



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:51 AM UTC

PDB ID : 2PBH / pdb_00002pbh
Title : CRYSTAL STRUCTURE OF HUMAN PROCATHEPSIN B AT 3.3
ANGSTROM RESOLUTION
Authors : Podobnik, M.; Turk, D.; Kuhelj, R.; Turk, V.
Deposited on : 1997-02-14
Resolution : 3.30 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

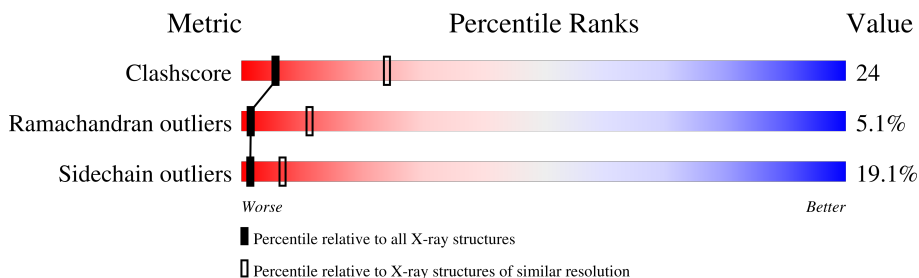
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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))
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	317	43% 41% 13% .

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 3015 atoms, of which 545 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 PROCATHEPSIN B.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	317	Total	C	H	N	O	S	673	0	0
			3015	1552	545	428	468	22			

There are 17 discrepancies between the modelled and reference sequences:

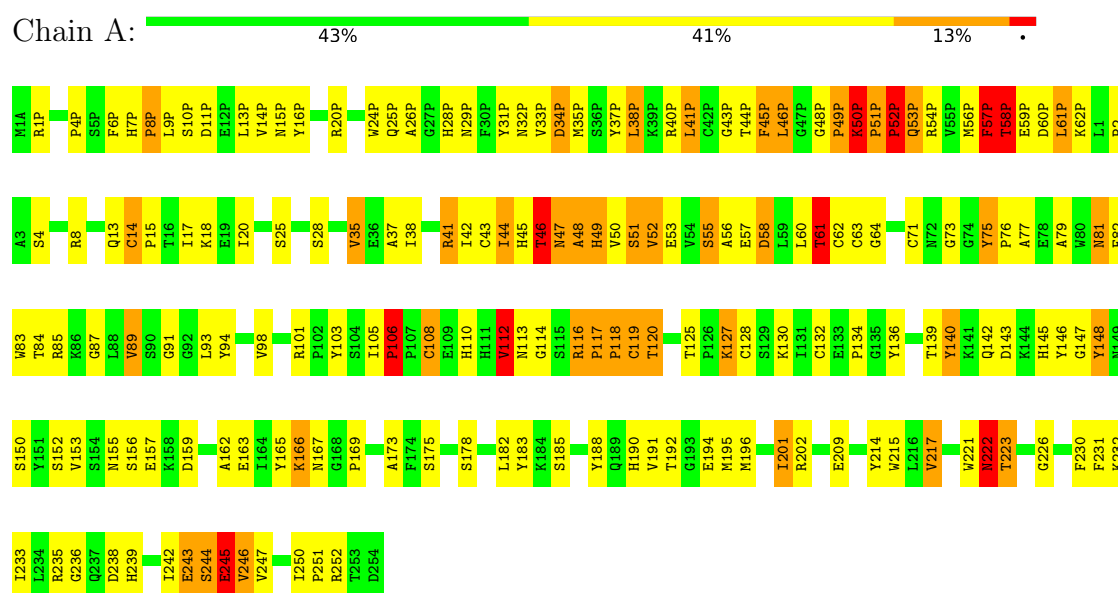
Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	TRP	deletion	UNP P07858
A	?	-	GLN	deletion	UNP P07858
A	?	-	LEU	deletion	UNP P07858
A	?	-	TRP	deletion	UNP P07858
A	?	-	ALA	deletion	UNP P07858
A	?	-	SER	deletion	UNP P07858
A	?	-	LEU	deletion	UNP P07858
A	?	-	CYS	deletion	UNP P07858
A	?	-	CYS	deletion	UNP P07858
A	?	-	LEU	deletion	UNP P07858
A	?	-	LEU	deletion	UNP P07858
A	?	-	VAL	deletion	UNP P07858
A	?	-	LEU	deletion	UNP P07858
A	?	-	ALA	deletion	UNP P07858
A	?	-	ASN	deletion	UNP P07858
A	?	-	ALA	deletion	UNP P07858
A	9P	LEU	VAL	conflict	UNP P07858

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. 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. 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.

