



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 30, 2023 – 06:56 AM EDT

PDB ID : 3NO9
Title : Crystal Structure of apo fumarate hydratase from Mycobacterium tuberculosis
Authors : Li, H.; Swanson, S.; Yu, M.; Hung, L.-W.; Sacchettini, J.C.; TB Structural Genomics Consortium (TBSGC)
Deposited on : 2010-06-25
Resolution : 2.48 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 : 4.02b-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

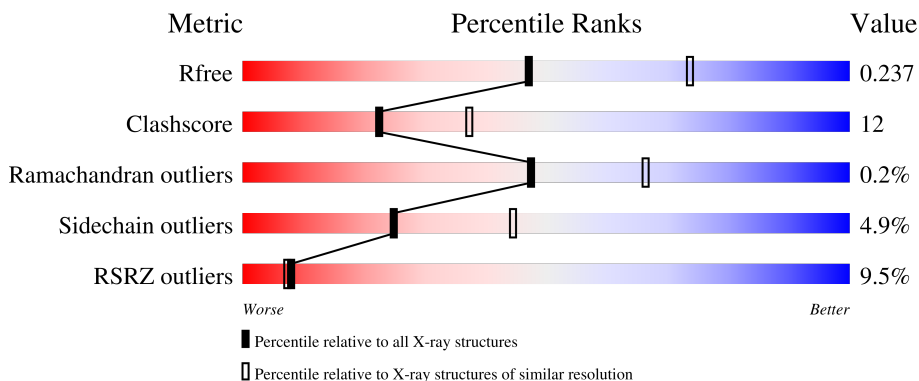
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.48 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	5857 (2.50-2.46)
Clashscore	141614	6594 (2.50-2.46)
Ramachandran outliers	138981	6469 (2.50-2.46)
Sidechain outliers	138945	6471 (2.50-2.46)
RSRZ outliers	127900	5738 (2.50-2.46)

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	475	 13% 78% 15% • 5%
1	B	475	 5% 79% 14% • 5%
1	C	475	 13% 75% 17% • 5%
1	D	475	 5% 78% 14% • 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13623 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 Fumarate hydratase class II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	453	3363	2095	609	648	11	38	1	0
1	B	452	3357	2092	608	646	11	16	1	0
1	C	453	3368	2098	611	648	11	13	1	0
1	D	452	3357	2092	608	646	11	13	1	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP O53446
B	0	SER	-	expression tag	UNP O53446
C	0	SER	-	expression tag	UNP O53446
D	0	SER	-	expression tag	UNP O53446

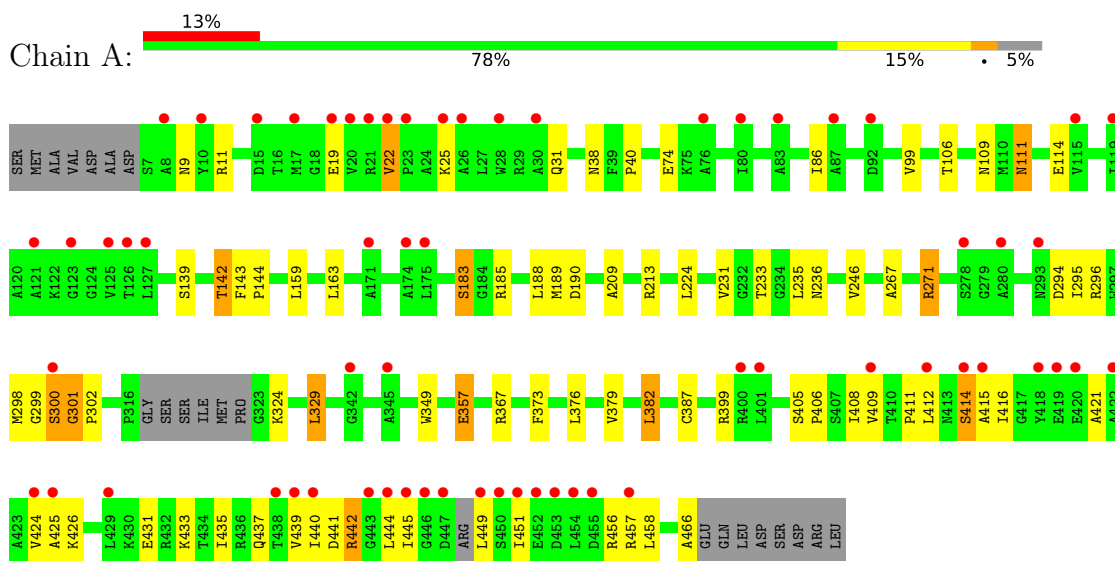
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	37	Total	O	0	0
			37	37		
2	B	55	Total	O	0	0
			55	55		
2	C	38	Total	O	0	0
			38	38		
2	D	48	Total	O	0	0
			48	48		

