



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:00 AM UTC

PDB ID : 5EGW / pdb_00005egw
Title : 2.70 Å crystal structure of the Amb a 11 cysteine protease, a major ragweed pollen allergen, in its proform
Authors : Briozzo, P.; Kopecny, D.
Deposited on : 2015-10-27
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

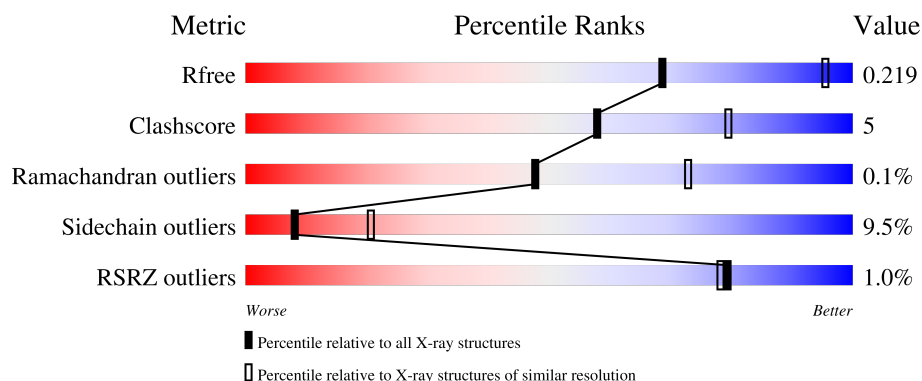
1 Overall quality at a glance

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	A	385	
1	B	385	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5423 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 Cysteine protease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	363	Total	C	N	O	S	0	0	0
			2858	1799	505	537	17			
1	B	325	Total	C	N	O	S	0	0	0
			2559	1611	454	477	17			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	MET	-	initiating methionine	UNP V5LU01
A	3	GLY	-	expression tag	UNP V5LU01
A	4	SER	-	expression tag	UNP V5LU01
A	5	SER	-	expression tag	UNP V5LU01
A	6	HIS	-	expression tag	UNP V5LU01
A	7	HIS	-	expression tag	UNP V5LU01
A	8	HIS	-	expression tag	UNP V5LU01
A	9	HIS	-	expression tag	UNP V5LU01
A	10	HIS	-	expression tag	UNP V5LU01
A	11	HIS	-	expression tag	UNP V5LU01
A	12	SER	-	expression tag	UNP V5LU01
A	13	SER	-	expression tag	UNP V5LU01
A	14	GLY	-	expression tag	UNP V5LU01
A	15	LEU	-	expression tag	UNP V5LU01
A	16	VAL	-	expression tag	UNP V5LU01
A	17	PRO	-	expression tag	UNP V5LU01
A	18	ARG	-	expression tag	UNP V5LU01
A	19	GLY	-	expression tag	UNP V5LU01
A	20	SER	-	expression tag	UNP V5LU01
A	21	HIS	-	expression tag	UNP V5LU01
A	22	MET	-	expression tag	UNP V5LU01
B	2	MET	-	initiating methionine	UNP V5LU01
B	3	GLY	-	expression tag	UNP V5LU01
B	4	SER	-	expression tag	UNP V5LU01
B	5	SER	-	expression tag	UNP V5LU01

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Chain	Residue	Modelled	Actual	Comment	Reference
B	6	HIS	-	expression tag	UNP V5LU01
B	7	HIS	-	expression tag	UNP V5LU01
B	8	HIS	-	expression tag	UNP V5LU01
B	9	HIS	-	expression tag	UNP V5LU01
B	10	HIS	-	expression tag	UNP V5LU01
B	11	HIS	-	expression tag	UNP V5LU01
B	12	SER	-	expression tag	UNP V5LU01
B	13	SER	-	expression tag	UNP V5LU01
B	14	GLY	-	expression tag	UNP V5LU01
B	15	LEU	-	expression tag	UNP V5LU01
B	16	VAL	-	expression tag	UNP V5LU01
B	17	PRO	-	expression tag	UNP V5LU01
B	18	ARG	-	expression tag	UNP V5LU01
B	19	GLY	-	expression tag	UNP V5LU01
B	20	SER	-	expression tag	UNP V5LU01
B	21	HIS	-	expression tag	UNP V5LU01
B	22	MET	-	expression tag	UNP V5LU01

