Full wwPDB X-ray Structure Validation Report (i)

Aug 15, 2023 – 06:10 AM EDT

PDB ID	:	1U2L
Title	:	Crystal structure of the C-terminal domain from the catalase-peroxidase KatG of Escherichia coli (P1)
Authors	:	Carpena, X.; Melik-Adamyan, W.; Loewen, P.C.; Fita, I.
Deposited on	:	2004-07-19
Resolution	:	2.30 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 (i) 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 (1)) were used in the production of this report:

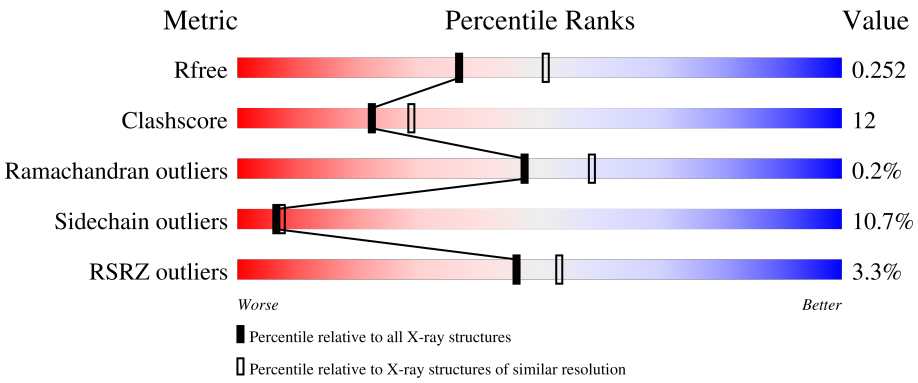
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
R_{free}	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 <=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	309	3% 70% 23% 6% •
1	B	309	4% 75% 22% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4931 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 Peroxidase/catalase HPI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	307	2360	1495	404	454	7	0	0	0
1	B	307	2360	1495	404	454	7	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	418	GLY	-	cloning artifact	UNP P13029
A	419	SER	-	cloning artifact	UNP P13029
A	420	HIS	-	cloning artifact	UNP P13029
A	421	MET	-	cloning artifact	UNP P13029
B	418	GLY	-	cloning artifact	UNP P13029
B	419	SER	-	cloning artifact	UNP P13029
B	420	HIS	-	cloning artifact	UNP P13029
B	421	MET	-	cloning artifact	UNP P13029

- Molecule 2 is water.

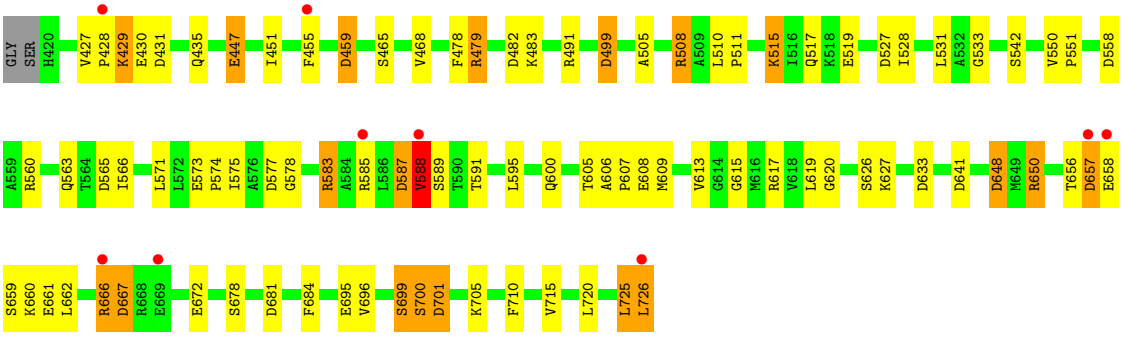
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	106	106	106	0	0
2	B	105	105	105	0	0

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.