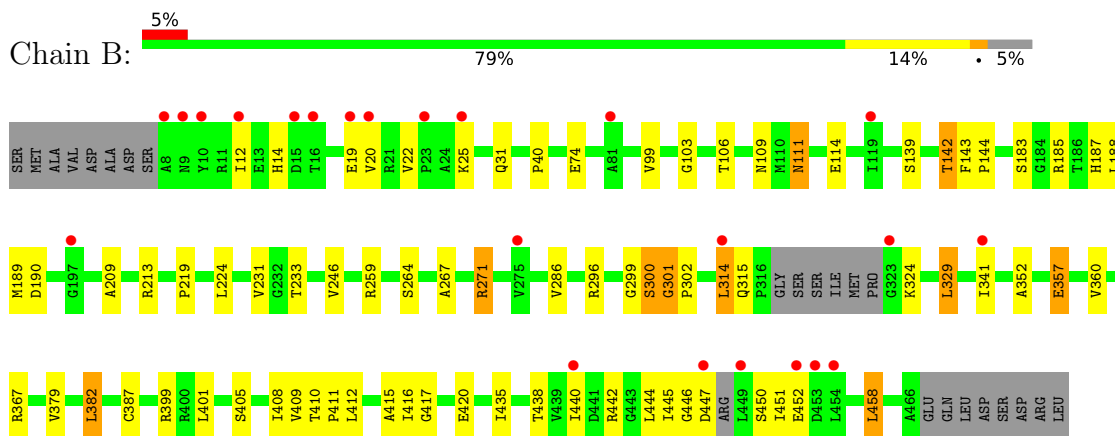
3 Residue-property plots i

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.

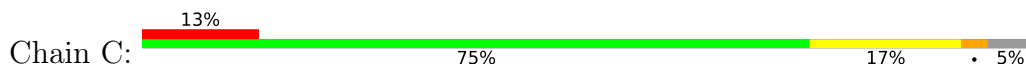
- Molecule 1: Fumarate hydratase class II

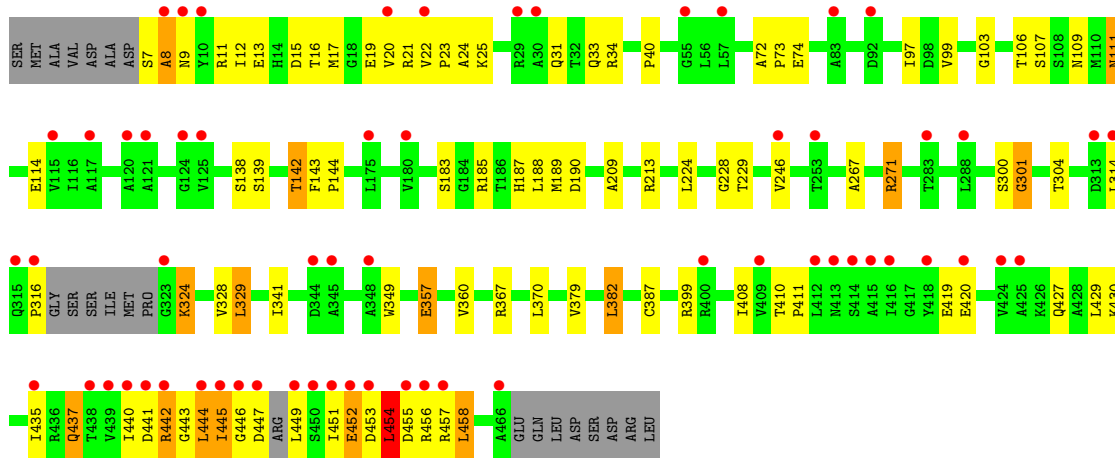


- Molecule 1: Fumarate hydratase class II

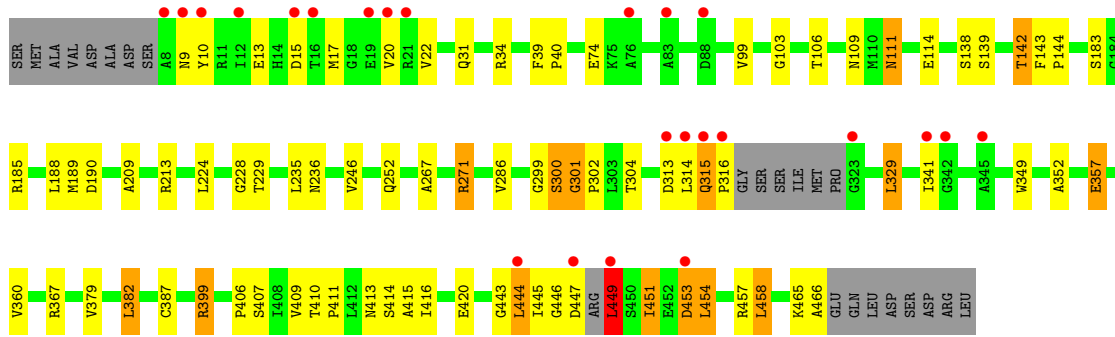
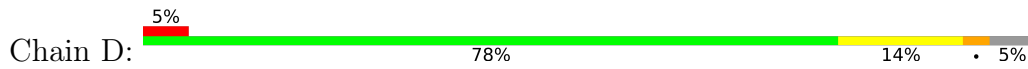