Note EDS was not executed.

• Molecule 1: PROCATHEPSIN B



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	78.75Å 78.75Å 153.64Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 3.30	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-3.30)	Depositor
R_{merge}	0.19	Depositor
R_{sym}	0.08	Depositor
Refinement program	MAIN	Depositor
R, R_{free}	0.231 , 0.269	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3015	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

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	0.69	0/2545	1.28	42/3457 (1.2%)

There are no bond length outliers.

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	45(P)	PHE	N-CA-C	10.32	123.99	111.40
1	A	47	ASN	N-CA-C	-9.86	93.41	108.00
1	A	116	ARG	CA-C-N	-9.10	111.01	120.38
1	A	116	ARG	C-N-CA	-9.10	111.01	120.38
1	A	50(P)	LYS	CA-C-N	8.37	129.00	120.38
1	A	50(P)	LYS	C-N-CA	8.37	129.00	120.38
1	A	106	PRO	CA-C-N	-7.76	112.70	120.31
1	A	106	PRO	C-N-CA	-7.76	112.70	120.31
1	A	113	ASN	N-CA-C	7.50	120.14	108.96
1	A	51	SER	N-CA-C	-7.21	97.49	109.24
1	A	48(P)	GLY	CA-C-N	-7.12	110.94	119.84
1	A	48(P)	GLY	C-N-CA	-7.12	110.94	119.84
1	A	61	THR	N-CA-C	7.02	122.01	113.38
1	A	127	LYS	N-CA-C	6.90	121.29	110.32
1	A	125	THR	CA-C-N	6.83	127.23	119.92
1	A	125	THR	C-N-CA	6.83	127.23	119.92
1	A	60(P)	ASP	N-CA-C	6.76	121.55	113.16
1	A	46	THR	N-CA-C	-6.73	102.11	111.81
1	A	112	VAL	N-CA-C	6.53	122.93	109.34
1	A	35	VAL	N-CA-C	6.46	117.25	110.72
1	A	14	CYS	CA-C-N	6.32	126.23	119.28
1	A	14	CYS	C-N-CA	6.32	126.23	119.28
1	A	59(P)	GLU	N-CA-C	6.22	117.77	108.31
1	A	48	ALA	N-CA-C	-6.17	103.17	111.87
1	A	56	ALA	N-CA-C	-6.13	104.51	112.23
1	A	55	SER	N-CA-C	6.10	118.58	109.25
1	A	165	TYR	N-CA-C	6.09	118.00	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	245	GLU	N-CA-C	6.00	118.22	110.65
1	A	196	MET	N-CA-C	-6.00	105.45	112.89
1	A	217	VAL	N-CA-C	5.94	117.04	108.48
1	A	98	VAL	N-CA-C	5.93	116.15	107.37
1	A	178	SER	N-CA-C	5.90	119.30	111.75
1	A	117	PRO	CA-C-N	5.77	127.05	119.84
1	A	117	PRO	C-N-CA	5.77	127.05	119.84
1	A	185	SER	N-CA-C	5.75	116.99	108.60
1	A	145	HIS	N-CA-C	-5.73	99.91	109.24
1	A	29(P)	ASN	N-CA-C	-5.36	106.80	113.55
1	A	75	TYR	CA-C-N	5.35	125.17	119.28
1	A	75	TYR	C-N-CA	5.35	125.17	119.28
1	A	34(P)	ASP	N-CA-C	-5.27	102.48	110.28
1	A	64	GLY	N-CA-C	5.16	121.20	112.22
1	A	15(P)	ASN	N-CA-C	5.09	116.91	111.36

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	2470	545	2305	106	5
All	All	2470	545	2305	106	5

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 24.