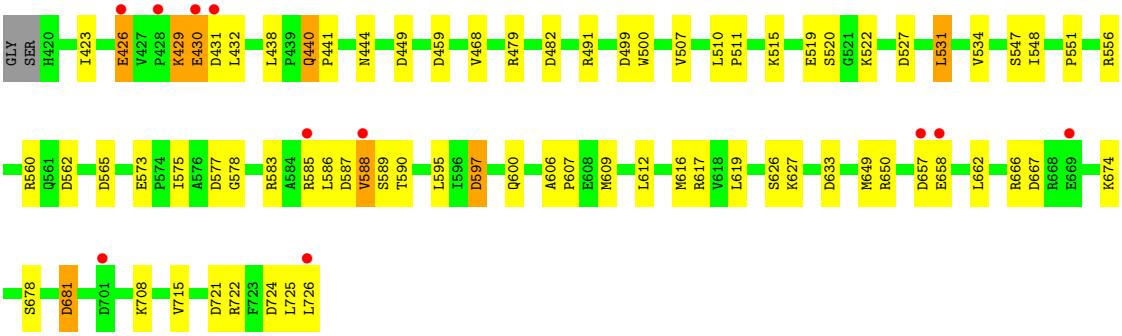
• Molecule 1: Peroxidase/catalase HPI

Chain A:



• Molecule 1: Peroxidase/catalase HPI

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	52.86Å 52.76Å 55.99Å 93.83° 90.71° 103.27°	Depositor
Resolution (Å)	25.00 – 2.30 24.84 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (25.00-2.30) 97.9 (24.84-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	3.81 (at 2.31Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.175 , 0.251 0.176 , 0.252	Depositor DCC
R_{free} test set	1304 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å^2)	19.4	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 52.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.016 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4931	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.39% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	0.43	1/2403 (0.0%)	0.73	14/3258 (0.4%)
1	B	0.42	1/2403 (0.0%)	0.72	13/3258 (0.4%)
All	All	0.43	2/4806 (0.0%)	0.73	27/6516 (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
1	A	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	430	GLU	CD-OE2	7.12	1.33	1.25
1	A	430	GLU	CD-OE2	6.88	1.33	1.25

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	681	ASP	CB-CG-OD2	6.49	124.14	118.30
1	A	577	ASP	CB-CG-OD2	5.90	123.61	118.30
1	B	724	ASP	CB-CG-OD2	5.82	123.54	118.30
1	A	701	ASP	CB-CG-OD2	5.78	123.50	118.30
1	A	641	ASP	CB-CG-OD2	5.72	123.45	118.30
1	B	449	ASP	CB-CG-OD2	5.72	123.45	118.30
1	A	633	ASP	CB-CG-OD2	5.70	123.43	118.30
1	B	565	ASP	CB-CG-OD2	5.63	123.37	118.30
1	A	667	ASP	CB-CG-OD2	5.53	123.28	118.30
1	A	459	ASP	CB-CG-OD2	5.50	123.25	118.30
1	A	431	ASP	CB-CG-OD2	5.50	123.25	118.30
1	B	459	ASP	CB-CG-OD2	5.48	123.23	118.30

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	721	ASP	CB-CG-OD2	5.46	123.21	118.30
1	B	633	ASP	CB-CG-OD2	5.44	123.20	118.30
1	A	657	ASP	CB-CG-OD2	5.40	123.16	118.30
1	A	499	ASP	CB-CG-OD2	5.39	123.15	118.30
1	B	562	ASP	CB-CG-OD2	5.39	123.15	118.30
1	B	431	ASP	CB-CG-OD2	5.35	123.11	118.30
1	A	558	ASP	CB-CG-OD2	5.30	123.07	118.30
1	A	565	ASP	CB-CG-OD2	5.28	123.05	118.30
1	B	587	ASP	CB-CG-OD2	5.23	123.01	118.30
1	B	667	ASP	CB-CG-OD2	5.19	122.97	118.30
1	B	597	ASP	CB-CG-OD2	5.17	122.95	118.30
1	B	577	ASP	CB-CG-OD2	5.17	122.95	118.30
1	A	648	ASP	CB-CG-OD2	5.13	122.92	118.30
1	A	587	ASP	CB-CG-OD2	5.07	122.87	118.30
1	A	527	ASP	CB-CG-OD2	5.04	122.83	118.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	588	VAL	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	2360	0	2341	65	0
1	B	2360	0	2341	45	0
2	A	106	0	0	15	0
2	B	105	0	0	7	0
All	All	4931	0	4682	110	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 12.

