



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 02:15 am BST

PDB ID : 1C8B
Title : CRYSTAL STRUCTURE OF A NOVEL GERMINATION PROTEASE FROM SPORES OF BACILLUS MEGATERIUM: STRUCTURAL REARRANGEMENTS AND ZYMOGEN ACTIVATION
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Deposited on : 2000-05-03
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

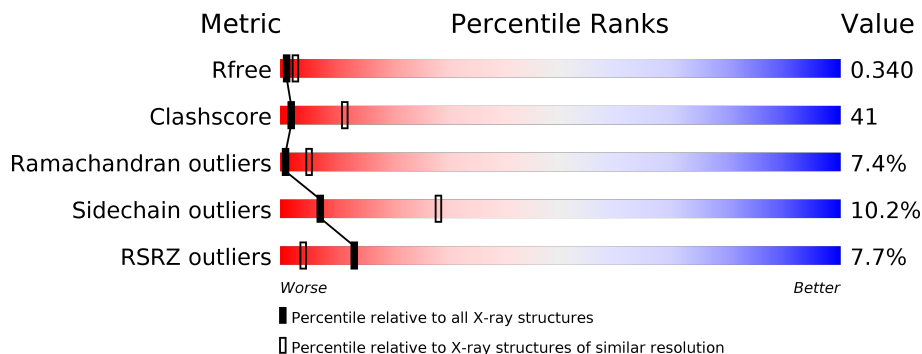
1 Overall quality at a glance

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}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (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 on 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	371	 7% 40% 38% 8% 14%
1	B	371	 6% 41% 38% 8% 14%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4454 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 SPORE PROTEASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	320	2164	1354	372	430	8	0	0	0
1	B	320	2164	1354	372	430	8	0	0	0

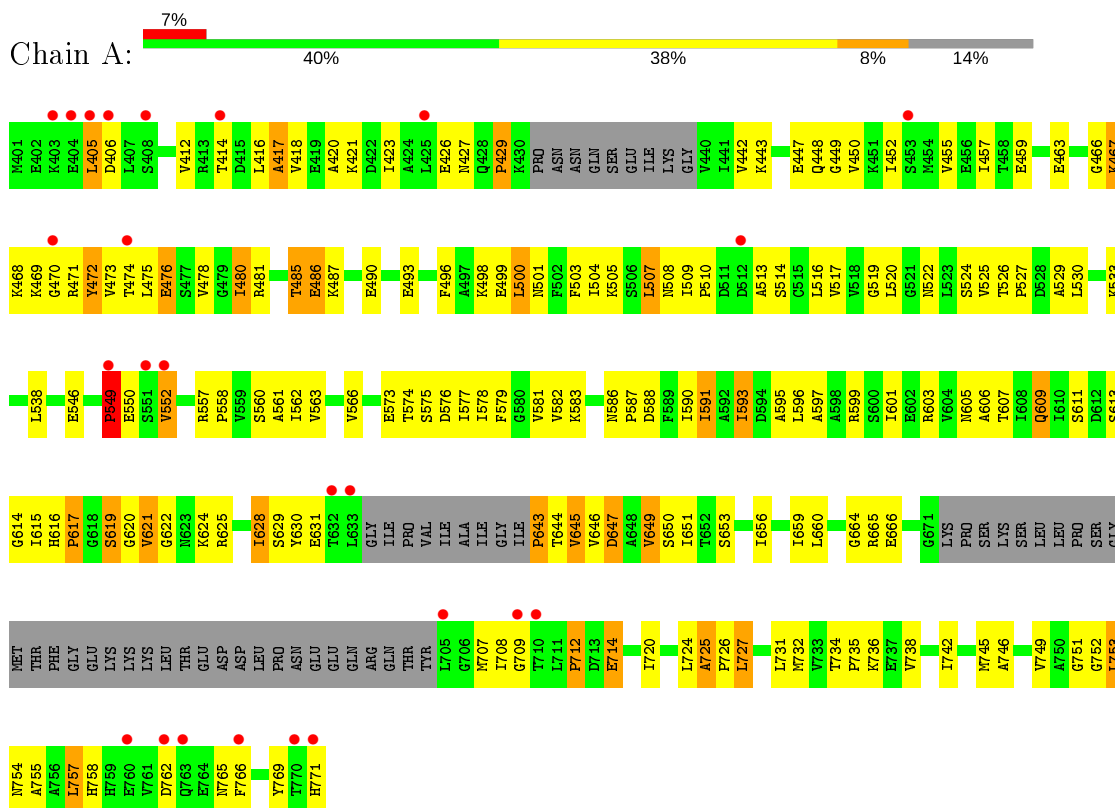
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	63	63	63	0	0
2	B	63	63	63	0	0