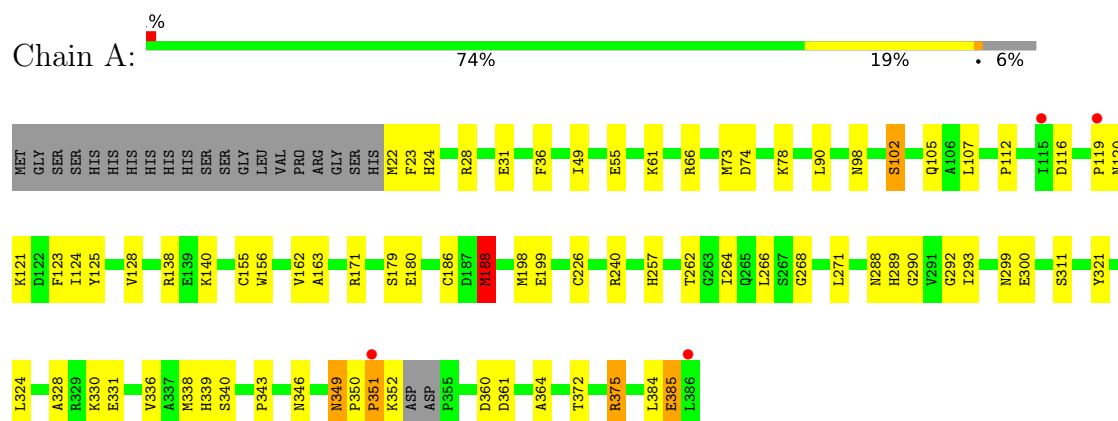
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	4	Total O 4 4	0	0
2	B	2	Total O 2 2	0	0

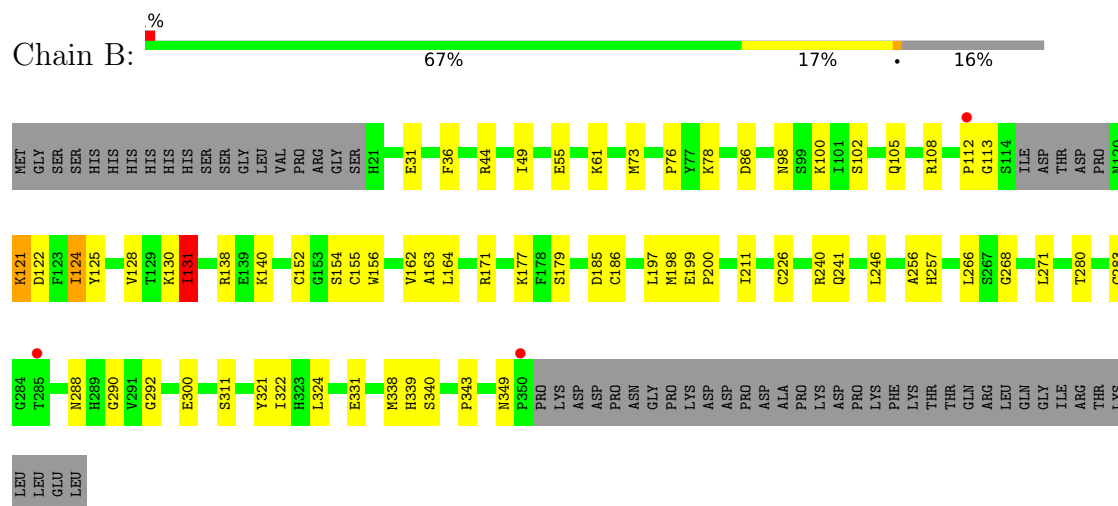
3 Residue-property plots

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: Cysteine protease



• Molecule 1: Cysteine protease



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.45Å 89.55Å 104.06Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.78 – 2.70 44.78 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.0 (44.78-2.70) 99.0 (44.78-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 2.69Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.10.0	Depositor
R, R_{free}	0.179 , 0.218 (Not available) , 0.219	Depositor DCC
R_{free} test set	1153 reflections (5.28%)	wwPDB-VP
Wilson B-factor (Å ²)	65.3	Xtriage
Anisotropy	0.553	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 61.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5423	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.79% 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	0.89	2/2928 (0.1%)	1.42	25/3948 (0.6%)
1	B	0.83	0/2622	1.38	15/3534 (0.4%)
All	All	0.86	2/5550 (0.0%)	1.40	40/7482 (0.5%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	188	MET	SD-CE	5.16	1.92	1.79
1	A	121	LYS	C-N	5.08	1.37	1.33

