



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 15, 2026 – 12:48 AM UTC

PDB ID : 9VDI / pdb_00009vdi
Title : Crystal structure of the recombinant A1-antitrypsin F51L/M351V/M358V triple mutant
Authors : Wang, B.; Lan, T.J.; Lin, M.; Wang, Y.J.; Li, L.F.; Lin, X.L.
Deposited on : 2025-06-08
Resolution : 2.60 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

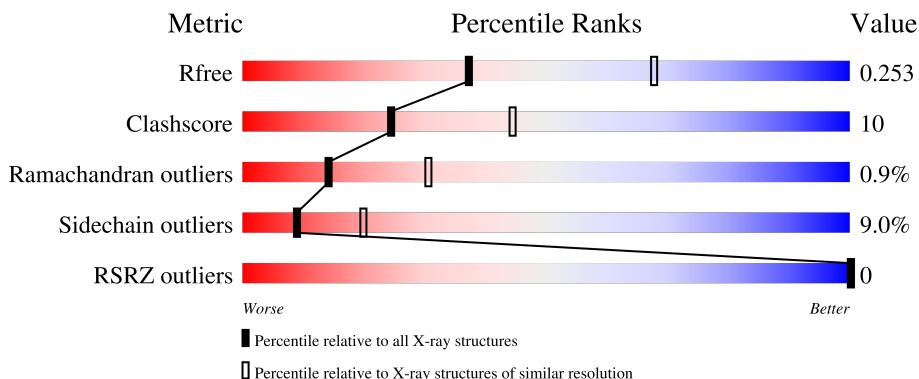
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	372	 71% 25% .
1	B	372	 71% 27% .

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5884 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 Alpha-1-antitrypsin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	372	Total	C	N	O	S	0	0	0
			2948	1899	481	560	8			
1	B	372	Total	C	N	O	S	0	0	0
			2936	1893	475	560	8			

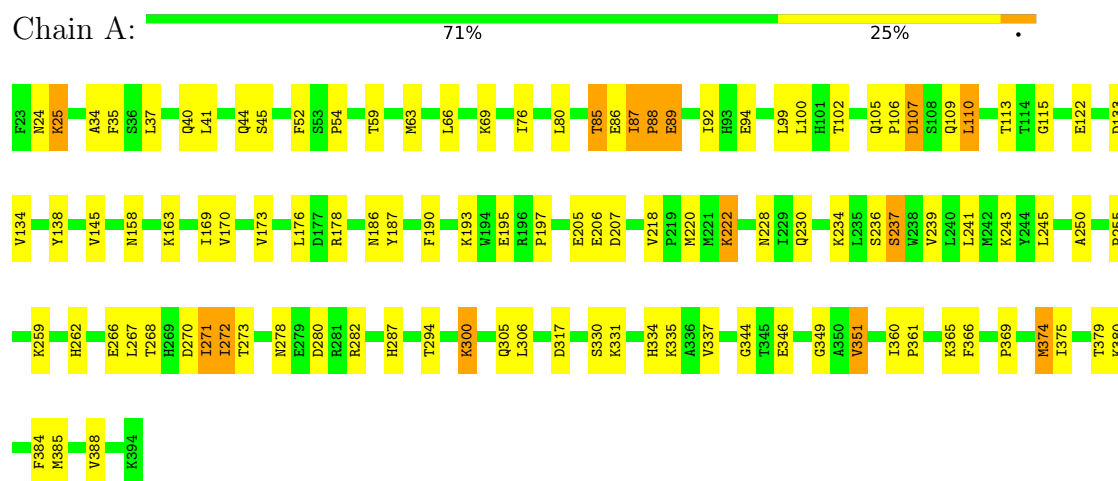
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	51	LEU	PHE	engineered mutation	UNP P01009
A	70	GLY	ALA	conflict	UNP P01009
A	101	HIS	ARG	conflict	UNP P01009
A	351	VAL	MET	engineered mutation	UNP P01009
A	358	VAL	MET	engineered mutation	UNP P01009
A	376	ASP	GLU	conflict	UNP P01009
B	51	LEU	PHE	engineered mutation	UNP P01009
B	70	GLY	ALA	conflict	UNP P01009
B	101	HIS	ARG	conflict	UNP P01009
B	351	VAL	MET	engineered mutation	UNP P01009
B	358	VAL	MET	engineered mutation	UNP P01009
B	376	ASP	GLU	conflict	UNP P01009