All (110) 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:479:ARG:HH11	1:A:479:ARG:CG	1.76	0.98
1:A:479:ARG:HH11	1:A:479:ARG:CB	1.81	0.93
1:A:447:GLU:HG3	2:A:152:HOH:O	1.69	0.90
1:B:595:LEU:HD11	1:B:609:MET:SD	2.16	0.86
1:A:479:ARG:HH11	1:A:479:ARG:HG2	1.41	0.85
1:A:505:ALA:O	1:A:508:ARG:HG3	1.82	0.79
1:A:666:ARG:HG2	1:A:666:ARG:HH11	1.49	0.77
1:A:701:ASP:CA	2:A:131:HOH:O	2.32	0.77
1:A:605:THR:OG1	1:A:608:GLU:HG3	1.85	0.76
1:B:595:LEU:CD1	1:B:609:MET:SD	2.75	0.75
1:A:701:ASP:N	2:A:131:HOH:O	2.21	0.73
1:A:701:ASP:HA	2:A:131:HOH:O	1.88	0.72
1:A:479:ARG:HG2	1:A:479:ARG:NH1	2.03	0.70
1:B:479:ARG:NH2	1:B:482:ASP:OD1	2.24	0.68
1:B:595:LEU:C	1:B:595:LEU:HD13	2.13	0.68
1:A:587:ASP:O	1:A:588:VAL:HB	1.94	0.67
1:B:726:LEU:HD21	2:B:99:HOH:O	1.95	0.65
1:B:468:VAL:HG21	1:B:619:LEU:HD13	1.80	0.64
1:A:479:ARG:CB	1:A:479:ARG:NH1	2.59	0.63
1:A:578:GLY:HA3	1:A:595:LEU:HD23	1.82	0.60
1:A:479:ARG:HH11	1:A:479:ARG:HB3	1.65	0.60
1:A:666:ARG:HH11	1:A:666:ARG:CG	2.15	0.59
1:B:666:ARG:HG3	1:B:666:ARG:NH1	2.18	0.58
1:A:726:LEU:HG	2:A:194:HOH:O	2.01	0.58
1:A:588:VAL:CG1	1:A:588:VAL:O	2.51	0.57
1:B:666:ARG:HG3	1:B:666:ARG:HH11	1.68	0.57
1:A:595:LEU:CD1	1:A:609:MET:SD	2.93	0.57
1:A:662:LEU:HD23	2:A:109:HOH:O	2.04	0.57
1:A:551:PRO:HD2	1:A:715:VAL:HG21	1.85	0.57
1:B:491:ARG:HG3	2:B:51:HOH:O	2.04	0.56
1:B:573:GLU:OE2	2:B:186:HOH:O	2.18	0.56
1:A:491:ARG:NH1	2:A:115:HOH:O	2.39	0.56
1:B:507:VAL:HG12	1:B:507:VAL:O	2.04	0.56
1:B:578:GLY:HA3	1:B:595:LEU:HD23	1.89	0.55
1:A:726:LEU:HA	2:A:128:HOH:O	2.06	0.55
2:A:181:HOH:O	1:B:590:THR:HG23	2.06	0.55
1:B:438:LEU:CD2	2:B:93:HOH:O	2.55	0.55
1:B:678:SER:HB2	2:B:182:HOH:O	2.06	0.54
1:B:430:GLU:HG2	2:B:204:HOH:O	2.06	0.54
1:B:595:LEU:HD13	1:B:595:LEU:O	2.08	0.54
1:A:661:GLU:OE1	2:A:27:HOH:O	2.19	0.53
1:B:588:VAL:CG1	1:B:589:SER:N	2.70	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:612:LEU:O	1:B:616:MET:HG3	2.09	0.53
1:A:650:ARG:NE	1:A:667:ASP:OD2	2.43	0.52
1:B:520:SER:O	1:B:522:LYS:HG2	2.10	0.52
1:A:515:LYS:HE3	1:A:519:GLU:OE2	2.09	0.51
1:A:510:LEU:N	1:A:511:PRO:CD	2.73	0.51
1:A:560:ARG:HG2	1:A:563:GLN:HG3	1.91	0.51
1:A:666:ARG:CG	1:A:666:ARG:NH1	2.73	0.51
1:B:657:ASP:C	1:B:657:ASP:OD2	2.49	0.51
1:B:500:TRP:CD2	1:B:573:GLU:HG3	2.46	0.50
1:A:606:ALA:HB3	1:A:607:PRO:HD3	1.94	0.50
1:A:725:LEU:O	1:A:726:LEU:C	2.49	0.49
1:B:507:VAL:O	1:B:507:VAL:CG1	2.60	0.49
1:A:620:GLY:HA2	2:A:14:HOH:O	2.12	0.49
1:A:609:MET:O	1:A:613:VAL:HG23	2.13	0.49
1:A:617:ARG:NE	1:A:681:ASP:OD1	2.45	0.48
1:A:678:SER:HB2	2:A:27:HOH:O	2.13	0.48
