN:OD1	2:B:426:ASN:C	2.62	0.43
1:A:56:VAL:HB	2:B:189:THR:HG23	2.00	0.43
2:B:204:LEU:CB	2:B:205:PRO:HD3	2.49	0.42
2:B:217:ILE:HD12	2:B:217:ILE:HA	1.93	0.42
2:B:487:GLN:O	2:B:488:TRP:C	2.62	0.42
2:B:414:VAL:O	2:B:439:CYS:HA	2.19	0.42
2:B:181:LEU:HD11	2:B:187:VAL:HG21	2.01	0.42
2:B:210:GLN:HA	2:B:211:SER:C	2.45	0.42
1:A:83:LEU:HA	1:A:86:TYR:HB3	2.01	0.41
2:B:204:LEU:O	2:B:206:ILE:N	2.52	0.41
2:B:405:SER:HB2	2:B:452:VAL:HG21	2.01	0.41
2:B:270:GLN:HG2	2:B:309:ILE:CD1	2.49	0.41
1:A:69:CYS:SG	1:A:72:THR:HG22	2.60	0.41
1:A:98:GLN:HE21	1:A:98:GLN:N	2.17	0.41
1:A:80:LYS:O	1:A:83:LEU:HB3	2.20	0.41
2:B:204:LEU:CB	2:B:205:PRO:CD	2.99	0.41
2:B:204:LEU:C	2:B:206:ILE:H	2.28	0.41
2:B:187:VAL:HG23	2:B:187:VAL:O	2.21	0.41
2:B:197:ASN:O	2:B:201:LYS:N	2.53	0.41
2:B:327:LYS:O	2:B:328:GLU:C	2.63	0.41
2:B:479:ASP:OD1	2:B:481:LEU:HB2	2.21	0.41
2:B:512:LEU:HD23	2:B:513:LEU:N	2.36	0.41
2:B:203:LEU:O	2:B:207:LEU:N	2.47	0.40
2:B:482:VAL:HG21	2:B:503:LEU:HD13	2.04	0.40
2:B:507:ARG:CZ	2:B:507:ARG:HB2	2.50	0.40
1:A:72:THR:OG1	1:A:73:ASP:N	2.54	0.40
2:B:497:GLU:O	2:B:498:LYS:C	2.64	0.40
2:B:220:VAL:O	2:B:221:ILE:C	2.63	0.40
2:B:533:LYS:O	2:B:536:GLU:O	2.40	0.40
2:B:198:TYR:C	2:B:198:TYR:CD1	2.99	0.40
2:B:331:ASN:HB2	3:B:616:HOH:O	2.22	0.40
2:B:487:GLN:NE2	2:B:490:ALA:HB2	2.37	0.40
2:B:536:GLU:OE2	2:B:538:VAL:HG22	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:214:ILE:O	2:B:214:ILE:O[14_444]	2.00	0.20

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	76/82 (93%)	60 (79%)	10 (13%)	6 (8%)	1	1
2	B	406/414 (98%)	326 (80%)	60 (15%)	20 (5%)	1	5
All	All	482/496 (97%)	386 (80%)	70 (14%)	26 (5%)	1	4

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	64	ILE
1	A	67	ASN
1	A	83	LEU
1	A	96	LEU
1	A	97	MET
2	B	161	GLU
2	B	172	LEU
2	B	224	GLN
2	B	355	ALA
2	B	485	SER
2	B	504	ALA
1	A	74	ALA
2	B	501	GLN
2	B	505	PHE
2	B	510	ASP
2	B	206	ILE
2	B	362	SER
2	B	462	GLN
2	B	500	ASN
2	B	541	SER

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Mol	Chain	Res	Type
2	B	202	GLN
2	B	216	ASN
2	B	540	LEU
2	B	205	PRO
2	B	294	GLU
2	B	518	GLY

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	70/73 (96%)	53 (76%)	17 (24%)	1	2
2	B	369/373 (99%)	306 (83%)	63 (17%)	2	7
All	All	439/446 (98%)	359 (82%)	80 (18%)	2	5

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	35	SER
1	A	41	SER
1	A	62	SER
1	A	63	ASN
1	A	66	GLU
1	A	68	LYS
1	A	69	CYS
1	A	77	LYS
1	A	78	LEU
1	A	79	ILE
1	A	80	LYS
1	A	81	GLN
1	A	83	LEU
1	A	94	GLN
1	A	98	GLN
1	A	99	SER

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Mol	Chain	Res	Type
2	B	144	VAL
2	B	163	GLU
2	B	165	ASN
2	B	168	LYS
2	B	172	LEU
2	B	176	LYS
2	B	180	SER
2	B	183	ASN
2	B	189	THR
2	B	201	LYS
2	B	202	GLN
2	B	203	LEU
2	B	204	LEU
2	B	206	ILE
2	B	207	LEU
2	B	213	SER
2	B	214	ILE
2	B	219	THR
2	B	255	SER
2	B	259	SER
2	B	262	ASN
2	B	267	THR
2	B	272	LYS
2	B	279	GLN
2	B	291	ILE
2	B	292	ILE
2	B	296	VAL
2	B	302	GLN
2	B	305	LEU
2	B	319	SER
2	B	326	THR
2	B	334	LEU
2	B	354	GLN
2	B	356	GLU
2	B	359	LYS
2	B	361	GLN
2	B	365	VAL
2	B	381	LEU
2	B	394	LYS
2	B	423	THR
2	B	426	ASN
2	B	429	ARG

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Mol	Chain	Res	Type
2	B	449	THR
2	B	450	VAL
2	B	459	VAL
2	B	463	GLU
2	B	472	GLU
2	B	481	LEU
2	B	487	GLN
2	B	492	ILE
2	B	503	LEU
2	B	506	ILE
2	B	507	ARG
2	B	509	SER
2	B	510	ASP
2	B	512	LEU
2	B	513	LEU
2	B	516	ILE
2	B	522	GLU
2	B	528	GLN
2	B	531	VAL
2	B	538	VAL
2	B	544	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	81	GLN
1	A	98	GLN
2	B	202	GLN
2	B	276	ASN
2	B	277	ASN
2	B	486	HIS
2	B	487	GLN
2	B	494	GLN

5.3.3 RNA ⓘ

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 [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

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	A	78/82 (95%)	0.30	7 (8%) 15 11	64, 99, 207, 237	0
2	B	408/414 (98%)	0.09	17 (4%) 40 32	55, 84, 216, 287	0
All	All	486/496 (97%)	0.13	24 (4%) 35 27	55, 87, 215, 287	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	102	ALA	7.4
2	B	544	LEU	6.2
1	A	101	PRO	5.1
2	B	540	LEU	4.5
2	B	204	LEU	4.1
2	B	488	TRP	4.0
2	B	541	SER	3.4
2	B	296	VAL	3.2
1	A	103	THR	3.1
2	B	203	LEU	2.9
2	B	187	VAL	2.9
2	B	486	HIS	2.8
2	B	539	LEU	2.7
2	B	505	PHE	2.7
2	B	542	THR	2.6
1	A	99	SER	2.6
2	B	485	SER	2.6
2	B	543	PHE	2.5
1	A	100	THR	2.4
2	B	516	ILE	2.3
1	A	75	LYS	2.2
2	B	214	ILE	2.1
1	A	79	ILE	2.1
2	B	512	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 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.