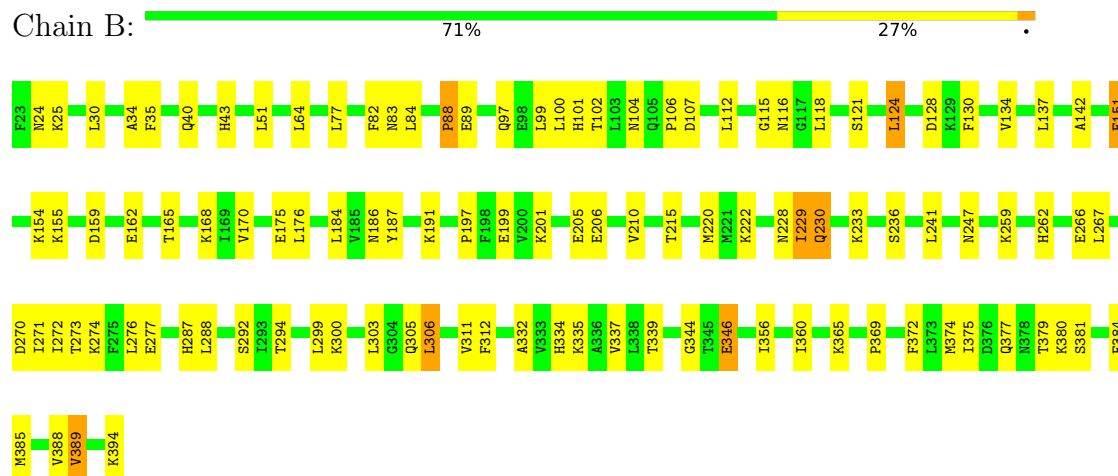
3 Residue-property plots

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: Alpha-1-antitrypsin



• Molecule 1: Alpha-1-antitrypsin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.85Å 36.37Å 115.30Å 90.00° 90.08° 90.00°	Depositor
Resolution (Å)	47.16 – 2.60 47.16 – 2.60	Depositor EDS
% Data completeness (in resolution range)	95.8 (47.16-2.60) 95.8 (47.16-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 2.61Å)	Xtriage
Refinement program	PHENIX 1.21	Depositor
R, R_{free}	0.210 , 0.254 0.214 , 0.253	Depositor DCC
R_{free} test set	2001 reflections (9.29%)	wwPDB-VP
Wilson B-factor (Å ²)	53.8	Xtriage
Anisotropy	0.493	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 41.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.457 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5884	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.32% 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	A	0.65	1/3009 (0.0%)	0.97	4/4068 (0.1%)
1	B	0.54	0/2997	0.82	2/4054 (0.0%)
All	All	0.60	1/6006 (0.0%)	0.90	6/8122 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	222	LYS	C-N	9.31	1.45	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	222	LYS	CA-C-N	-5.92	111.52	121.94
1	A	222	LYS	C-N-CA	-5.92	111.52	121.94
1	A	351	VAL	N-CA-CB	-5.57	104.71	110.95
1	A	109	GLN	N-CA-C	-5.26	106.25	113.30
1	B	389	VAL	N-CA-C	-5.16	108.47	113.53
1	B	346	GLU	N-CA-C	-5.03	103.41	110.50

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	2948	0	2967	56	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	2936	0	2945	62	0
All	All	5884	0	5912	117	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 10.