- Molecule 1: Fumarate hydratase class II





- Molecule 1: Fumarate hydratase class II



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	271.22Å 96.56Å 89.89Å 90.00° 102.99° 90.00°	Depositor
Resolution (Å)	48.52 – 2.48 48.52 – 2.48	Depositor EDS
% Data completeness (in resolution range)	88.4 (48.52-2.48) 88.4 (48.52-2.48)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.87 (at 2.48Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.6.1_357)	Depositor
R, R_{free}	0.206 , 0.241 0.207 , 0.237	Depositor DCC
R_{free} test set	3534 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	41.0	Xtrriage
Anisotropy	0.341	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13623	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.10% 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	0.41	0/3411	0.54	1/4635 (0.0%)
1	B	0.42	0/3405	0.56	3/4627 (0.1%)
1	C	0.42	1/3417 (0.0%)	0.58	3/4643 (0.1%)
1	D	0.43	0/3405	0.56	2/4627 (0.0%)
All	All	0.42	1/13638 (0.0%)	0.56	9/18532 (0.0%)

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	3
1	B	0	4
1	C	0	3
1	D	0	3
All	All	0	13

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	441	ASP	C-N	-5.83	1.20	1.34

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	12	ILE	CB-CA-C	5.98	123.56	111.60
1	C	452	GLU	CB-CA-C	5.97	122.34	110.40
1	A	456	ARG	N-CA-CB	-5.96	99.88	110.60
1	B	314	LEU	CA-CB-CG	-5.84	101.86	115.30
1	D	449	LEU	CA-CB-CG	-5.78	102.02	115.30
1	C	441	ASP	O-C-N	-5.58	113.77	122.70
1	C	441	ASP	CB-CA-C	5.54	121.48	110.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	12	ILE	N-CA-C	-5.36	96.54	111.00
1	D	449	LEU	CB-CG-CD2	-5.01	102.47	111.00

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	299	GLY	Peptide
1	A	300	SER	Peptide
1	A	301	GLY	Peptide
1	B	299	GLY	Peptide
1	B	300	SER	Peptide
1	B	301	GLY	Peptide
1	B	415	ALA	Peptide
1	C	300	SER	Peptide
1	C	301	GLY	Peptide
1	C	8	ALA	Peptide
1	D	299	GLY	Peptide
1	D	300	SER	Peptide
1	D	301	GLY	Peptide

