



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:29 PM UTC

PDB ID : 1ZVU / pdb_00001zvu
Title : Structure of the full-length E. coli ParC subunit
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Deposited on : 2005-06-02
Resolution : 3.00 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

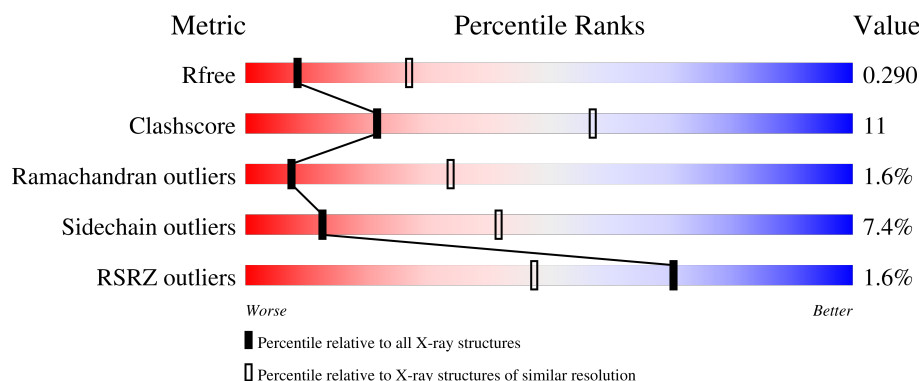
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	716	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5402 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 Topoisomerase IV subunit A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	685	Total	C	N	O	S	0	0	0
			5367	3369	964	1006	28			

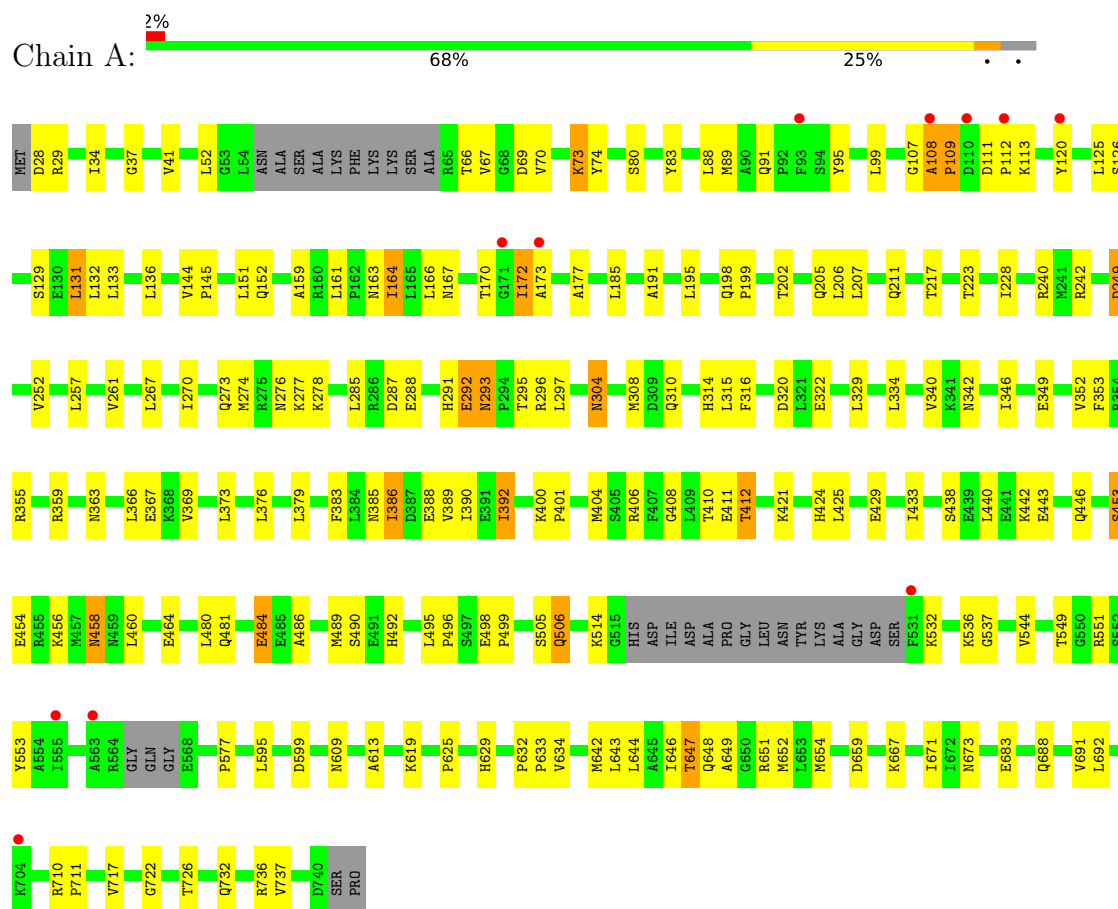
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	35	Total	O	0	0
			35	35		