All (117) 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:158:ASN:HD21	1:A:173:VAL:H	1.21	0.86
1:B:162:GLU:HG3	1:B:170:VAL:HG12	1.58	0.83
1:A:85:THR:HG23	1:B:292:SER:HB3	1.62	0.82
1:B:154:LYS:HE2	1:B:175:GLU:HA	1.69	0.74
1:B:165:THR:O	1:B:168:LYS:HG2	1.89	0.73
1:A:237:SER:HB3	1:A:255:PRO:HA	1.74	0.69
1:B:191:LYS:HD3	1:B:339:THR:HB	1.77	0.67
1:B:276:LEU:HD11	1:B:375:ILE:HD11	1.78	0.65
1:A:115:GLY:HA3	1:A:187:TYR:CZ	2.32	0.65
1:B:372:PHE:CD1	1:B:374:MET:HE2	2.32	0.65
1:A:89:GLU:HA	1:A:92:ILE:HD12	1.80	0.64
1:B:199:GLU:HG2	1:B:201:LYS:HG2	1.80	0.64
1:A:280:ASP:HB3	1:A:282:ARG:HH12	1.63	0.63
1:A:294:THR:HG22	1:A:337:VAL:HG22	1.81	0.62
1:A:52:PHE:HE2	1:A:54:PRO:HG3	1.67	0.59
1:B:300:LYS:HD2	1:B:312:PHE:HB3	1.85	0.59
1:B:115:GLY:HA3	1:B:187:TYR:CZ	2.38	0.59
1:A:280:ASP:HB3	1:A:282:ARG:NH1	2.18	0.58
1:A:106:PRO:O	1:A:107:ASP:C	2.46	0.58
1:A:158:ASN:ND2	1:A:173:VAL:H	1.98	0.58
1:B:118:LEU:HD13	1:B:142:ALA:HB2	1.86	0.57
1:B:294:THR:HG22	1:B:337:VAL:HG22	1.87	0.56
1:A:236:SER:OG	1:A:259:LYS:HD3	2.05	0.56
1:B:35:PHE:CD1	1:B:385:MET:HE1	2.42	0.55
1:B:77:LEU:HB2	1:B:84:LEU:HD21	1.89	0.55
1:A:52:PHE:CE2	1:A:54:PRO:HG3	2.42	0.55
1:B:97:GLN:HG3	1:B:137:LEU:O	2.07	0.55
1:B:247:ASN:OD1	1:B:377:GLN:HG2	2.07	0.54
1:A:63:MET:HB2	1:A:138:TYR:CG	2.42	0.54
1:A:268:THR:HB	1:A:271:ILE:HG13	1.88	0.54
1:A:66:LEU:HD13	1:A:133:ASP:HB3	1.90	0.53
1:A:193:LYS:HD3	1:A:344:GLY:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:LEU:HG	1:A:190:PHE:HE1	1.73	0.52
1:A:287:HIS:CG	1:A:365:LYS:HG2	2.44	0.52
1:B:299:LEU:HD11	1:B:334:HIS:HB2	1.91	0.52
1:B:228:ASN:HB3	1:B:241:LEU:HB3	1.91	0.51
1:A:169:ILE:HD12	1:A:187:TYR:HB2	1.93	0.51
1:A:99:LEU:HD13	1:A:379:THR:HG21	1.93	0.51
1:B:151:GLU:CD	1:B:151:GLU:H	2.19	0.50
1:B:270:ASP:HA	1:B:273:THR:HG22	1.92	0.50
1:A:287:HIS:HB2	1:A:365:LYS:HA	1.94	0.50
1:B:34:ALA:HB1	1:B:385:MET:HE2	1.92	0.50
1:B:121:SER:HB2	1:B:124:LEU:HD11	1.94	0.50
1:A:287:HIS:CE1	1:A:365:LYS:HE3	2.48	0.49
1:B:186:ASN:O	1:B:334:HIS:HA	2.13	0.49
1:B:369:PRO:HA	1:B:388:VAL:O	2.13	0.49
1:B:379:THR:HG22	1:B:381:SER:N	2.28	0.49
1:A:250:ALA:HB1	1:A:374:MET:HE2	1.94	0.49
1:B:104:ASN:C	1:B:106:PRO:HD3	2.37	0.49
1:A:195:GLU:HG2	1:A:245:LEU:HD23	1.96	0.47
1:B:262:HIS:CE1	1:B:266:GLU:HG3	2.49	0.47
1:A:262:HIS:CE1	1:A:266:GLU:HG3	2.50	0.47
1:B:372:PHE:HD1	1:B:374:MET:HE2	1.77	0.47
1:B:184:LEU:O	1:B:332:ALA:HA	2.15	0.47
1:B:303:LEU:HA	1:B:306:LEU:HB2	1.97	0.47
1:A:99:LEU:O	1:A:102:THR:HB	2.14	0.47
1:A:99:LEU:CD1	1:A:379:THR:HG21	2.44	0.47
1:A:186:ASN:O	1:A:334:HIS:HA	2.15	0.47
1:A:250:ALA:CB	1:A:374:MET:HE2	2.45	0.47
1:B:287:HIS:CG	1:B:365:LYS:HG2	2.49	0.47
1:A:220:MET:HE1	1:A:366:PHE:O	2.15	0.46