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	3363	0	3399	100	0
1	B	3357	0	3394	66	0
1	C	3368	0	3403	109	0
1	D	3357	0	3394	91	0
2	A	37	0	0	3	0
2	B	55	0	0	4	0
2	C	38	0	0	2	0
2	D	48	0	0	5	0
All	All	13623	0	13590	313	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 (313) 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:C:22:VAL:HG12	1:C:33:GLN:OE1	1.24	1.26
1:C:437:GLN:NE2	1:C:440:ILE:HD12	1.59	1.16
1:C:324:LYS:HZ2	1:C:324:LYS:HB3	1.19	1.05
1:A:415:ALA:HB1	1:A:449:LEU:HD13	1.39	1.04
1:C:22:VAL:CG1	1:C:33:GLN:OE1	2.08	1.02
1:C:324:LYS:HB3	1:C:324:LYS:NZ	1.74	0.99
1:D:449:LEU:N	1:D:449:LEU:HD23	1.72	0.99
1:D:449:LEU:N	1:D:449:LEU:CD2	2.29	0.95
1:C:13:GLU:HB3	1:C:20:VAL:CG2	1.97	0.94
1:C:437:GLN:HE22	1:C:440:ILE:HD12	1.27	0.94
1:C:22:VAL:HB	1:C:23:PRO:HD2	1.49	0.92
1:A:295:ILE:HA	1:A:298:MET:CE	2.01	0.91
1:C:17:MET:SD	1:D:315:GLN:O	2.28	0.91
1:A:295:ILE:HA	1:A:298:MET:HE3	1.53	0.90
1:C:22:VAL:HG12	1:C:33:GLN:CD	1.93	0.89
1:D:449:LEU:HD23	1:D:449:LEU:O	1.72	0.88
1:B:189:MET:CE	1:C:301:GLY:HA2	2.04	0.87
1:A:99:VAL:HG11	1:A:367:ARG:HD2	1.58	0.85
1:A:294:ASP:O	1:A:298:MET:HG3	1.77	0.85
1:D:314:LEU:C	1:D:316:PRO:CD	2.45	0.85
1:C:21:ARG:O	1:C:21:ARG:HG2	1.78	0.84
1:B:99:VAL:HG11	1:B:367:ARG:HD2	1.59	0.82
1:D:99:VAL:HG11	1:D:367:ARG:HD2	1.61	0.82
1:C:446:GLY:N	1:C:449:LEU:O	2.10	0.82
1:D:315:GLN:N	1:D:316:PRO:CD	2.43	0.81
1:D:449:LEU:CD2	1:D:449:LEU:O	2.30	0.80
1:D:314:LEU:C	1:D:316:PRO:HD2	2.00	0.80
1:D:445:ILE:HG21	1:D:451:ILE:HG13	1.64	0.80
1:A:439:VAL:O	1:A:444:LEU:HD13	1.82	0.80
1:C:437:GLN:HE21	1:C:437:GLN:HA	1.48	0.79
1:D:314:LEU:O	1:D:316:PRO:HD2	1.83	0.79
1:B:14:HIS:HA	1:B:19:GLU:HG2	1.65	0.78
1:A:415:ALA:CB	1:A:449:LEU:HD13	2.13	0.78
1:C:445:ILE:H	1:C:445:ILE:HD12	1.49	0.77
1:B:302:PRO:HD2	1:C:190:ASP:OD2	1.85	0.77
1:C:99:VAL:HG11	1:C:367:ARG:HD2	1.67	0.76
1:A:324:LYS:NZ	1:A:324:LYS:HB3	2.02	0.75
1:B:445:ILE:HD13	1:B:451:ILE:HG12	1.69	0.75
1:D:209:ALA:O	1:D:213:ARG:HG2	1.85	0.75
1:C:34:ARG:NE	2:C:497:HOH:O	2.20	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:445:ILE:HD11	1:D:454:LEU:HD12	1.70	0.73
1:A:190:ASP:OD2	1:D:302:PRO:HD2	1.87	0.73
1:A:324:LYS:HB3	1:A:324:LYS:HZ3	1.54	0.72
1:D:106:THR:HA	1:D:139:SER:OG	1.90	0.72
1:B:189:MET:HE1	1:C:301:GLY:HA2	1.71	0.72
1:B:106:THR:HA	1:B:139:SER:OG	1.90	0.72
1:A:324:LYS:NZ	1:A:324:LYS:CB	2.53	0.71
1:A:439:VAL:HG13	1:A:444:LEU:HD22	1.72	0.71
1:A:106:THR:HA	1:A:139:SER:OG	1.89	0.71
1:D:314:LEU:C	1:D:316:PRO:HD3	2.11	0.70
1:A:159:LEU:HD22	1:A:373:PHE:CD1	2.27	0.70
1:C:106:THR:HA	1:C:139:SER:OG	1.91	0.69
1:A:300:SER:HB2	1:D:188:LEU:HB3	1.74	0.69
1:C:209:ALA:O	1:C:213:ARG:HG2	1.91	0.69
1:C:442:ARG:HB3	1:C:444:LEU:HD23	1.73	0.69
1:A:442:ARG:HB2	1:A:444:LEU:HD11	1.75	0.67
1:A:301:GLY:CA	1:D:189:MET:CE	2.72	0.67
1:A:444:LEU:N	1:A:444:LEU:HD12	2.09	0.67