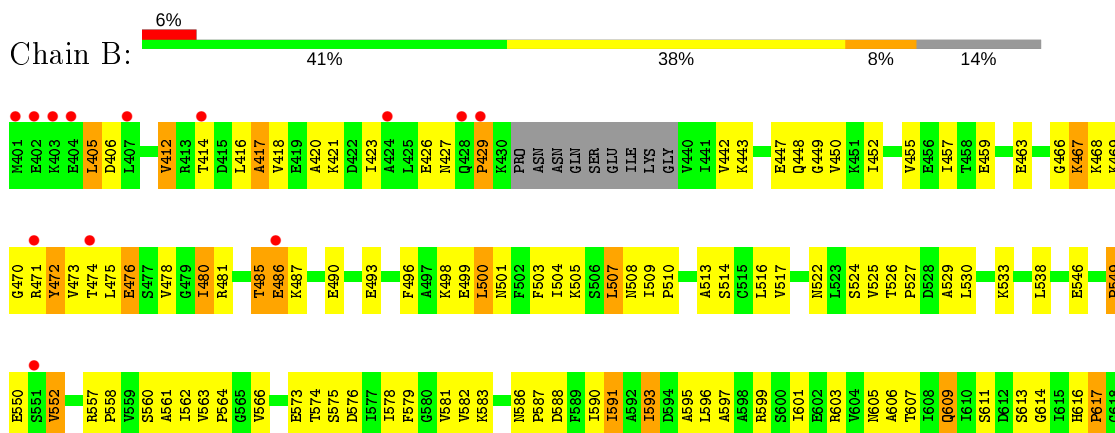
3 Residue-property plots

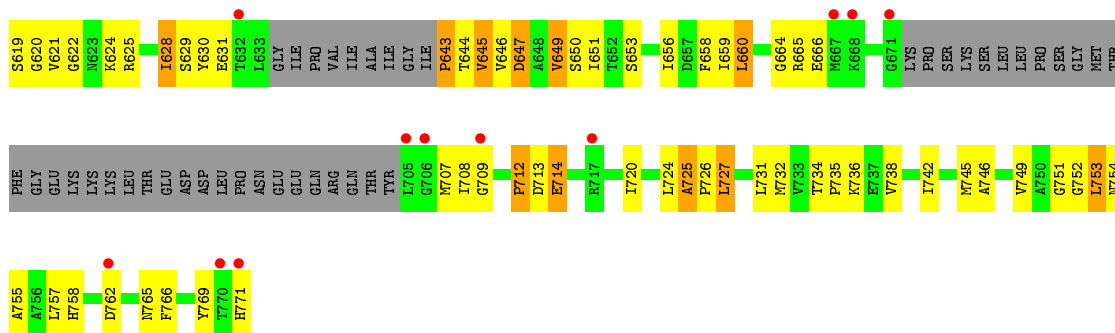
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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: SPORE PROTEASE



- Molecule 1: SPORE PROTEASE





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	77.41Å 77.41Å 313.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 3.00 48.75 – 3.00	Depositor EDS
% Data completeness (in resolution range)	88.0 (15.00-3.00) 92.2 (48.75-3.00)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.25 (at 3.01Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.308 , 0.330 0.332 , 0.340	Depositor DCC
R_{free} test set	1927 reflections (9.81%)	wwPDB-VP
Wilson B-factor (Å ²)	69.5	Xtrriage
Anisotropy	0.089	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 160.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	4454	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 24.33 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 3.9133e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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	0.63	0/2184	0.85	4/2976 (0.1%)
1	B	0.69	0/2184	0.87	3/2976 (0.1%)
All	All	0.66	0/4368	0.86	7/5952 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	712	PRO	N-CA-CB	6.68	111.31	103.30
1	A	712	PRO	N-CA-CB	6.58	111.19	103.30
1	A	643	PRO	N-CA-CB	5.73	110.18	103.30
1	B	643	PRO	N-CA-CB	5.47	109.86	103.30
1	B	660	LEU	CA-CB-CG	-5.19	103.36	115.30
1	A	549	PRO	N-CA-CB	5.16	109.49	103.30
1	A	660	LEU	CA-CB-CG	-5.16	103.44	115.30

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	2164	0	1943	173	5
1	B	2164	0	1943	173	8
2	A	63	0	0	12	14
2	B	63	0	0	10	11

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4454	0	3886	338	19

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 41.