1:B:116:ASN:ND2	1:B:186:ASN:ND2	2.63	0.46
1:B:116:ASN:ND2	1:B:186:ASN:HD22	2.13	0.46
1:B:97:GLN:HE22	1:B:101:HIS:CD2	2.33	0.46
1:B:210:VAL:HG22	1:B:389:VAL:O	2.15	0.46
1:A:59:THR:HG21	1:A:100:LEU:CD1	2.45	0.46
1:A:100:LEU:HD23	1:A:100:LEU:HA	1.65	0.46
1:A:25:LYS:HE3	1:A:25:LYS:HB3	1.62	0.46
1:B:40:GLN:HG2	1:B:305:GLN:HG3	1.97	0.45
1:B:273:THR:O	1:B:277:GLU:HG2	2.15	0.45
1:B:77:LEU:CB	1:B:84:LEU:HD21	2.46	0.45
1:A:278:ASN:OD1	1:A:280:ASP:HB2	2.16	0.45
1:B:294:THR:HG22	1:B:337:VAL:HG13	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:369:PRO:HA	1:A:388:VAL:O	2.17	0.45
1:A:59:THR:HG21	1:A:100:LEU:HG	1.98	0.45
1:A:41:LEU:HD23	1:A:41:LEU:HA	1.77	0.44
1:A:374:MET:HB2	1:A:384:PHE:HB2	1.99	0.44
1:A:270:ASP:HA	1:A:273:THR:HG22	1.98	0.44
1:B:267:LEU:HD23	1:B:271:ILE:HD11	2.00	0.44
1:B:155:LYS:O	1:B:159:ASP:HB2	2.18	0.44
1:A:228:ASN:HB3	1:A:241:LEU:HB3	2.00	0.43
1:B:64:LEU:HD21	1:B:311:VAL:HG21	2.00	0.43
1:B:197:PRO:HB3	1:B:344:GLY:N	2.32	0.43
1:B:379:THR:O	1:B:380:LYS:HB2	2.17	0.43
1:B:187:TYR:HA	1:B:335:LYS:O	2.18	0.43
1:B:220:MET:HE3	1:B:287:HIS:C	2.44	0.43
1:B:287:HIS:HB2	1:B:365:LYS:HA	2.00	0.43
1:A:86:GLU:HB3	1:A:87:ILE:H	1.73	0.43
1:A:230:GLN:HG2	1:A:239:VAL:HG22	2.01	0.43
1:A:379:THR:O	1:A:380:LYS:HB2	2.19	0.43
1:B:130:PHE:O	1:B:134:VAL:HG13	2.19	0.43
1:A:40:GLN:HG2	1:A:305:GLN:HG3	2.00	0.42
1:B:99:LEU:CD1	1:B:379:THR:HG21	2.48	0.42
1:A:76:ILE:O	1:A:80:LEU:HG	2.20	0.42
1:B:247:ASN:O	1:B:377:GLN:HB3	2.19	0.42
1:A:35:PHE:CD1	1:A:385:MET:HE1	2.55	0.42
1:B:374:MET:CG	1:B:384:PHE:HB2	2.49	0.42
1:A:34:ALA:HB1	1:A:385:MET:HE2	2.02	0.42
1:A:267:LEU:HD22	1:A:272:ILE:HD11	2.02	0.42
1:A:234:LYS:NZ	1:A:271:ILE:HD11	2.35	0.42
1:B:360:ILE:HD12	1:B:360:ILE:O	2.20	0.42
1:A:69:LYS:HB2	1:A:317:ASP:OD2	2.20	0.41
1:A:300:LYS:HE2	1:A:300:LYS:HB2	1.70	0.41
1:A:360:ILE:HA	1:A:361:PRO:HD3	1.89	0.41
1:B:77:LEU:O	1:B:82:PHE:HB2	2.21	0.41
1:A:197:PRO:HB3	1:A:344:GLY:HA3	2.03	0.41
1:A:205:GLU:HA	1:A:218:VAL:O	2.20	0.41
1:B:102:THR:HG21	1:B:379:THR:OG1	2.21	0.41
1:B:229:ILE:C	1:B:230:GLN:HG3	2.46	0.41
1:B:100:LEU:HD23	1:B:100:LEU:HA	1.77	0.41
1:B:220:MET:HE3	1:B:288:LEU:N	2.36	0.41
1:B:51:LEU:HD11	1:B:384:PHE:CD2	2.56	0.40
1:B:99:LEU:HD13	1:B:379:THR:HG21	2.04	0.40
1:B:236:SER:O	1:B:236:SER:OG	2.38	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:GLN:CG	1:B:137:LEU:O	2.69	0.40
1:A:187:TYR:HA	1:A:335:LYS:O	2.21	0.40
1:B:184:LEU:HD12	1:B:184:LEU:HA	1.90	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	370/372 (100%)	351 (95%)	15 (4%)	4 (1%)	11	25
1	B	370/372 (100%)	358 (97%)	9 (2%)	3 (1%)	16	34
All	All	740/744 (100%)	709 (96%)	24 (3%)	7 (1%)	14	30