1:A:209:ALA:O	1:A:213:ARG:HG2	1.95	0.66
1:C:454:LEU:O	1:C:456:ARG:N	2.29	0.66
1:B:189:MET:CE	1:C:301:GLY:CA	2.74	0.66
1:B:189:MET:HA	1:C:301:GLY:HA3	1.76	0.65
1:C:442:ARG:HB3	1:C:444:LEU:CD2	2.25	0.65
1:C:7:SER:OG	1:C:8:ALA:N	2.30	0.65
1:C:453:ASP:O	1:C:456:ARG:HB3	1.97	0.64
1:D:313:ASP:O	1:D:316:PRO:HD3	1.97	0.64
1:A:295:ILE:HA	1:A:298:MET:HE2	1.78	0.64
1:D:454:LEU:HD22	1:D:458:LEU:HD22	1.79	0.64
1:C:329:LEU:HD23	1:C:387:CYS:HB2	1.80	0.63
1:B:209:ALA:O	1:B:213:ARG:HG2	1.98	0.63
1:A:382:LEU:HD21	1:B:40:PRO:HD2	1.80	0.63
1:A:301:GLY:HA2	1:D:189:MET:CE	2.29	0.63
1:B:267:ALA:O	1:B:271:ARG:NH2	2.32	0.62
1:A:445:ILE:HG23	1:A:449:LEU:O	2.00	0.62
1:C:9:ASN:OD1	1:C:9:ASN:C	2.34	0.62
1:C:437:GLN:HE21	1:C:437:GLN:CA	2.13	0.62
1:C:440:ILE:HA	1:C:445:ILE:HD11	1.82	0.62
1:D:267:ALA:O	1:D:271:ARG:NH2	2.31	0.62
1:D:449:LEU:O	1:D:449:LEU:CG	2.48	0.62
1:D:9:ASN:OD1	1:D:10:TYR:CE2	2.53	0.61
1:C:437:GLN:NE2	1:C:437:GLN:HA	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:185:ARG:O	1:D:357:GLU:HG2	2.01	0.61
1:A:267:ALA:O	1:A:271:ARG:NH2	2.34	0.60
1:C:437:GLN:HE22	1:C:440:ILE:CD1	2.08	0.60
1:A:442:ARG:C	1:A:444:LEU:HD12	2.21	0.60
1:C:13:GLU:HB3	1:C:20:VAL:HG23	1.83	0.60
1:A:442:ARG:HG3	1:A:444:LEU:HD11	1.83	0.59
1:B:187:HIS:O	1:B:188:LEU:HB2	2.02	0.59
1:A:301:GLY:N	1:D:189:MET:SD	2.75	0.59
1:B:189:MET:SD	1:C:301:GLY:CA	2.91	0.59
1:A:440:ILE:HG12	1:A:445:ILE:HD12	1.84	0.59
1:B:219:PRO:HA	2:B:527:HOH:O	2.02	0.59
1:A:442:ARG:CB	1:A:444:LEU:HD11	2.31	0.59
1:D:444:LEU:O	1:D:449:LEU:HD21	2.03	0.59
1:B:357:GLU:HG2	1:D:185:ARG:O	2.03	0.59
1:D:446:GLY:O	1:D:447:ASP:HB2	2.03	0.58
1:C:453:ASP:HA	1:C:456:ARG:HE	1.67	0.58
1:A:426:LYS:O	1:A:426:LYS:HD3	2.04	0.58
1:D:329:LEU:HD23	1:D:387:CYS:HB2	1.86	0.58
1:D:454:LEU:CD2	1:D:458:LEU:HD22	2.34	0.58
1:A:302:PRO:HD2	1:D:190:ASP:OD2	2.04	0.57
1:A:444:LEU:H	1:A:444:LEU:CD1	2.17	0.57
1:C:324:LYS:NZ	1:C:324:LYS:CB	2.49	0.57
1:C:382:LEU:HD21	1:D:40:PRO:HD2	1.86	0.57
1:C:454:LEU:C	1:C:456:ARG:H	2.08	0.57
1:B:213:ARG:NH1	2:B:478:HOH:O	2.37	0.57
1:C:17:MET:HG3	1:C:34:ARG:NH1	2.19	0.57
1:D:20:VAL:CG1	1:D:34:ARG:HE	2.16	0.57
1:A:301:GLY:HA2	1:D:189:MET:HE1	1.85	0.57
1:B:329:LEU:HD23	1:B:387:CYS:HB2	1.86	0.57
1:A:444:LEU:HD12	1:A:444:LEU:H	1.69	0.56
1:A:142:THR:HG22	1:A:143:PHE:N	2.19	0.56
1:C:328:VAL:HG11	1:D:39:PHE:HZ	1.71	0.56
1:C:349:TRP:CZ3	1:D:341:ILE:HD13	2.40	0.56
1:D:315:GLN:N	1:D:316:PRO:HD3	2.20	0.56
1:A:189:MET:HE1	1:D:301:GLY:HA2	1.87	0.56
1:A:301:GLY:HA3	1:D:189:MET:HE3	1.86	0.56
1:B:408:ILE:O	1:B:411:PRO:HD2	2.05	0.56
1:A:324:LYS:CB	1:A:324:LYS:HZ2	2.16	0.56
1:B:31:GLN:HB2	1:B:114:GLU:OE2	2.06	0.56
1:C:454:LEU:C	1:C:456:ARG:N	2.59	0.56
1:A:295:ILE:CG1	1:A:298:MET:HE2	2.36	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:21:ARG:O	1:C:21:ARG:CG	2.51	0.56
1:D:9:ASN:OD1	1:D:10:TYR:CD2	2.59	0.56