All (338) 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:442:VAL:HG12	1:B:443:LYS:H	1.12	1.13
1:A:442:VAL:HG12	1:A:443:LYS:H	1.13	1.09
1:A:510:PRO:HG2	1:A:513:ALA:HB2	1.39	1.04
1:B:510:PRO:HG2	1:B:513:ALA:HB2	1.39	1.03
1:B:416:LEU:HA	2:B:106:HOH:O	1.61	0.99
1:A:463:GLU:HA	1:A:473:VAL:HG11	1.41	0.97
1:B:463:GLU:HA	1:B:473:VAL:HG11	1.43	0.97
1:B:481:ARG:O	1:B:606:ALA:HB1	1.66	0.95
1:A:481:ARG:O	1:A:606:ALA:HB1	1.66	0.95
1:B:530:LEU:CD1	1:B:593:ILE:HD11	1.96	0.95
1:A:530:LEU:CD1	1:A:593:ILE:HD11	1.98	0.92
1:A:725:ALA:HB1	1:A:726:PRO:HD3	1.53	0.89
1:B:725:ALA:HB1	1:B:726:PRO:HD3	1.52	0.88
1:B:522:ASN:HD21	1:B:524:SER:HB2	1.38	0.87
1:A:522:ASN:HD21	1:A:524:SER:HB2	1.38	0.86
1:B:442:VAL:HG12	1:B:443:LYS:N	1.91	0.85
1:A:442:VAL:HG12	1:A:443:LYS:N	1.92	0.84
1:A:516:LEU:HB3	1:A:590:ILE:HG22	1.59	0.84
1:A:579:PHE:O	1:A:582:VAL:HG22	1.79	0.82
1:B:725:ALA:HB1	1:B:726:PRO:CD	2.08	0.82
1:B:516:LEU:HB3	1:B:590:ILE:HG22	1.62	0.82
1:B:579:PHE:O	1:B:582:VAL:HG22	1.79	0.82
1:A:725:ALA:HB1	1:A:726:PRO:CD	2.09	0.81
1:A:522:ASN:ND2	1:A:524:SER:HB2	1.96	0.80
1:A:563:VAL:O	1:A:566:VAL:HG22	1.83	0.78
1:B:522:ASN:ND2	1:B:524:SER:HB2	1.97	0.78
1:B:563:VAL:O	1:B:566:VAL:HG22	1.84	0.77
1:A:476:GLU:HG3	1:A:611:SER:HB3	1.66	0.76
1:B:476:GLU:HG3	1:B:611:SER:HB3	1.66	0.76
1:B:575:SER:HA	1:B:578:ILE:HG12	1.67	0.75
1:B:614:GLY:HA3	1:B:628:ILE:HG22	1.69	0.75
1:A:651:ILE:HG21	1:A:745:MET:HG2	1.69	0.74
1:A:514:SER:HA	1:A:558:PRO:HG2	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:591:ILE:HB	1:A:647:ASP:HB3	1.70	0.74
1:A:414:THR:HA	2:A:101:HOH:O	1.87	0.74
1:B:414:THR:HA	2:B:1101:HOH:O	1.86	0.74
1:A:575:SER:HA	1:A:578:ILE:HG12	1.67	0.73
1:A:417:ALA:HB2	1:A:609:GLN:OE1	1.89	0.73
1:B:417:ALA:HB2	1:B:609:GLN:OE1	1.89	0.73
1:B:514:SER:HA	1:B:558:PRO:HG2	1.70	0.73
1:A:614:GLY:HA3	1:A:628:ILE:HG22	1.69	0.73
1:B:651:ILE:HG21	1:B:745:MET:HG2	1.71	0.72
1:B:530:LEU:HD13	1:B:593:ILE:HD11	1.71	0.72
1:B:476:GLU:HG3	1:B:611:SER:CB	2.20	0.72
1:A:476:GLU:HG3	1:A:611:SER:CB	2.20	0.72
1:B:505:LYS:O	1:B:508:ASN:N	2.21	0.71
1:B:562:ILE:HD13	1:B:581:VAL:HG21	1.71	0.71
1:B:591:ILE:HB	1:B:647:ASP:HB3	1.73	0.71
1:A:562:ILE:HD13	1:A:581:VAL:HG21	1.72	0.71
1:A:505:LYS:O	1:A:508:ASN:N	2.21	0.71
1:A:530:LEU:HD13	1:A:593:ILE:HD11	1.72	0.71
1:A:651:ILE:HD12	1:A:651:ILE:N	2.06	0.71
1:A:659:ILE:HB	1:A:720:ILE:HD12	1.73	0.70
1:B:651:ILE:N	1:B:651:ILE:HD12	2.06	0.70
1:B:653:SER:O	1:B:734:THR:HG22	1.92	0.70
1:A:517:VAL:HB	1:A:561:ALA:HB2	1.73	0.69
1:B:448:GLN:HB3	1:B:457:ILE:HG23	1.74	0.69
1:B:517:VAL:HB	1:B:561:ALA:HB2	1.75	0.68
1:A:653:SER:O	1:A:734:THR:HG22	1.92	0.68
1:A:448:GLN:HB3	1:A:457:ILE:HG23	1.75	0.68
1:B:738:VAL:O	1:B:742:ILE:HG12	1.93	0.67
1:A:738:VAL:O	1:A:742:ILE:HG12	1.95	0.67
1:B:593:ILE:HD13	1:B:749:VAL:HG11	1.76	0.67
1:A:513:ALA:HB3	1:A:557:ARG:HH12	1.60	0.66
1:A:517:VAL:HB	1:A:561:ALA:CB	2.25	0.66
1:A:659:ILE:HA	2:A:116:HOH:O	1.95	0.66
1:A:593:ILE:HD13	1:A:749:VAL:HG11	1.76	0.66
1:B:659:ILE:HB	1:B:720:ILE:HD12	1.76	0.66
1:B:480:ILE:HG22	2:B:1107:HOH:O	1.95	0.65
1:A:480:ILE:HG22	2:A:107:HOH:O	1.95	0.65
1:B:442:VAL:CG1	1:B:443:LYS:H	1.97	0.65
1:A:563:VAL:O	1:A:563:VAL:HG13	1.96	0.65
1:A:727:LEU:HD12	1:A:731:LEU:HD13	1.79	0.64
1:A:498:LYS:HE3	1:A:499:GLU:OE2	1.97	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:538:LEU:HD11	1:A:753:LEU:HD12	1.78	0.64
1:B:517:VAL:HB	1:B:561:ALA:CB	2.27	0.64
1:B:659:ILE:HA	2:B:1116:HOH:O	1.97	0.64
1:B:727:LEU:HD12	1:B:731:LEU:HD13	1.79	0.64
1:B:498:LYS:HE3	1:B:499:GLU:OE2	1.98	0.64
1:B:449:GLY:HA3	2:B:1110:HOH:O	1.97	0.64
1:B:507:LEU:HD13	1:B:509:ILE:HD11	1.80	0.64
1:B:513:ALA:HB3	1:B:557:ARG:HH12	1.60	0.64
1:A:591:ILE:HD11	1:A:753:LEU:HD11	1.80	0.64
1:A:427:ASN:O	1:A:429:PRO:HD3	1.98	0.63
1:B:563:VAL:O	1:B:563:VAL:HG13	1.96	0.63
1:A:480:ILE:HG21	1:A:603:ARG:NH2	2.15	0.62
1:B:605:ASN:ND2	1:B:735:PRO:HG2	2.15	0.62
1:A:507:LEU:HD13	1:A:509:ILE:HD11	1.82	0.62
1:B:427:ASN:O	1:B:429:PRO:HD3	2.00	0.62
1:B:480:ILE:HG21	1:B:603:ARG:NH2	2.15	0.62
1:A:619:SER:HA	2:A:1131:HOH:O	1.99	0.61
1:B:591:ILE:HD11	1:B:753:LEU:HD11	1.82	0.61
1:A:449:GLY:HA3	2:A:110:HOH:O	1.99	0.61
1:B:725:ALA:CB	1:B:726:PRO:CD	2.79	0.61
1:A:442:VAL:CG1	1:A:443:LYS:H	1.98	0.61
1:B:538:LEU:HD11	1:B:753:LEU:HD12	1.81	0.61
1:A:455:VAL:HB	1:A:480:ILE:HG23	1.83	0.60
1:A:605:ASN:ND2	1:A:735:PRO:HG2	2.17	0.60
1:A:725:ALA:CB	1:A:726:PRO:CD	2.80	0.60
1:A:504:ILE:HG22	1:A:509:ILE:HB	1.84	0.60
1:B:504:ILE:HG22	1:B:509:ILE:HB	1.83	0.60
1:A:628:ILE:HD13	1:A:628:ILE:H	1.67	0.59
1:B:455:VAL:HB	1:B:480:ILE:HG23	1.83	0.59
1:B:575:SER:OG	1:B:625:ARG:HD2	2.02	0.59
1:B:423:ILE:O	1:B:426:GLU:HB3	2.03	0.59
1:B:596:LEU:HG	1:B:650:SER:HB2	1.84	0.59
1:A:726:PRO:O	1:A:727:LEU:HG	2.03	0.59
1:B:644:THR:O	1:B:646:VAL:HG23	2.03	0.59
1:B:628:ILE:HD13	1:B:631:GLU:CB	2.34	0.58
1:B:405:LEU:HD23	1:B:405:LEU:H	1.67	0.58
1:B:582:VAL:HG23	1:B:583:LYS:N	2.18	0.58