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: Topoisomerase IV subunit A



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	257.99Å 62.14Å 64.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 20.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	89.6 (20.00-3.00) 89.2 (20.00-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.68 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.240 , 0.296 0.235 , 0.290	Depositor DCC
R_{free} test set	994 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	47.0	Xtriage
Anisotropy	0.542	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 43.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.032 for -h,l,k	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	5402	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.12% 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	0.49	0/5454	0.85	8/7368 (0.1%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	495	LEU	CA-C-N	6.41	127.85	119.84
1	A	495	LEU	C-N-CA	6.41	127.85	119.84
1	A	498	GLU	CA-C-N	6.23	127.63	119.84
1	A	498	GLU	C-N-CA	6.23	127.63	119.84
1	A	292	GLU	N-CA-C	-5.80	104.38	112.45
1	A	671	ILE	N-CA-C	-5.48	107.29	111.62
1	A	304	ASN	N-CA-C	-5.09	107.62	113.88
1	A	537	GLY	N-CA-C	5.08	117.47	111.63

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	5367	0	5487	117	0
2	A	35	0	0	1	0
All	All	5402	0	5487	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 11.

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:164:ILE:H	1:A:164:ILE:HD12	1.26	1.01
1:A:108:ALA:HB3	1:A:109:PRO:HD3	1.47	0.95
1:A:647:THR:HG22	1:A:651:ARG:H	1.33	0.93
1:A:111:ASP:HB3	1:A:112:PRO:CD	2.11	0.81
1:A:369:VAL:HG23	1:A:440:LEU:HD22	1.65	0.79
1:A:89:MET:HG2	1:A:95:TYR:HE2	1.49	0.76
1:A:355:ARG:HH12	1:A:458:ASN:HD21	1.33	0.76
1:A:108:ALA:CB	1:A:109:PRO:HD3	2.17	0.74
1:A:288:GLU:OE1	1:A:296:ARG:NH1	2.21	0.74
1:A:164:ILE:H	1:A:164:ILE:CD1	1.99	0.72
1:A:202:THR:H	1:A:205:GLN:HE21	1.38	0.70
1:A:484:GLU:CD	1:A:484:GLU:H	2.02	0.67
1:A:270:ILE:O	1:A:274:MET:HG3	1.94	0.67
1:A:388:GLU:O	1:A:392:ILE:HG22	1.96	0.66
1:A:270:ILE:HG22	1:A:274:MET:HE2	1.79	0.64
1:A:267:LEU:HD21	1:A:287:ASP:HB2	1.79	0.64
1:A:223:THR:HG23	1:A:240:ARG:H	1.62	0.64
1:A:276:ASN:C	1:A:278:LYS:H	2.07	0.63
1:A:355:ARG:NH1	1:A:458:ASN:HD21	1.97	0.63
1:A:66:THR:HG22	1:A:69:ASP:CG	2.24	0.63
1:A:400:LYS:HB3	1:A:401:PRO:HD3	1.80	0.62
1:A:89:MET:HG2	1:A:95:TYR:CE2	2.35	0.61
1:A:595:LEU:HB3	1:A:634:VAL:HB	1.83	0.61
1:A:421:LYS:H	1:A:424:HIS:HD2	1.49	0.60
1:A:242:ARG:HD3	2:A:765:HOH:O	2.00	0.60
1:A:386:ILE:HA	1:A:389:VAL:HG12	1.84	0.60
1:A:163:ASN:HD21	1:A:167:ASN:HD22	1.50	0.59
1:A:202:THR:H	1:A:205:GLN:NE2	1.99	0.59
1:A:386:ILE:HA	1:A:389:VAL:CG1	2.33	0.58
1:A:642:MET:HE3	1:A:654:MET:HG2	1.85	0.58
1:A:107:GLY:CA	1:A:113:LYS:HB2	2.34	0.58
1:A:274:MET:HE1	1:A:285:LEU:HB2	1.85	0.57
1:A:505:SER:HA	1:A:532:LYS:HB3	1.85	0.57
1:A:111:ASP:HB3	1:A:112:PRO:HD3	1.85	0.56
1:A:108:ALA:HB3	1:A:109:PRO:CD	2.31	0.55
1:A:314:HIS:HD2	1:A:489:MET:H	1.54	0.55
1:A:111:ASP:HB3	1:A:112:PRO:HD2	1.89	0.54
1:A:252:VAL:HG23	1:A:308:MET:HG2	1.90	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:373:LEU:HD21	1:A:440:LEU:HB2	1.88	0.54
1:A:410:THR:C	1:A:412:THR:H	2.15	0.54
1:A:673:ASN:HB2	1:A:722:GLY:O	2.08	0.54
1:A:66:THR:HG22	1:A:69:ASP:H	1.73	0.53