All (106) 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:155:ASN:OD1	1:A:244:SER:HA	1.58	1.01
1:A:188:TYR:HD2	1:A:233:ILE:HG13	1.41	0.85
1:A:52(P):PRO:HD2	1:A:150:SER:O	1.79	0.83
1:A:81:ASN:HD22	1:A:85:ARG:HE	1.30	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:153:VAL:HB	1:A:246:VAL:HB	1.65	0.77
1:A:40(P):ARG:HH21	1:A:71:CYS:HB3	1.51	0.76
1:A:20(P):ARG:O	1:A:20(P):ARG:HG3	1.88	0.74
1:A:20(P):ARG:NH1	1:A:191:VAL:HG13	2.05	0.72
1:A:56(P):MET:HE2	1:A:162:ALA:HB1	1.71	0.72
1:A:108:CYS:SG	1:A:110:HIS:HB3	2.29	0.72
1:A:10(P):SER:O	1:A:13(P):LEU:HB2	1.90	0.71
1:A:13:GLN:HG3	1:A:48:ALA:HB1	1.71	0.70
1:A:175:SER:HB3	1:A:195:MET:SD	2.33	0.68
1:A:24(P):TRP:CZ3	1:A:26(P):ALA:HB2	2.32	0.65
1:A:43(P):GLY:O	1:A:73:GLY:HA2	1.97	0.64
1:A:8:ARG:HG2	1:A:17:ILE:HG22	1.80	0.63
1:A:20:ILE:HD11	1:A:230:PHE:CE1	2.34	0.63
1:A:46:THR:O	1:A:47:ASN:HB2	2.01	0.60
1:A:46:THR:O	1:A:49:HIS:HB2	2.02	0.59
1:A:34(P):ASP:HB3	1:A:120:THR:HA	1.83	0.59
1:A:76:PRO:O	1:A:79:ALA:HB3	2.02	0.59
1:A:13:GLN:O	1:A:15:PRO:HD3	2.03	0.58
1:A:222:ASN:HD22	1:A:223:THR:H	1.53	0.57
1:A:45(P):PHE:CZ	1:A:73:GLY:HA3	2.40	0.57
1:A:9(P):LEU:HB3	1:A:13(P):LEU:HD12	1.87	0.57
1:A:7(P):HIS:O	1:A:8(P):PRO:C	2.47	0.56
1:A:103:TYR:CE2	1:A:105:ILE:HB	2.40	0.56
1:A:215:TRP:CD1	1:A:235:ARG:HB2	2.41	0.56
1:A:40(P):ARG:NH2	1:A:71:CYS:HB3	2.21	0.56
1:A:190:HIS:H	1:A:239:HIS:CE1	2.23	0.55
1:A:37:ALA:O	1:A:41:ARG:HG3	2.06	0.54
1:A:221:TRP:O	1:A:222:ASN:HB2	2.06	0.54
1:A:63:CYS:HA	1:A:82:PHE:CE2	2.43	0.53
1:A:46:THR:O	1:A:46:THR:HG23	2.08	0.53
1:A:13(P):LEU:HD13	1:A:182:LEU:HD11	1.91	0.53
1:A:61(P):LEU:HD13	1:A:166:LYS:HE2	1.91	0.53
1:A:51(P):PRO:HG2	1:A:247:VAL:HG23	1.91	0.52
1:A:156:SER:HB3	1:A:159:ASP:OD2	2.10	0.51
1:A:46(P):LEU:HA	1:A:245:GLU:OE2	2.09	0.51
1:A:118:PRO:O	1:A:120:THR:N	2.43	0.51
1:A:93:LEU:HD12	1:A:93:LEU:N	2.26	0.51
1:A:7(P):HIS:O	1:A:7(P):HIS:CG	2.64	0.50
1:A:94:TYR:HA	1:A:103:TYR:O	2.12	0.50
1:A:38(P):LEU:HA	1:A:41(P):LEU:HD22	1.92	0.49
1:A:2:PRO:HB2	1:A:4:SER:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:CYS:SG	1:A:140:TYR:HA	2.52	0.49
1:A:31(P):TYR:O	1:A:32(P):ASN:HB3	2.11	0.49
1:A:183:TYR:O	1:A:226:GLY:HA2	2.12	0.49
1:A:139:THR:HG23	1:A:142:GLN:OE1	2.13	0.49
1:A:140:TYR:C	1:A:140:TYR:CD1	2.90	0.49
1:A:201:ILE:HB	1:A:217:VAL:CG1	2.43	0.49
1:A:32(P):ASN:OD1	1:A:118:PRO:HG3	2.13	0.49