1:B:510:LEU:HB2	1:B:511:PRO:HD3	1.94	0.48
1:B:429:LYS:HD2	1:B:432:LEU:CD1	2.43	0.48
1:B:510:LEU:N	1:B:511:PRO:CD	2.76	0.48
1:A:595:LEU:HD11	1:A:609:MET:SD	2.53	0.47
1:A:451:ILE:HG23	1:A:455:PHE:CE1	2.50	0.47
1:B:429:LYS:HA	1:B:429:LYS:HD3	1.56	0.47
1:A:468:VAL:HG21	1:A:619:LEU:HD13	1.96	0.47
1:B:500:TRP:CG	1:B:573:GLU:HG3	2.50	0.47
1:A:505:ALA:O	1:A:508:ARG:CG	2.59	0.47
1:B:510:LEU:N	1:B:511:PRO:HD2	2.29	0.47
1:A:613:VAL:HG21	1:A:684:PHE:CZ	2.51	0.46
1:A:510:LEU:HB2	1:A:511:PRO:HD3	1.98	0.46
1:B:440:GLN:NE2	1:B:441:PRO:O	2.46	0.46
1:A:666:ARG:CZ	1:A:666:ARG:HB3	2.47	0.45
1:B:627:LYS:N	1:B:627:LYS:HD2	2.30	0.45
1:B:429:LYS:HD2	1:B:432:LEU:HD11	1.99	0.45
1:B:527:ASP:OD1	1:B:556:ARG:HB2	2.17	0.45
1:A:429:LYS:HA	1:A:429:LYS:HD2	1.68	0.45
1:B:531:LEU:O	1:B:534:VAL:HB	2.17	0.45
1:B:595:LEU:CD1	1:B:595:LEU:C	2.84	0.45
1:B:551:PRO:HD2	1:B:715:VAL:HG21	1.99	0.44
1:A:575:ILE:HD13	2:A:83:HOH:O	2.17	0.44
1:A:595:LEU:HD13	1:A:609:MET:SD	2.57	0.44
1:A:619:LEU:HD23	1:A:619:LEU:HA	1.75	0.44
1:A:427:VAL:HG13	2:A:199:HOH:O	2.17	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:648:ASP:OD2	1:A:650:ARG:HB3	2.18	0.44
1:A:656:THR:O	1:A:657:ASP:HB3	2.18	0.44
1:A:615:GLY:HA3	1:A:710:PHE:CE1	2.53	0.43
1:B:606:ALA:HB3	1:B:607:PRO:HD3	2.00	0.43
1:A:583:ARG:HH21	1:A:591:THR:N	2.16	0.43
1:A:700:SER:O	1:A:701:ASP:HB2	2.17	0.43
1:A:427:VAL:HA	1:A:428:PRO:HD3	1.88	0.43
1:B:626:SER:HB3	2:B:150:HOH:O	2.18	0.43
1:A:451:ILE:CG2	1:A:455:PHE:HE1	2.32	0.43
1:B:606:ALA:N	1:B:607:PRO:CD	2.82	0.43
1:B:426:GLU:O	1:B:426:GLU:HG2	2.18	0.42
1:A:482:ASP:O	1:A:483:LYS:HB2	2.18	0.42
1:A:573:GLU:HA	1:A:574:PRO:HD3	1.85	0.42
1:B:588:VAL:HG12	1:B:589:SER:N	2.35	0.42
1:A:479:ARG:NH1	1:A:482:ASP:OD2	2.53	0.42
1:A:533:GLY:HA3	2:A:140:HOH:O	2.19	0.42
1:A:696:VAL:O	1:A:699:SER:HB2	2.19	0.42
1:A:478:PHE:CD2	1:A:720:LEU:HD13	2.54	0.42
1:B:438:LEU:N	1:B:438:LEU:HD23	2.33	0.41
1:B:617:ARG:NE	1:B:681:ASP:OD1	2.47	0.41
1:A:666:ARG:HG2	1:A:666:ARG:NH1	2.25	0.41
1:A:606:ALA:N	1:A:607:PRO:CD	2.84	0.41
1:A:517:GLN:CG	1:A:528:ILE:HD12	2.51	0.41
1:B:722:ARG:HB3	1:B:725:LEU:HD12	2.02	0.41
1:A:659:SER:O	1:A:660:LYS:HB2	2.19	0.41
1:A:550:VAL:HA	1:A:551:PRO:HD3	1.92	0.40
1:B:657:ASP:OD2	1:B:658:GLU:N	2.54	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	305/309 (99%)	300 (98%)	4 (1%)	1 (0%)	41	50
1	B	305/309 (99%)	297 (97%)	8 (3%)	0	100	100
All	All	610/618 (99%)	597 (98%)	12 (2%)	1 (0%)	47	58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	588	VAL