1:A:478:VAL:HG22	1:A:609:GLN:HG3	1.86	0.58
1:B:629:SER:O	1:B:643:PRO:HA	2.04	0.57
1:B:749:VAL:O	1:B:753:LEU:HB2	2.04	0.57
1:A:405:LEU:HD23	1:A:405:LEU:H	1.68	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:510:PRO:HG2	1:A:513:ALA:CB	2.26	0.57
1:A:596:LEU:HG	1:A:650:SER:HB2	1.86	0.57
1:A:516:LEU:HD23	1:A:582:VAL:HG12	1.86	0.57
1:B:414:THR:O	1:B:599:ARG:HB2	2.05	0.57
1:B:478:VAL:HG22	1:B:609:GLN:HG3	1.85	0.57
1:A:644:THR:O	1:A:646:VAL:HG23	2.04	0.57
1:B:628:ILE:H	1:B:628:ILE:HD13	1.68	0.57
1:A:714:GLU:OE1	1:A:714:GLU:HA	2.05	0.57
1:B:487:LYS:O	1:B:490:GLU:HG2	2.05	0.57
1:B:714:GLU:HA	1:B:714:GLU:OE1	2.04	0.57
1:B:470:GLY:O	1:B:472:TYR:N	2.38	0.57
1:A:575:SER:OG	1:A:625:ARG:HD2	2.05	0.57
1:A:582:VAL:HG23	1:A:583:LYS:N	2.19	0.57
1:B:463:GLU:HA	1:B:473:VAL:CG1	2.27	0.56
1:A:423:ILE:O	1:A:426:GLU:HB3	2.05	0.56
1:A:470:GLY:O	1:A:472:TYR:N	2.39	0.56
1:B:726:PRO:O	1:B:727:LEU:HG	2.04	0.56
1:A:529:ALA:O	1:A:533:LYS:HG3	2.06	0.56
1:B:516:LEU:HD23	1:B:582:VAL:HG12	1.87	0.56
1:A:628:ILE:HD13	1:A:631:GLU:CB	2.36	0.56
1:A:629:SER:O	1:A:643:PRO:HA	2.05	0.56
1:A:414:THR:O	1:A:599:ARG:HB2	2.05	0.56
1:B:510:PRO:HG2	1:B:513:ALA:CB	2.27	0.56
1:A:463:GLU:HA	1:A:473:VAL:CG1	2.25	0.55
1:B:582:VAL:O	1:B:586:ASN:N	2.37	0.55
1:A:664:GLY:C	1:A:666:GLU:H	2.09	0.55
1:B:664:GLY:C	1:B:666:GLU:H	2.10	0.55
1:B:562:ILE:CD1	1:B:581:VAL:HG21	2.35	0.55
1:A:749:VAL:O	1:A:753:LEU:HB2	2.07	0.55
1:A:487:LYS:O	1:A:490:GLU:HG2	2.07	0.54
1:B:613:SER:HB3	1:B:616:HIS:NE2	2.22	0.54
1:A:735:PRO:HA	1:B:731:LEU:HD12	1.88	0.54
1:A:562:ILE:CD1	1:A:581:VAL:HG21	2.37	0.54
1:A:514:SER:HB3	1:A:588:ASP:H	1.72	0.54
1:A:517:VAL:HG21	1:A:538:LEU:HD21	1.90	0.54
1:B:514:SER:HB3	1:B:588:ASP:H	1.72	0.54
1:A:522:ASN:HD21	1:A:524:SER:CB	2.16	0.54
1:A:582:VAL:O	1:A:586:ASN:N	2.37	0.54
1:B:529:ALA:O	1:B:533:LYS:HG3	2.08	0.54
1:B:574:THR:O	1:B:578:ILE:HG23	2.08	0.54
1:B:599:ARG:HG2	1:B:599:ARG:HH11	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:530:LEU:HD11	1:B:593:ILE:HD11	1.88	0.53
1:A:513:ALA:HB3	1:A:557:ARG:NH1	2.22	0.53
1:A:499:GLU:O	1:A:503:PHE:HB2	2.08	0.53
1:A:613:SER:HB3	1:A:616:HIS:NE2	2.24	0.53
1:A:496:PHE:HD2	1:A:752:GLY:HA3	1.73	0.53
1:A:485:THR:O	1:A:486:GLU:HB3	2.09	0.53
1:A:651:ILE:CD1	1:A:651:ILE:N	2.72	0.53
1:A:574:THR:O	1:A:578:ILE:HG23	2.08	0.53
1:B:485:THR:O	1:B:486:GLU:HB3	2.08	0.53
1:B:499:GLU:O	1:B:503:PHE:HB2	2.09	0.53
1:B:751:GLY:HA2	1:B:754:ASN:HD22	1.74	0.53
1:B:563:VAL:HG13	1:B:566:VAL:CG2	2.39	0.53
1:B:450:VAL:HG22	1:B:455:VAL:HG22	1.91	0.53
1:A:726:PRO:C	1:A:727:LEU:HG	2.29	0.53
1:A:448:GLN:HB3	1:A:457:ILE:HG12	1.91	0.52
1:A:450:VAL:HG22	1:A:455:VAL:HG22	1.91	0.52
1:B:513:ALA:HB3	1:B:557:ARG:NH1	2.23	0.52
1:B:550:GLU:N	2:B:1100:HOH:O	2.43	0.52
1:A:563:VAL:HG13	1:A:566:VAL:CG2	2.39	0.52
1:B:496:PHE:HD2	1:B:752:GLY:HA3	1.74	0.52
1:A:599:ARG:HH11	1:A:599:ARG:HG2	1.75	0.52
1:A:664:GLY:C	1:A:666:GLU:N	2.63	0.52
1:B:412:VAL:HB	2:B:190:HOH:O	2.09	0.52
1:A:550:GLU:N	2:A:100:HOH:O	2.42	0.52
1:B:726:PRO:C	1:B:727:LEU:HG	2.31	0.52
1:A:751:GLY:HA2	1:A:754:ASN:HD22	1.76	0.51
1:B:651:ILE:CD1	1:B:651:ILE:N	2.72	0.51
1:A:546:GLU:O	1:A:549:PRO:N	2.44	0.51
1:A:557:ARG:HG2	1:A:558:PRO:N	2.25	0.51
1:B:522:ASN:HD21	1:B:524:SER:CB	2.18	0.51
1:B:595:ALA:HA	1:B:651:ILE:O	2.11	0.51
1:B:448:GLN:HB3	1:B:457:ILE:HG12	1.92	0.51
1:B:517:VAL:HG21	1:B:538:LEU:HD21	1.92	0.51
1:B:587:PRO:HG3	1:B:590:ILE:HG22	1.92	0.51
1:B:557:ARG:HG2	1:B:558:PRO:N	2.26	0.51
1:B:596:LEU:HD11	1:B:650:SER:OG	2.11	0.50
1:B:664:GLY:C	1:B:666:GLU:N	2.64	0.50
1:B:546:GLU:O	1:B:549:PRO:N	2.44	0.50
1:B:724:LEU:O	1:B:725:ALA:C	2.50	0.50
1:A:591:ILE:HB	1:A:647:ASP:CB	2.40	0.50
1:A:525:VAL:HG12	1:A:527:PRO:HD2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:582:VAL:CG2	1:B:583:LYS:N	2.74	0.50
1:A:452:ILE:HG22	1:A:452:ILE:O	2.11	0.50
1:B:504:ILE:CG2	1:B:509:ILE:HB	2.42	0.50
1:A:595:ALA:HA	1:A:651:ILE:O	2.12	0.49
1:A:621:VAL:HA	1:B:736:LYS:HD3	1.94	0.49
1:A:480:ILE:HG21	1:A:603:ARG:CZ	2.42	0.49
1:A:418:VAL:C	1:A:420:ALA:N	2.66	0.49
1:A:617:PRO:HA	1:A:624:LYS:O	2.12	0.49
1:B:513:ALA:H	1:B:557:ARG:NH1	2.10	0.49
1:A:418:VAL:O	1:A:421:LYS:N	2.44	0.49
1:A:659:ILE:HD12	1:A:724:LEU:HG	1.94	0.49
1:A:582:VAL:CG2	1:A:583:LYS:N	2.76	0.49
1:A:738:VAL:HG22	1:A:738:VAL:O	2.12	0.49
1:B:452:ILE:HG22	1:B:452:ILE:O	2.12	0.49
1:B:525:VAL:HG12	1:B:527:PRO:HD2	1.94	0.49
1:B:738:VAL:HG22	1:B:738:VAL:O	2.13	0.49
1:B:480:ILE:HG21	1:B:603:ARG:CZ	2.42	0.49
1:A:596:LEU:HD11	1:A:650:SER:OG	2.13	0.48
1:A:650:SER:C	1:A:651:ILE:HD12	2.33	0.48
1:A:724:LEU:O	1:A:725:ALA:C	2.50	0.48
1:B:573:GLU:O	1:B:576:ASP:HB2	2.14	0.48
1:A:504:ILE:CG2	1:A:509:ILE:HB	2.43	0.48
1:B:591:ILE:HB	1:B:647:ASP:CB	2.43	0.48
1:A:557:ARG:H	1:A:757:LEU:HD23	1.78	0.48
1:B:659:ILE:HD12	1:B:724:LEU:HG	1.93	0.48
1:A:587:PRO:HG3	1:A:590:ILE:HG22	1.95	0.48
1:B:617:PRO:HA	1:B:624:LYS:O	2.13	0.48
1:A:480:ILE:HG21	1:A:603:ARG:HH22	1.77	0.48