1:A:454:GLU:O	1:A:458:ASN:ND2	2.42	0.53
1:A:647:THR:HG23	1:A:649:ALA:H	1.74	0.53
1:A:37:GLY:CA	1:A:164:ILE:HG13	2.39	0.53
1:A:710:ARG:HB3	1:A:711:PRO:HD2	1.92	0.52
1:A:88:LEU:HD21	1:A:170:THR:OG1	2.09	0.52
1:A:646:ILE:HB	1:A:688:GLN:HB2	1.91	0.52
1:A:211:GLN:HB3	1:A:480:LEU:HD12	1.92	0.51
1:A:276:ASN:O	1:A:278:LYS:N	2.42	0.51
1:A:385:ASN:O	1:A:389:VAL:HG12	2.11	0.50
1:A:52:LEU:HD12	1:A:136:LEU:HD23	1.92	0.50
1:A:29:ARG:HH11	1:A:177:ALA:HB2	1.76	0.50
1:A:223:THR:CG2	1:A:240:ARG:H	2.25	0.50
1:A:66:THR:CG2	1:A:69:ASP:H	2.25	0.50
1:A:273:GLN:HE22	1:A:314:HIS:CE1	2.30	0.49
1:A:443:GLU:O	1:A:446:GLN:HB3	2.11	0.49
1:A:126:SER:O	1:A:129:SER:HB2	2.12	0.49
1:A:406:ARG:C	1:A:408:GLY:H	2.19	0.49
1:A:267:LEU:CD2	1:A:287:ASP:HB2	2.42	0.49
1:A:273:GLN:HE22	1:A:314:HIS:HE1	1.60	0.48
1:A:334:LEU:HD21	1:A:353:PHE:HB2	1.96	0.48
1:A:366:LEU:HA	1:A:369:VAL:HG12	1.96	0.47
1:A:458:ASN:HD22	1:A:458:ASN:N	2.12	0.47
1:A:166:LEU:HD22	1:A:185:LEU:HA	1.95	0.47
1:A:107:GLY:N	1:A:113:LYS:HD2	2.30	0.47
1:A:506:GLN:HG3	1:A:532:LYS:HB2	1.96	0.47
1:A:551:ARG:HA	1:A:619:LYS:O	2.14	0.47
1:A:66:THR:H	1:A:69:ASP:HB2	1.80	0.47
1:A:421:LYS:H	1:A:424:HIS:CD2	2.30	0.47
1:A:108:ALA:CB	1:A:109:PRO:CD	2.90	0.47
1:A:460:LEU:O	1:A:464:GLU:HG3	2.14	0.47
1:A:544:VAL:HA	1:A:553:TYR:O	2.15	0.47
1:A:646:ILE:HG12	1:A:652:MET:HG2	1.98	0.46
1:A:379:LEU:O	1:A:383:PHE:HD1	1.99	0.46
1:A:625:PRO:HG3	1:A:667:LYS:HE3	1.98	0.46
1:A:70:VAL:HG21	1:A:120:TYR:CD1	2.51	0.46
1:A:107:GLY:HA2	1:A:113:LYS:HB2	1.98	0.46
1:A:73:LYS:HD2	1:A:74:TYR:CE2	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:549:THR:HG21	1:A:577:PRO:HG2	1.99	0.45
1:A:144:VAL:HB	1:A:145:PRO:HD2	1.98	0.45
1:A:316:PHE:HB2	1:A:486:ALA:HB2	1.98	0.45
1:A:89:MET:HE3	1:A:99:LEU:HB2	1.98	0.44
1:A:267:LEU:HD13	1:A:297:LEU:HD13	1.99	0.44
1:A:376:LEU:HD22	1:A:433:ILE:HG23	1.98	0.44
1:A:644:LEU:HD21	1:A:652:MET:HE3	1.99	0.44
1:A:273:GLN:NE2	1:A:314:HIS:HE1	2.15	0.44
1:A:355:ARG:O	1:A:359:ARG:HG2	2.18	0.44
1:A:270:ILE:HG22	1:A:274:MET:CE	2.48	0.43
1:A:438:SER:O	1:A:442:LYS:HG3	2.19	0.43
1:A:172:ILE:HG23	1:A:172:ILE:O	2.19	0.43
1:A:632:PRO:HA	1:A:633:PRO:HD3	1.92	0.43
1:A:73:LYS:HD3	1:A:151:LEU:HD11	2.00	0.43
1:A:195:LEU:HD23	1:A:199:PRO:HA	2.00	0.43
1:A:292:GLU:HB2	1:A:293:ASN:ND2	2.34	0.43
1:A:643:LEU:C	1:A:643:LEU:HD23	2.44	0.43
1:A:66:THR:HG23	1:A:67:VAL:N	2.33	0.42
1:A:211:GLN:HB3	1:A:480:LEU:CD1	2.49	0.42
1:A:366:LEU:O	1:A:369:VAL:HG12	2.20	0.42
1:A:736:ARG:HG2	1:A:737:VAL:N	2.34	0.42
1:A:191:ALA:HB1	1:A:206:LEU:HD22	2.02	0.42
1:A:717:VAL:HG12	1:A:726:THR:HB	2.01	0.42
1:A:599:ASP:CG	1:A:629:HIS:H	2.28	0.41
1:A:342:ASN:O	1:A:346:ILE:HG13	2.21	0.41
1:A:349:GLU:HA	1:A:352:VAL:HG12	2.02	0.41
1:A:453:SER:HB3	1:A:456:LYS:HB2	2.02	0.41
1:A:648:GLN:O	1:A:732:GLN:HG2	2.20	0.41
1:A:424:HIS:HA	1:A:429:GLU:HG3	2.03	0.41
1:A:642:MET:O	1:A:691:VAL:HA	2.19	0.41
1:A:613:ALA:HB2	1:A:619:LYS:HB2	2.02	0.41
1:A:34:ILE:O	1:A:164:ILE:HD11	2.21	0.41
1:A:34:ILE:HD13	1:A:349:GLU:HB3	2.03	0.41
1:A:132:LEU:HD23	1:A:159:ALA:HB2	2.02	0.41
1:A:131:LEU:HD21	1:A:161:LEU:HD12	2.03	0.40
1:A:83:TYR:OH	1:A:113:LYS:N	2.55	0.40
1:A:320:ASP:C	1:A:322:GLU:N	2.80	0.40
1:A:490:SER:C	1:A:492:HIS:N	2.79	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	A	677/716 (95%)	621 (92%)	45 (7%)	11 (2%)	7 34