1:C:267:ALA:O	1:C:271:ARG:NH2	2.39	0.55
1:D:224:LEU:HD12	1:D:246:VAL:HG22	1.87	0.55
1:C:411:PRO:HB2	1:C:457:ARG:HB3	1.87	0.55
1:D:31:GLN:HB2	1:D:114:GLU:OE2	2.07	0.54
1:A:295:ILE:HG12	1:A:298:MET:CE	2.37	0.54
1:C:12:ILE:HA	1:C:21:ARG:HA	1.89	0.54
1:B:445:ILE:HG21	1:B:451:ILE:HG12	1.88	0.54
1:A:442:ARG:CG	1:A:444:LEU:HD11	2.38	0.54
1:D:406:PRO:O	1:D:409:VAL:HG22	2.08	0.54
1:A:11:ARG:HH21	1:A:22:VAL:HG21	1.73	0.54
1:A:31:GLN:HB2	1:A:114:GLU:OE2	2.07	0.54
1:A:329:LEU:HD23	1:A:387:CYS:HB2	1.88	0.54
1:B:324:LYS:O	1:B:324:LYS:HG2	2.06	0.54
1:A:349:TRP:CZ3	1:B:341:ILE:HD13	2.43	0.54
1:D:315:GLN:NE2	2:D:513:HOH:O	2.40	0.54
1:B:189:MET:SD	1:C:301:GLY:N	2.81	0.54
1:D:449:LEU:N	1:D:449:LEU:HD22	2.22	0.54
1:B:25:LYS:HD2	1:B:25:LYS:O	2.08	0.53
1:A:189:MET:CE	1:D:301:GLY:HA2	2.38	0.53
1:B:264:SER:HB2	2:B:492:HOH:O	2.08	0.53
1:A:411:PRO:HB2	1:A:457:ARG:HB3	1.91	0.53
1:A:159:LEU:HD22	1:A:373:PHE:HD1	1.69	0.53
1:A:440:ILE:HG13	1:A:445:ILE:HD11	1.91	0.53
1:C:15:ASP:OD2	1:C:16:THR:N	2.39	0.53
1:C:19:GLU:O	1:C:19:GLU:HG3	2.09	0.53
1:A:349:TRP:CH2	1:B:341:ILE:HD13	2.44	0.52
1:B:224:LEU:HD12	1:B:246:VAL:HG22	1.92	0.52
1:C:31:GLN:HB2	1:C:114:GLU:OE2	2.10	0.52
1:C:408:ILE:O	1:C:411:PRO:HD2	2.09	0.52
1:D:415:ALA:HB2	1:D:457:ARG:CZ	2.40	0.52
1:B:143:PHE:N	1:B:144:PRO:HD2	2.24	0.52
1:D:252:GLN:NE2	2:D:515:HOH:O	2.23	0.52
1:C:11:ARG:N	1:C:22:VAL:O	2.42	0.51
1:C:143:PHE:N	1:C:144:PRO:HD2	2.25	0.51
1:C:22:VAL:HB	1:C:23:PRO:CD	2.31	0.51
1:A:231:VAL:HG23	1:A:233:THR:HG23	1.91	0.51
1:A:406:PRO:O	1:A:409:VAL:HG22	2.11	0.51
1:A:444:LEU:N	1:A:444:LEU:CD1	2.72	0.51
1:C:349:TRP:CH2	1:D:341:ILE:HD13	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:416:ILE:HB	1:B:420:GLU:HB2	1.93	0.51
1:D:111:ASN:C	1:D:111:ASN:HD22	2.13	0.51
1:A:301:GLY:HA3	1:D:189:MET:CE	2.39	0.51
1:A:301:GLY:CA	1:D:189:MET:SD	2.99	0.51
1:A:224:LEU:HD12	1:A:246:VAL:HG22	1.91	0.51
1:B:111:ASN:C	1:B:111:ASN:HD22	2.14	0.51
1:B:259:ARG:NH2	2:B:527:HOH:O	2.42	0.51
1:C:13:GLU:HB3	1:C:20:VAL:HG22	1.88	0.50
1:C:109:ASN:OD1	1:C:142:THR:HG21	2.11	0.50
1:B:442:ARG:HB2	1:B:444:LEU:CD2	2.41	0.50
1:C:111:ASN:C	1:C:111:ASN:HD22	2.12	0.50
1:D:413:ASN:O	1:D:416:ILE:O	2.29	0.50
1:A:405:SER:O	1:A:408:ILE:HG12	2.11	0.50
1:D:445:ILE:HD13	1:D:451:ILE:HG13	1.93	0.50
1:A:189:MET:CE	1:D:301:GLY:CA	2.90	0.50
1:A:367:ARG:HG2	1:A:367:ARG:HH11	1.77	0.49
1:C:34:ARG:CZ	2:C:497:HOH:O	2.58	0.49
1:C:446:GLY:O	1:C:447:ASP:HB2	2.12	0.49
1:B:187:HIS:O	1:B:188:LEU:CB	2.60	0.49
1:D:142:THR:HG22	1:D:143:PHE:N	2.26	0.49
1:A:440:ILE:HG12	1:A:445:ILE:CD1	2.42	0.49
1:A:301:GLY:CA	1:D:189:MET:HE3	2.41	0.49
1:B:438:THR:O	1:B:442:ARG:HG3	2.13	0.49
1:D:143:PHE:N	1:D:144:PRO:HD2	2.27	0.49
1:B:109:ASN:OD1	1:B:142:THR:HG21	2.13	0.49
1:A:445:ILE:HD13	1:A:451:ILE:HG12	1.95	0.48
1:A:40:PRO:HD2	1:B:382:LEU:HD21	1.93	0.48