1:A:573:GLU:O	1:A:576:ASP:HB2	2.14	0.48
1:B:557:ARG:H	1:B:757:LEU:HD23	1.78	0.48
1:B:628:ILE:HG12	1:B:630:TYR:N	2.28	0.48
1:A:516:LEU:HA	1:A:560:SER:O	2.14	0.47
1:A:601:ILE:HB	1:B:725:ALA:HB3	1.96	0.47
1:B:476:GLU:HG3	1:B:611:SER:HB2	1.97	0.47
1:A:732:MET:CE	1:B:736:LYS:HA	2.44	0.47
1:B:480:ILE:HG21	1:B:603:ARG:HH22	1.78	0.47
1:A:496:PHE:CD2	1:A:752:GLY:HA3	2.48	0.47
1:B:496:PHE:CD2	1:B:752:GLY:HA3	2.49	0.47
1:A:527:PRO:HB3	1:A:738:VAL:HG21	1.97	0.47
1:B:614:GLY:CA	1:B:628:ILE:HG22	2.42	0.47
1:B:480:ILE:CG2	1:B:603:ARG:NH1	2.78	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:650:SER:C	1:B:651:ILE:HD12	2.35	0.46
1:A:552:VAL:O	1:A:552:VAL:HG12	2.16	0.46
1:A:513:ALA:H	1:A:557:ARG:NH1	2.13	0.46
1:A:480:ILE:CG2	1:A:603:ARG:NH1	2.78	0.46
1:B:418:VAL:C	1:B:420:ALA:N	2.66	0.46
1:B:664:GLY:O	1:B:666:GLU:N	2.49	0.46
1:A:573:GLU:HG2	2:A:1153:HOH:O	2.16	0.46
1:A:417:ALA:HB2	1:A:609:GLN:HG3	1.97	0.46
1:A:530:LEU:HD11	1:A:651:ILE:HD13	1.97	0.46
1:A:628:ILE:HG12	1:A:630:TYR:N	2.30	0.46
1:A:664:GLY:O	1:A:666:GLU:N	2.48	0.46
1:B:418:VAL:O	1:B:421:LYS:N	2.46	0.46
1:A:731:LEU:HD21	1:B:601:ILE:HD13	1.98	0.45
1:B:467:LYS:HB2	2:B:1169:HOH:O	2.17	0.45
1:A:644:THR:O	1:A:645:VAL:C	2.54	0.45
1:A:736:LYS:HG2	1:B:732:MET:HB3	1.98	0.45
1:A:500:LEU:O	1:A:501:ASN:C	2.55	0.45
1:B:557:ARG:CG	1:B:558:PRO:HD2	2.46	0.45
1:A:735:PRO:CA	1:B:731:LEU:HD12	2.47	0.45
1:B:516:LEU:HA	1:B:560:SER:O	2.16	0.45
1:B:644:THR:O	1:B:645:VAL:C	2.54	0.45
1:B:507:LEU:HB3	1:B:509:ILE:HG13	1.98	0.45
1:A:473:VAL:HG23	1:A:473:VAL:O	2.17	0.45
1:B:527:PRO:HB3	1:B:738:VAL:HG21	1.98	0.45
1:A:628:ILE:HD13	1:A:628:ILE:N	2.32	0.45
1:A:732:MET:HE1	1:B:736:LYS:HA	1.99	0.45
1:A:614:GLY:CA	1:A:628:ILE:HG22	2.42	0.44
1:B:597:ALA:HB1	1:B:656:ILE:HG13	2.00	0.44
1:A:476:GLU:HG3	1:A:611:SER:HB2	1.97	0.44
1:B:480:ILE:HG21	1:B:603:ARG:NH1	2.32	0.44
1:B:530:LEU:HD11	1:B:651:ILE:HD13	1.99	0.44
1:B:552:VAL:O	1:B:552:VAL:HG12	2.17	0.44
1:B:742:ILE:O	1:B:746:ALA:HB2	2.18	0.44
1:A:466:GLY:O	1:A:468:LYS:N	2.50	0.44
1:B:656:ILE:O	1:B:658:PHE:N	2.47	0.44
1:B:480:ILE:CG2	1:B:603:ARG:HH12	2.31	0.44
1:A:480:ILE:CG2	1:A:603:ARG:HH12	2.30	0.44
1:A:507:LEU:HB3	1:A:509:ILE:HG13	1.99	0.44
1:B:466:GLY:O	1:B:468:LYS:N	2.50	0.44
1:B:599:ARG:HG2	1:B:599:ARG:NH1	2.33	0.44
1:B:418:VAL:C	1:B:420:ALA:H	2.21	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:490:GLU:O	1:B:493:GLU:HB3	2.18	0.43
1:B:591:ILE:HD11	1:B:753:LEU:HD21	2.00	0.43
1:A:485:THR:OG1	1:A:486:GLU:N	2.51	0.43
1:A:563:VAL:HG13	1:A:566:VAL:HG21	2.00	0.43
1:B:620:GLY:HA2	2:B:1146:HOH:O	2.18	0.43
1:B:500:LEU:O	1:B:501:ASN:C	2.55	0.43
1:B:563:VAL:HA	1:B:564:PRO:HD2	1.84	0.43
1:B:605:ASN:HD21	1:B:735:PRO:HG2	1.82	0.43
1:B:755:ALA:O	1:B:758:HIS:HA	2.19	0.43
1:A:416:LEU:HD12	2:A:1106:HOH:O	2.19	0.43
1:A:520:LEU:HD11	1:A:615:ILE:HG13	2.00	0.43
1:B:417:ALA:HB2	1:B:609:GLN:HG3	2.00	0.43
1:B:563:VAL:HG13	1:B:566:VAL:HG22	2.00	0.43
1:A:490:GLU:O	1:A:493:GLU:HB3	2.18	0.43
1:A:599:ARG:HG2	1:A:599:ARG:NH1	2.34	0.43
1:A:620:GLY:HA2	2:A:146:HOH:O	2.19	0.43
1:A:418:VAL:C	1:A:420:ALA:H	2.21	0.42
1:A:480:ILE:HG21	1:A:603:ARG:NH1	2.33	0.42
1:B:563:VAL:HG13	1:B:566:VAL:HG21	2.01	0.42
1:A:563:VAL:HG13	1:A:566:VAL:HG22	2.00	0.42
1:A:593:ILE:HA	1:A:649:VAL:O	2.20	0.42
1:B:593:ILE:HA	1:B:649:VAL:O	2.20	0.42
1:A:557:ARG:CG	1:A:558:PRO:HD2	2.49	0.42
1:A:467:LYS:HB2	2:A:169:HOH:O	2.18	0.42
1:A:526:THR:HB	1:A:527:PRO:HD3	2.01	0.42
1:A:597:ALA:HB1	1:A:656:ILE:HG13	2.02	0.42
1:B:550:GLU:CB	2:B:1100:HOH:O	2.68	0.42
1:B:755:ALA:O	1:B:758:HIS:N	2.53	0.42
1:B:473:VAL:HG23	1:B:473:VAL:O	2.19	0.42
1:B:485:THR:OG1	1:B:486:GLU:N	2.51	0.42
1:B:416:LEU:O	1:B:417:ALA:C	2.58	0.42
1:A:480:ILE:HB	1:A:603:ARG:NH1	2.34	0.41
1:A:742:ILE:O	1:A:746:ALA:HB2	2.20	0.41
1:A:755:ALA:O	1:A:758:HIS:HA	2.20	0.41
1:B:478:VAL:HG22	1:B:609:GLN:CG	2.49	0.41
1:A:651:ILE:HG21	1:A:745:MET:CE	2.51	0.41
1:A:724:LEU:O	1:A:725:ALA:O	2.39	0.41
1:B:707:MET:O	1:B:709:GLY:N	2.53	0.41
1:A:550:GLU:CB	2:A:100:HOH:O	2.68	0.41
1:A:591:ILE:HD11	1:A:753:LEU:HD21	2.01	0.41
1:B:507:LEU:HD13	1:B:509:ILE:CD1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:470:GLY:HA3	2:A:173:HOH:O	2.20	0.41
1:A:478:VAL:HG22	1:A:609:GLN:CG	2.50	0.41
1:B:628:ILE:N	1:B:628:ILE:HD13	2.33	0.41
1:B:660:LEU:HD23	1:B:660:LEU:HA	1.72	0.41
1:A:519:GLY:O	1:A:563:VAL:HG23	2.20	0.41
1:A:651:ILE:HG21	1:A:745:MET:HE2	2.03	0.41
1:A:577:ILE:C	1:A:579:PHE:N	2.74	0.41
1:B:466:GLY:C	1:B:468:LYS:H	2.25	0.41
1:B:480:ILE:HB	1:B:603:ARG:NH1	2.36	0.41
1:A:416:LEU:O	1:A:417:ALA:C	2.59	0.41
1:A:516:LEU:CD2	1:A:582:VAL:HG12	2.51	0.40
1:A:530:LEU:HD11	1:A:593:ILE:HD11	1.90	0.40
1:B:753:LEU:HD22	1:B:753:LEU:HA	1.89	0.40
1:B:575:SER:CA	1:B:578:ILE:HG12	2.46	0.40
1:B:651:ILE:HG21	1:B:745:MET:CE	2.51	0.40
1:A:707:MET:O	1:A:709:GLY:N	2.54	0.40
1:A:418:VAL:O	1:A:420:ALA:N	2.55	0.40
1:B:526:THR:HG21	1:B:738:VAL:HG13	2.03	0.40