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	88	PRO
1	A	89	GLU
1	B	89	GLU
1	A	107	ASP
1	B	88	PRO
1	B	107	ASP
1	A	349	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	329/329 (100%)	295 (90%)	34 (10%)	7	15
1	B	327/329 (99%)	302 (92%)	25 (8%)	12	27
All	All	656/658 (100%)	597 (91%)	59 (9%)	9	20

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	ASN
1	A	25	LYS
1	A	37	LEU
1	A	44	GLN
1	A	45	SER
1	A	85	THR
1	A	87	ILE
1	A	88	PRO
1	A	94	GLU
1	A	105	GLN
1	A	110	LEU
1	A	113	THR
1	A	122	GLU
1	A	134	VAL
1	A	145	VAL
1	A	163	LYS
1	A	170	VAL
1	A	176	LEU
1	A	178	ARG
1	A	206	GLU
1	A	207	ASP
1	A	222	LYS
1	A	237	SER
1	A	243	LYS
1	A	271	ILE
1	A	272	ILE
1	A	300	LYS
1	A	306	LEU
1	A	330	SER
1	A	331	LYS
1	A	346	GLU
1	A	351	VAL
1	A	374	MET
1	A	375	ILE

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Mol	Chain	Res	Type
1	B	24	ASN
1	B	25	LYS
1	B	30	LEU
1	B	43	HIS
1	B	83	ASN
1	B	88	PRO
1	B	112	LEU
1	B	124	LEU
1	B	128	ASP
1	B	151	GLU
1	B	176	LEU
1	B	205	GLU
1	B	206	GLU
1	B	215	THR
1	B	222	LYS
1	B	229	ILE
1	B	230	GLN
1	B	233	LYS
1	B	259	LYS
1	B	272	ILE
1	B	274	LYS
1	B	306	LEU
1	B	346	GLU
1	B	356	ILE
1	B	394	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	83	ASN
1	A	111	GLN
1	A	139	HIS
1	A	156	GLN
1	A	158	ASN
1	A	166	GLN
1	A	269	HIS
1	B	44	GLN
1	B	97	GLN
1	B	109	GLN
1	B	116	ASN
1	B	139	HIS
1	B	230	GLN

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Mol	Chain	Res	Type
1	B	269	HIS
1	B	287	HIS

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 [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	372/372 (100%)	-1.29	0 100 100	32, 57, 106, 137	0
1	B	372/372 (100%)	-1.25	0 100 100	32, 59, 106, 165	0
All	All	744/744 (100%)	-1.27	0 100 100	32, 58, 106, 165	0

There are no RSRZ outliers to report.

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.