1:A:42:ILE:O	1:A:46:THR:HG22	2.12	0.48
1:A:105:ILE:HG23	1:A:106:PRO:HD2	1.95	0.48
1:A:89:VAL:HG12	1:A:143:ASP:HB3	1.93	0.48
1:A:146:TYR:N	1:A:146:TYR:CD1	2.81	0.48
1:A:9(P):LEU:HD13	1:A:182:LEU:HD21	1.96	0.48
1:A:50:VAL:HG22	1:A:52:VAL:HG12	1.96	0.47
1:A:147:GLY:C	1:A:252:ARG:HG3	2.39	0.47
1:A:44:ILE:HG21	1:A:167:ASN:O	2.14	0.47
1:A:221:TRP:O	1:A:222:ASN:CB	2.61	0.47
1:A:49(P):PRO:CG	1:A:77:ALA:HB3	2.44	0.47
1:A:173:ALA:HB3	1:A:245:GLU:CB	2.44	0.47
1:A:222:ASN:ND2	1:A:223:THR:H	2.13	0.47
1:A:61:THR:O	1:A:128:CYS:HB2	2.14	0.47
1:A:57(P):PHE:CD1	1:A:57(P):PHE:N	2.76	0.46
1:A:128:CYS:SG	1:A:130:LYS:HE2	2.56	0.46
1:A:209:GLU:HB2	1:A:214:TYR:HE2	1.82	0.45
1:A:61(P):LEU:HD13	1:A:166:LYS:CE	2.46	0.45
1:A:43:CYS:HB2	1:A:51:SER:HA	1.99	0.44
1:A:55:SER:HB2	1:A:91:GLY:N	2.31	0.44
1:A:148:TYR:HB2	1:A:250:ILE:O	2.17	0.44
1:A:130:LYS:HA	1:A:140:TYR:CE2	2.52	0.44
1:A:235:ARG:HG3	1:A:236:GLY:N	2.32	0.44
1:A:209:GLU:HB2	1:A:214:TYR:CE2	2.53	0.43
1:A:49(P):PRO:HG2	1:A:77:ALA:HB3	2.00	0.43
1:A:44:ILE:HG22	1:A:45:HIS:ND1	2.34	0.43
1:A:153:VAL:HG22	1:A:163:GLU:HG3	2.01	0.42
1:A:46:THR:HG21	1:A:50:VAL:O	2.19	0.42
1:A:58:ASP:OD1	1:A:101:ARG:NH1	2.53	0.42
1:A:173:ALA:HB3	1:A:245:GLU:HB2	2.02	0.42
1:A:45(P):PHE:O	1:A:46(P):LEU:HB2	2.20	0.42
1:A:139:THR:O	1:A:142:GLN:N	2.53	0.42
1:A:24(P):TRP:HZ3	1:A:26(P):ALA:HB2	1.82	0.42
1:A:45:HIS:C	1:A:47:ASN:H	2.27	0.42
1:A:53(P):GLN:NE2	1:A:152:SER:HB3	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:CYS:C	1:A:110:HIS:H	2.28	0.42
1:A:45:HIS:HD2	1:A:251:PRO:HD2	1.84	0.41
1:A:63:CYS:SG	1:A:82:PHE:CD2	3.14	0.41
1:A:146:TYR:HB2	1:A:252:ARG:HH11	1.84	0.41
1:A:159:ASP:O	1:A:162:ALA:HB3	2.20	0.41
1:A:38(P):LEU:HA	1:A:41(P):LEU:CD2	2.50	0.41
1:A:217:VAL:HB	1:A:231:PHE:CZ	2.56	0.41
1:A:11(P):ASP:OD1	1:A:28(P):HIS:HE1	2.03	0.41
1:A:28:SER:HB3	1:A:60:LEU:HD13	2.03	0.41
1:A:16(P):TYR:CD1	1:A:16(P):TYR:C	2.96	0.41
1:A:75:TYR:HA	1:A:76:PRO:HD3	1.81	0.41
1:A:83:TRP:O	1:A:87:GLY:HA2	2.21	0.41
1:A:242:ILE:HG23	1:A:243:GLU:N	2.36	0.40
1:A:34(P):ASP:O	1:A:37(P):TYR:HB3	2.22	0.40
1:A:209:GLU:OE1	1:A:232:LYS:NZ	2.54	0.40
1:A:50(P):LYS:HA	1:A:51(P):PRO:HD3	1.80	0.40
1:A:54(P):ARG:HA	1:A:54(P):ARG:HD3	1.83	0.40
1:A:132:CYS:HB3	1:A:136:TYR:HD2	1.87	0.40
1:A:37(P):TYR:O	1:A:41(P):LEU:HD22	2.22	0.40
1:A:91:GLY:HA3	1:A:101:ARG:O	2.22	0.40