1:B:301:GLY:HA2	1:C:189:MET:CE	2.42	0.48
1:D:449:LEU:O	1:D:449:LEU:HG	2.12	0.48
1:C:138:SER:HB3	1:C:229:THR:HA	1.94	0.48
1:A:445:ILE:HG21	1:A:451:ILE:HG13	1.95	0.48
1:C:410:THR:N	1:C:411:PRO:CD	2.77	0.48
1:A:295:ILE:HG13	1:A:298:MET:HE2	1.95	0.47
1:A:185:ARG:O	1:C:357:GLU:HG2	2.14	0.47
1:B:352:ALA:HA	1:D:286:VAL:HG13	1.96	0.47
1:D:20:VAL:HG13	1:D:34:ARG:HE	1.79	0.47
1:D:329:LEU:HD12	1:D:329:LEU:HA	1.80	0.47
1:A:143:PHE:N	1:A:144:PRO:HD2	2.30	0.47
1:C:224:LEU:HD12	1:C:246:VAL:HG22	1.95	0.47
1:C:420:GLU:OE2	1:C:444:LEU:HD21	2.14	0.47
1:D:315:GLN:O	1:D:316:PRO:C	2.51	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:367:ARG:HG2	1:C:367:ARG:HH11	1.80	0.46
1:A:409:VAL:O	1:A:412:LEU:HB2	2.14	0.46
1:A:466:ALA:N	2:A:495:HOH:O	2.47	0.46
1:B:324:LYS:O	1:B:324:LYS:CG	2.63	0.46
1:C:20:VAL:HB	1:C:33:GLN:HG2	1.98	0.46
1:A:295:ILE:CG1	1:A:298:MET:CE	2.94	0.46
1:C:103:GLY:HA3	1:C:360:VAL:HB	1.97	0.46
1:D:109:ASN:OD1	1:D:142:THR:HG21	2.16	0.46
1:D:410:THR:OG1	1:D:411:PRO:HD3	2.16	0.46
1:A:433:LYS:HE2	1:A:441:ASP:OD2	2.16	0.46
1:B:296:ARG:HG3	1:C:188:LEU:HD12	1.96	0.46
1:D:235:LEU:O	1:D:236:ASN:HB2	2.15	0.46
1:B:409:VAL:O	1:B:412:LEU:HB2	2.16	0.46
1:C:451:ILE:H	1:C:451:ILE:HG13	1.51	0.46
1:B:440:ILE:HG12	1:B:445:ILE:HD12	1.97	0.45
1:C:187:HIS:O	1:C:188:LEU:HB2	2.15	0.45
1:C:454:LEU:O	1:C:457:ARG:N	2.49	0.45
1:D:445:ILE:CG2	1:D:451:ILE:HG13	2.42	0.45
1:A:235:LEU:O	1:A:236:ASN:HB2	2.15	0.45
1:A:440:ILE:CG1	1:A:445:ILE:HD11	2.47	0.45
1:B:142:THR:HG22	1:B:143:PHE:N	2.31	0.45
1:D:15:ASP:OD1	1:D:34:ARG:NH1	2.49	0.45
1:A:38:ASN:HD21	1:B:315:GLN:HG3	1.82	0.45
1:A:433:LYS:HE3	1:A:437:GLN:HG3	1.98	0.45
1:B:367:ARG:HG2	1:B:367:ARG:HH11	1.81	0.45
1:D:453:ASP:O	1:D:457:ARG:HG3	2.17	0.45
1:A:412:LEU:O	1:A:416:ILE:HG12	2.17	0.45
1:B:301:GLY:HA2	1:C:189:MET:HE1	1.99	0.45
1:D:414:SER:OG	1:D:457:ARG:NH2	2.50	0.44
1:C:22:VAL:CB	1:C:23:PRO:HD2	2.24	0.44
1:C:328:VAL:HG11	1:D:39:PHE:CZ	2.51	0.44
1:A:439:VAL:HG13	1:A:444:LEU:HB2	1.99	0.44
1:C:329:LEU:HD12	1:C:329:LEU:HA	1.79	0.44
1:C:454:LEU:HD23	1:C:454:LEU:HA	1.75	0.44
1:A:38:ASN:ND2	1:B:315:GLN:HE21	2.16	0.44
1:B:286:VAL:HG13	1:D:352:ALA:HA	2.00	0.44
1:C:40:PRO:HD2	1:D:382:LEU:HD21	2.00	0.44
1:A:111:ASN:HD22	1:A:111:ASN:C	2.21	0.44
1:A:25:LYS:HD2	1:A:25:LYS:O	2.18	0.44
1:B:231:VAL:HG23	1:B:233:THR:HG23	1.99	0.44
1:B:435:ILE:HG22	1:B:458:LEU:HD21	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:13:GLU:HB3	1:D:20:VAL:HG23	1.99	0.44
1:D:410:THR:N	1:D:411:PRO:CD	2.81	0.44
1:D:443:GLY:O	2:D:518:HOH:O	2.21	0.44
1:A:424:VAL:CG1	1:A:435:ILE:HG23	2.48	0.43
1:D:213:ARG:NH1	2:D:477:HOH:O	2.51	0.43
1:A:86:ILE:HA	2:A:476:HOH:O	2.19	0.43
1:A:424:VAL:HG11	1:A:435:ILE:HG23	1.99	0.43
1:B:189:MET:HE3	1:C:301:GLY:HA2	1.90	0.43
1:A:109:ASN:OD1	1:A:142:THR:HG21	2.18	0.43
1:A:431:GLU:OE1	1:A:433:LYS:HD2	2.18	0.43