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	120	ASN	CA-C-N	7.00	134.30	121.70
1	A	120	ASN	C-N-CA	7.00	134.30	121.70
1	A	385	GLU	CA-C-N	6.99	134.29	121.70
1	A	385	GLU	C-N-CA	6.99	134.29	121.70
1	B	113	GLY	CA-C-N	6.56	133.50	121.70
1	B	113	GLY	C-N-CA	6.56	133.50	121.70
1	B	86	ASP	CA-CB-CG	6.53	119.13	112.60
1	B	122	ASP	CA-CB-CG	6.39	118.99	112.60
1	B	179	SER	N-CA-C	6.27	118.50	108.41
1	B	121	LYS	N-CA-C	6.22	119.35	111.69
1	A	179	SER	N-CA-C	6.16	118.33	108.41
1	B	268	GLY	CA-C-N	6.11	128.78	120.54
1	B	268	GLY	C-N-CA	6.11	128.78	120.54
1	A	351	PRO	CA-C-N	6.02	132.53	121.70
1	A	351	PRO	C-N-CA	6.02	132.53	121.70
1	A	268	GLY	CA-C-N	5.97	128.61	120.54
1	A	268	GLY	C-N-CA	5.97	128.61	120.54
1	B	185	ASP	CA-CB-CG	5.92	118.52	112.60
1	A	98	ASN	N-CA-C	5.86	118.04	110.65
1	A	102	SER	CA-C-N	5.80	128.30	120.65

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	102	SER	C-N-CA	5.80	128.30	120.65
1	B	44	ARG	CA-C-N	5.71	127.93	120.28
1	B	44	ARG	C-N-CA	5.71	127.93	120.28
1	A	299	ASN	CA-CB-CG	5.65	118.25	112.60
1	A	116	ASP	CA-C-N	5.59	127.70	120.44
1	A	116	ASP	C-N-CA	5.59	127.70	120.44
1	A	328	ALA	N-CA-C	5.54	118.50	111.69
1	B	98	ASN	N-CA-C	5.51	117.59	110.65
1	A	123	PHE	CA-CB-CG	5.46	119.27	113.80
1	B	124	ILE	N-CA-C	5.45	117.86	111.05
1	A	124	ILE	N-CA-C	5.34	118.53	111.17
1	A	180	GLU	N-CA-C	-5.24	105.63	112.23
1	A	74	ASP	CA-CB-CG	5.23	117.83	112.60
1	B	131	ILE	CA-CB-CG2	5.17	119.29	110.50
1	B	288	ASN	N-CA-C	5.13	119.39	113.18
1	A	346	ASN	CA-CB-CG	5.11	117.70	112.60
1	A	349	ASN	CA-CB-CG	5.08	117.68	112.60
1	A	288	ASN	CA-CB-CG	5.08	117.68	112.60
1	A	66	ARG	CA-C-N	5.04	127.01	120.56
1	A	66	ARG	C-N-CA	5.04	127.01	120.56

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	2858	0	2774	27	0
1	B	2559	0	2461	20	0
2	A	4	0	0	0	0
2	B	2	0	0	0	0
All	All	5423	0	5235	46	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 5.