All (19) 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)
2:A:177:HOH:O	2:B:1174:HOH:O[1_655]	0.20	2.00
2:A:174:HOH:O	2:B:1177:HOH:O[1_655]	0.36	1.84
2:A:112:HOH:O	2:B:1107:HOH:O[1_655]	0.44	1.76
2:A:144:HOH:O	2:B:1105:HOH:O[8_555]	0.48	1.72
2:A:105:HOH:O	2:B:1144:HOH:O[8_555]	0.61	1.59
2:A:132:HOH:O	2:B:1110:HOH:O[1_655]	0.62	1.58
2:A:110:HOH:O	2:B:1132:HOH:O[1_655]	0.67	1.53
2:A:107:HOH:O	2:B:1112:HOH:O[1_655]	0.70	1.50
1:A:769:TYR:CB	1:B:769:TYR:CB[1_655]	1.14	1.06
2:A:108:HOH:O	2:B:1113:HOH:O[8_555]	1.31	0.89
2:A:113:HOH:O	2:B:1108:HOH:O[8_555]	1.54	0.66
1:A:771:HIS:CB	1:B:475:LEU:CB[1_655]	1.77	0.43
1:A:475:LEU:CB	1:B:771:HIS:CB[1_655]	1.84	0.36
1:B:755:ALA:O	2:A:149:HOH:O[1_455]	1.95	0.25
1:B:755:ALA:C	2:A:149:HOH:O[1_455]	1.99	0.21
1:B:713:ASP:OD1	2:A:193:HOH:O[5_455]	1.99	0.21
2:A:188:HOH:O	2:B:1158:HOH:O[1_655]	2.12	0.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:771:HIS:CB	1:B:475:LEU:CG[1_655]	2.14	0.06
1:A:475:LEU:CG	1:B:771:HIS:CB[1_655]	2.16	0.04