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	248/250 (99%)	219 (88%)	29 (12%)	5	6
1	B	248/250 (99%)	224 (90%)	24 (10%)	8	9
All	All	496/500 (99%)	443 (89%)	53 (11%)	6	7

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	429	LYS
1	A	435	GLN
1	A	447	GLU
1	A	459	ASP
1	A	465	SER
1	A	479	ARG
1	A	499	ASP
1	A	508	ARG
1	A	515	LYS
1	A	531	LEU
1	A	542	SER
1	A	566	ILE
1	A	571	LEU
1	A	583	ARG
1	A	585	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	589	SER
1	A	600	GLN
1	A	626	SER
1	A	627	LYS
1	A	650	ARG
1	A	658	GLU
1	A	666	ARG
1	A	672	GLU
1	A	695	GLU
1	A	699	SER
1	A	700	SER
1	A	705	LYS
1	A	725	LEU
1	A	726	LEU
1	B	423	ILE
1	B	426	GLU
1	B	429	LYS
1	B	440	GLN
1	B	444	ASN
1	B	499	ASP
1	B	515	LYS
1	B	519	GLU
1	B	531	LEU
1	B	547	SER
1	B	548	ILE
1	B	560	ARG
1	B	575	ILE
1	B	583	ARG
1	B	585	ARG
1	B	586	LEU
1	B	588	VAL
1	B	597	ASP
1	B	600	GLN
1	B	649	MET
1	B	650	ARG
1	B	662	LEU
1	B	674	LYS
1	B	708	LYS

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	517	GLN
1	A	600	GLN
1	A	601	GLN
1	A	628	ASN
1	B	444	ASN
1	B	600	GLN
1	B	601	GLN
1	B	628	ASN

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 monosaccharides 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	307/309 (99%)	0.07	9 (2%)	51	58	11, 24, 46, 58	0
1	B	307/309 (99%)	0.04	11 (3%)	42	49	10, 22, 47, 61	0
All	All	614/618 (99%)	0.06	20 (3%)	46	53	10, 23, 46, 61	0

All (20) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	588	VAL	4.3
1	B	657	ASP	4.0
1	A	658	GLU	3.3
1	A	669	GLU	3.1
1	A	726	LEU	3.1
1	A	588	VAL	3.0
1	B	726	LEU	2.9
1	B	669	GLU	2.7
1	B	430	GLU	2.6
1	B	431	ASP	2.6
1	B	428	PRO	2.6
1	A	455	PHE	2.6
1	A	428	PRO	2.5
1	B	585	ARG	2.5
1	B	658	GLU	2.4
1	A	585	ARG	2.2
1	A	657	ASP	2.2
1	B	426	GLU	2.1
1	B	701	ASP	2.1
1	A	666	ARG	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 monosaccharides 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.