All (5) 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)
1:A:116:ARG:NH1	1:A:238:ASP:OD2[9_875]	1.80	0.40
1:A:1(P):ARG:HH11	1:A:58(P):THR:CG2[9_875]	1.29	0.31
1:A:116:ARG:HH11	1:A:238:ASP:OD2[9_875]	1.32	0.28
1:A:18:LYS:HZ3	1:A:148:TYR:CE1[6_544]	1.36	0.24
1:A:112:VAL:O	1:A:114:GLY:C[10_775]	2.13	0.07

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	A	315/317 (99%)	263 (84%)	36 (11%)	16 (5%)	1	11

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8(P)	PRO
1	A	49(P)	PRO
1	A	50(P)	LYS
1	A	52(P)	PRO
1	A	49	HIS
1	A	119	CYS
1	A	6(P)	PHE
1	A	51(P)	PRO
1	A	58(P)	THR
1	A	57(P)	PHE
1	A	222	ASN
1	A	89	VAL
1	A	112	VAL
1	A	169	PRO
1	A	118	PRO
1	A	4(P)	PRO

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	267/267 (100%)	216 (81%)	51 (19%)	1	7

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	14(P)	VAL
1	A	25(P)	GLN
1	A	33(P)	VAL
1	A	35(P)	MET

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Mol	Chain	Res	Type
1	A	38(P)	LEU
1	A	41(P)	LEU
1	A	44(P)	THR
1	A	46(P)	LEU
1	A	52(P)	PRO
1	A	53(P)	GLN
1	A	57(P)	PHE
1	A	58(P)	THR
1	A	61(P)	LEU
1	A	62(P)	LYS
1	A	14	CYS
1	A	25	SER
1	A	35	VAL
1	A	38	ILE
1	A	41	ARG
1	A	44	ILE
1	A	46	THR
1	A	52	VAL
1	A	53	GLU
1	A	57	GLU
1	A	58	ASP
1	A	61	THR
1	A	62	CYS
1	A	81	ASN
1	A	84	THR
1	A	106	PRO
1	A	108	CYS
1	A	112	VAL
1	A	117	PRO
1	A	119	CYS
1	A	120	THR
1	A	127	LYS
1	A	134	PRO
1	A	140	TYR
1	A	148	TYR
1	A	157	GLU
1	A	166	LYS
1	A	192	THR
1	A	194	GLU
1	A	201	ILE
1	A	202	ARG
1	A	222	ASN

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Mol	Chain	Res	Type
1	A	223	THR
1	A	243	GLU
1	A	244	SER
1	A	245	GLU
1	A	246	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25(P)	GLN
1	A	28(P)	HIS
1	A	10	GLN
1	A	45	HIS
1	A	72	ASN
1	A	81	ASN
1	A	167	ASN
1	A	190	HIS
1	A	222	ASN
1	A	239	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

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.