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	312/371 (84%)	238 (76%)	51 (16%)	23 (7%)	1 5
1	B	312/371 (84%)	236 (76%)	53 (17%)	23 (7%)	1 5
All	All	624/742 (84%)	474 (76%)	104 (17%)	46 (7%)	1 5

All (46) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	406	ASP
1	A	417	ALA
1	A	429	PRO
1	A	469	LYS
1	A	471	ARG
1	A	485	THR
1	A	549	PRO
1	A	619	SER
1	A	708	ILE
1	A	725	ALA
1	A	762	ASP
1	B	406	ASP
1	B	417	ALA
1	B	429	PRO
1	B	469	LYS
1	B	471	ARG
1	B	485	THR
1	B	549	PRO

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Mol	Chain	Res	Type
1	B	619	SER
1	B	708	ILE
1	B	725	ALA
1	B	762	ASP
1	A	712	PRO
1	B	712	PRO
1	A	467	LYS
1	A	486	GLU
1	A	617	PRO
1	B	467	LYS
1	B	486	GLU
1	A	474	THR
1	A	622	GLY
1	A	765	ASN
1	A	766	PHE
1	B	474	THR
1	B	617	PRO
1	B	622	GLY
1	B	765	ASN
1	B	766	PHE
1	A	412	VAL
1	B	412	VAL
1	A	665	ARG
1	B	665	ARG
1	A	552	VAL
1	B	552	VAL
1	A	621	VAL
1	B	621	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	192/317 (61%)	172 (90%)	20 (10%)	7	27
1	B	192/317 (61%)	173 (90%)	19 (10%)	8	30
All	All	384/634 (61%)	345 (90%)	39 (10%)	7	28