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	109	PRO
1	A	496	PRO
1	A	173	ALA
1	A	249	ASP
1	A	108	ALA
1	A	277	LYS
1	A	411	GLU
1	A	659	ASP
1	A	683	GLU
1	A	291	HIS
1	A	499	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	583/605 (96%)	540 (93%)	43 (7%)	13 42

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ASP
1	A	41	VAL
1	A	73	LYS
1	A	80	SER
1	A	91	GLN
1	A	125	LEU
1	A	131	LEU
1	A	133	LEU
1	A	152	GLN
1	A	164	ILE
1	A	172	ILE
1	A	198	GLN
1	A	207	LEU
1	A	217	THR
1	A	228	ILE
1	A	249	ASP
1	A	257	LEU
1	A	261	VAL
1	A	293	ASN
1	A	295	THR
1	A	304	ASN
1	A	310	GLN
1	A	315	LEU
1	A	329	LEU
1	A	340	VAL
1	A	363	ASN
1	A	367	GLU
1	A	386	ILE
1	A	390	ILE
1	A	392	ILE
1	A	404	MET
1	A	412	THR
1	A	425	LEU
1	A	453	SER
1	A	458	ASN
1	A	481	GLN
1	A	484	GLU
1	A	506	GLN
1	A	514	LYS
1	A	536	LYS
1	A	609	ASN
1	A	647	THR
1	A	692	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	HIS
1	A	91	GLN
1	A	138	GLN
1	A	163	ASN
1	A	205	GLN
1	A	273	GLN
1	A	276	ASN
1	A	293	ASN
1	A	314	HIS
1	A	330	ASN
1	A	342	ASN
1	A	363	ASN
1	A	395	ASN
1	A	424	HIS
1	A	437	GLN
1	A	448	GLN
1	A	458	ASN
1	A	584	HIS
1	A	609	ASN
1	A	673	ASN
1	A	688	GLN
1	A	695	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 [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	685/716 (95%)	-0.09	11 (1%) 70 47	20, 37, 71, 78	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	112	PRO	3.6
1	A	120	TYR	3.4
1	A	563	ALA	3.0
1	A	171	GLY	2.8
1	A	531	PHE	2.6
1	A	108	ALA	2.5
1	A	110	ASP	2.5
1	A	555	ILE	2.2
1	A	173	ALA	2.2
1	A	93	PHE	2.1
1	A	704	LYS	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.