All (46) 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:199:GLU:HG2	1:B:200:PRO:HD3	1.51	0.92
1:A:155:CYS:SG	1:A:290:GLY:N	2.43	0.91
1:B:155:CYS:SG	1:B:290:GLY:N	2.46	0.89
1:A:186:CYS:HG	1:A:226:CYS:HG	1.30	0.77
1:A:188:MET:HE2	1:A:188:MET:HA	1.68	0.74
1:B:186:CYS:HG	1:B:226:CYS:HG	1.35	0.74
1:A:264:ILE:HD12	1:A:266:LEU:HD21	1.70	0.74
1:A:324:LEU:HD23	1:A:336:VAL:HG11	1.76	0.68
1:B:112:PRO:HG3	1:B:338:MET:HE3	1.84	0.58
1:B:128:VAL:HG21	1:B:257:HIS:HE1	1.67	0.58
1:A:188:MET:HA	1:A:188:MET:CE	2.32	0.58
1:A:155:CYS:SG	1:A:289:HIS:HA	2.47	0.55
1:B:266:LEU:HD22	1:B:271:LEU:HD22	1.90	0.54
1:A:155:CYS:HG	1:A:290:GLY:H	1.52	0.53
1:A:112:PRO:HG3	1:A:338:MET:HE3	1.91	0.52
1:A:266:LEU:HD22	1:A:271:LEU:HD22	1.92	0.52
1:A:264:ILE:CG2	1:A:336:VAL:HG13	2.41	0.51
1:B:76:PRO:HG2	1:B:280:THR:HG21	1.94	0.49
1:A:264:ILE:HD12	1:A:266:LEU:CD2	2.43	0.47
1:B:152:CYS:SG	1:B:154:SER:HB3	2.54	0.47
1:A:125:TYR:O	1:A:128:VAL:HG22	2.15	0.47
1:A:361:ASP:HB3	1:A:364:ALA:HB2	1.96	0.47
1:B:128:VAL:HG21	1:B:131:ILE:HD11	1.97	0.47
1:A:264:ILE:CG2	1:A:336:VAL:CG1	2.92	0.47
1:B:124:ILE:HG23	1:B:241:GLN:HE22	1.79	0.47
1:B:131:ILE:HG12	1:B:256:ALA:CB	2.44	0.47
1:A:156:TRP:O	1:A:198:MET:HG3	2.14	0.46
1:B:156:TRP:O	1:B:198:MET:HG3	2.14	0.46
1:B:125:TYR:O	1:B:128:VAL:HG22	2.15	0.46
1:A:138:ARG:HD3	1:A:321:TYR:CZ	2.51	0.46
1:A:262:THR:HG21	1:A:293:ILE:HD11	1.98	0.45
1:A:375:ARG:HD2	1:A:375:ARG:HA	1.67	0.44
1:B:138:ARG:HD3	1:B:321:TYR:CZ	2.51	0.44
1:A:128:VAL:HG21	1:A:257:HIS:HE1	1.83	0.43
1:A:22:MET:HG2	1:A:23:PHE:N	2.33	0.43
1:B:105:GLN:HG2	1:B:339:HIS:CE1	2.52	0.43
1:A:105:GLN:HG2	1:A:339:HIS:CE1	2.53	0.43
1:A:163:ALA:HB1	1:A:343:PRO:HD3	2.00	0.43
1:A:324:LEU:CD2	1:A:336:VAL:HG11	2.48	0.42
1:B:163:ALA:HB1	1:B:343:PRO:HD3	2.01	0.42
1:A:350:PRO:HA	1:A:351:PRO:HD3	1.95	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:VAL:HG11	1:A:292:GLY:HA3	2.02	0.41
1:B:162:VAL:HG11	1:B:292:GLY:HA3	2.03	0.41
1:B:131:ILE:HG12	1:B:256:ALA:HB1	2.02	0.40
1:A:90:LEU:HD21	1:B:100:LYS:HG2	2.04	0.40
1:B:164:LEU:HD13	1:B:211:ILE:HG21	2.03	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	359/385 (93%)	344 (96%)	14 (4%)	1 (0%)	36	60
1	B	321/385 (83%)	309 (96%)	12 (4%)	0	100	100
All	All	680/770 (88%)	653 (96%)	26 (4%)	1 (0%)	48	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	PRO

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	303/322 (94%)	275 (91%)	28 (9%)	8	22
1	B	268/322 (83%)	242 (90%)	26 (10%)	8	20
All	All	571/644 (89%)	517 (90%)	54 (10%)	8	20

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	HIS
1	A	28	ARG
1	A	31	GLU
1	A	36	PHE
1	A	49	ILE
1	A	55	GLU
1	A	61	LYS
1	A	73	MET
1	A	78	LYS
1	A	102	SER
1	A	107	LEU
1	A	140	LYS
1	A	171	ARG
1	A	188	MET
1	A	199	GLU
1	A	240	ARG
1	A	300	GLU
1	A	311	SER
1	A	330	LYS
1	A	331	GLU
1	A	340	SER
1	A	349	ASN
1	A	352	LYS
1	A	360	ASP
1	A	372	THR
1	A	375	ARG
1	A	384	LEU
1	A	385	GLU
1	B	31	GLU
1	B	36	PHE
1	B	49	ILE
1	B	55	GLU

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Mol	Chain	Res	Type
1	B	61	LYS
1	B	73	MET
1	B	78	LYS
1	B	102	SER
1	B	108	ARG
1	B	121	LYS
1	B	130	LYS
1	B	131	ILE
1	B	140	LYS
1	B	171	ARG
1	B	177	LYS
1	B	197	LEU
1	B	240	ARG
1	B	246	LEU
1	B	283	CYS
1	B	300	GLU
1	B	311	SER
1	B	322	ILE
1	B	324	LEU
1	B	331	GLU
1	B	340	SER
1	B	349	ASN

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	58	ASN
1	A	168	ASN
1	B	58	ASN
1	B	94	ASN
1	B	98	ASN
1	B	168	ASN
1	B	265	GLN
1	B	325	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 [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	363/385 (94%)	-0.20	4 (1%) 78 76	50, 70, 106, 141	0
1	B	325/385 (84%)	-0.16	3 (0%) 81 80	51, 74, 117, 138	0
All	All	688/770 (89%)	-0.18	7 (1%) 79 78	50, 71, 115, 141	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	386	LEU	3.2
1	A	119	PRO	2.7
1	A	115	ILE	2.4
1	A	351	PRO	2.3
1	B	112	PRO	2.3
1	B	285	THR	2.1
1	B	350	PRO	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 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.