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	405	LEU
1	A	447	GLU
1	A	459	GLU
1	A	472	TYR
1	A	476	GLU
1	A	480	ILE
1	A	500	LEU
1	A	507	LEU
1	A	591	ILE
1	A	593	ILE
1	A	607	THR
1	A	609	GLN
1	A	628	ILE
1	A	645	VAL
1	A	647	ASP
1	A	649	VAL
1	A	714	GLU
1	A	727	LEU
1	A	753	LEU
1	A	757	LEU
1	B	405	LEU
1	B	447	GLU
1	B	459	GLU
1	B	472	TYR
1	B	476	GLU
1	B	480	ILE
1	B	500	LEU
1	B	507	LEU
1	B	591	ILE
1	B	593	ILE
1	B	607	THR
1	B	609	GLN
1	B	628	ILE
1	B	645	VAL
1	B	647	ASP
1	B	649	VAL
1	B	714	GLU
1	B	727	LEU
1	B	753	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	448	GLN
1	A	522	ASN
1	A	543	HIS
1	A	605	ASN
1	A	747	ASN
1	A	765	ASN
1	B	522	ASN
1	B	543	HIS
1	B	605	ASN
1	B	747	ASN
1	B	765	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 carbohydrates 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

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	320/371 (86%)	0.37	25 (7%) 13 4	19, 59, 98, 103	0
1	B	320/371 (86%)	0.35	24 (7%) 14 4	19, 59, 98, 103	0
All	All	640/742 (86%)	0.36	49 (7%) 13 4	19, 59, 98, 103	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	551	SER	10.5
1	A	404	GLU	6.6
1	A	405	LEU	6.5
1	A	414	THR	6.4
1	B	709	GLY	6.3
1	A	770	THR	6.1
1	A	453	SER	5.9
1	B	403	LYS	5.3
1	A	762	ASP	5.3
1	A	632	THR	5.1
1	B	771	HIS	4.8
1	A	633	LEU	4.6
1	A	403	LYS	4.6
1	B	770	THR	4.4
1	A	771	HIS	4.4
1	A	552	VAL	4.3
1	B	424	ALA	4.3
1	B	404	GLU	4.3
1	B	428	GLN	4.1
1	A	406	ASP	3.9
1	A	763	GLN	3.8
1	B	706	GLY	3.8
1	B	671	GLY	3.8
1	B	402	GLU	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	668	LYS	3.6
1	B	429	PRO	3.4
1	B	717	ARG	3.4
1	B	705	LEU	3.4
1	B	551	SER	3.1
1	A	474	THR	3.1
1	A	549	PRO	2.9
1	B	401	MET	2.8
1	B	486	GLU	2.8
1	A	709	GLY	2.7
1	B	762	ASP	2.7
1	B	474	THR	2.7
1	B	414	THR	2.7
1	A	766	PHE	2.7
1	A	425	LEU	2.6
1	A	470	GLY	2.6
1	B	667	MET	2.6
1	A	705	LEU	2.4
1	A	710	THR	2.4
1	B	632	THR	2.3
1	B	407	LEU	2.2
1	B	471	ARG	2.2
1	A	760	GLU	2.2
1	A	408	SER	2.2
1	A	512	ASP	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 carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.