1:B:329:LEU:HD12	1:B:329:LEU:HA	1.80	0.43
1:B:446:GLY:O	1:B:447:ASP:HB2	2.19	0.43
1:D:103:GLY:HA3	1:D:360:VAL:HB	2.01	0.43
1:B:300:SER:HB2	1:C:188:LEU:HB3	2.01	0.43
1:B:452:GLU:OE1	1:B:452:GLU:HA	2.18	0.43
1:C:17:MET:HG3	1:C:34:ARG:HH12	1.83	0.43
1:C:22:VAL:CB	1:C:23:PRO:CD	2.93	0.43
1:C:341:ILE:HD13	1:D:349:TRP:CZ3	2.54	0.43
1:A:414:SER:OG	1:A:457:ARG:NH2	2.52	0.43
1:D:314:LEU:HD21	1:D:329:LEU:HD22	2.01	0.43
1:A:296:ARG:O	1:A:300:SER:HB3	2.19	0.43
1:A:367:ARG:NH1	2:A:475:HOH:O	2.52	0.43
1:C:445:ILE:H	1:C:445:ILE:CD1	2.11	0.43
1:C:419:GLU:HA	1:C:419:GLU:OE2	2.18	0.42
1:B:103:GLY:HA3	1:B:360:VAL:HB	2.01	0.42
1:C:443:GLY:C	1:C:445:ILE:HD12	2.39	0.42
1:D:453:ASP:OD2	1:D:457:ARG:NH1	2.51	0.42
1:C:11:ARG:NH1	1:C:24:ALA:O	2.53	0.42
1:B:417:GLY:N	1:B:420:GLU:OE1	2.43	0.42
1:A:159:LEU:O	1:A:163:LEU:HG	2.19	0.42
1:C:142:THR:HG22	1:C:143:PHE:N	2.33	0.42
1:C:452:GLU:O	1:C:456:ARG:HB2	2.20	0.42
1:C:429:LEU:HA	1:C:429:LEU:HD23	1.81	0.42
1:B:301:GLY:CA	1:C:189:MET:CE	2.98	0.42
1:D:367:ARG:HH11	1:D:367:ARG:HG2	1.84	0.42
1:A:376:LEU:HD23	1:A:376:LEU:HA	1.92	0.42
1:A:424:VAL:HG12	1:A:435:ILE:HD12	2.01	0.41
1:A:440:ILE:CG1	1:A:445:ILE:CD1	2.98	0.41
1:C:22:VAL:CG1	1:C:33:GLN:CD	2.72	0.41
1:D:138:SER:HB3	1:D:229:THR:HA	2.02	0.41
1:A:9:ASN:OD1	1:A:9:ASN:N	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:399:ARG:NH1	2:D:503:HOH:O	2.49	0.41
1:C:228:GLY:O	1:C:229:THR:OG1	2.30	0.41
1:C:314:LEU:C	1:C:316:PRO:HD3	2.41	0.41
1:C:454:LEU:HD22	1:C:458:LEU:HD22	2.02	0.41
1:A:357:GLU:HG2	1:C:185:ARG:O	2.20	0.41
1:B:190:ASP:OD1	1:C:304:THR:HG23	2.21	0.41
1:C:341:ILE:HD13	1:D:349:TRP:CH2	2.56	0.41
1:C:435:ILE:HG22	1:C:458:LEU:HD21	2.02	0.41
1:A:406:PRO:HB3	1:A:425:ALA:HB1	2.02	0.41
1:C:97:ILE:HD13	1:C:107:SER:HB3	2.01	0.41
1:D:465:LYS:O	1:D:466:ALA:C	2.57	0.41
1:B:401:LEU:HD23	1:B:401:LEU:HA	1.92	0.41
1:C:445:ILE:HD12	1:C:445:ILE:N	2.27	0.41
1:D:15:ASP:C	1:D:17:MET:H	2.23	0.41
1:A:183:SER:HB3	1:D:304:THR:HG21	2.04	0.41
1:A:188:LEU:HB3	1:D:300:SER:HB2	2.03	0.40
1:A:295:ILE:CA	1:A:298:MET:HE2	2.49	0.40
1:C:430:LYS:HB2	1:C:430:LYS:HE3	1.65	0.40
1:A:324:LYS:HZ2	1:A:324:LYS:HB2	1.86	0.40
1:A:421:ALA:O	1:A:424:VAL:N	2.55	0.40
1:B:405:SER:O	1:B:408:ILE:HG12	2.22	0.40
1:B:442:ARG:HB2	1:B:444:LEU:HD23	2.02	0.40
1:C:370:LEU:HD23	1:C:370:LEU:HA	1.96	0.40
1:B:296:ARG:HG3	1:C:188:LEU:CD1	2.52	0.40
1:D:228:GLY:O	1:D:229:THR:OG1	2.30	0.40
1:B:189:MET:HE3	1:C:301:GLY:CA	2.47	0.40
1:C:72:ALA:HA	1:C:73:PRO:HD2	1.90	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	448/475 (94%)	427 (95%)	21 (5%)	0	100	100
1	B	447/475 (94%)	432 (97%)	15 (3%)	0	100	100
1	C	448/475 (94%)	433 (97%)	12 (3%)	3 (1%)	22	36
1	D	447/475 (94%)	434 (97%)	12 (3%)	1 (0%)	47	66
All	All	1790/1900 (94%)	1726 (96%)	60 (3%)	4 (0%)	47	66

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	455	ASP
1	C	454	LEU
1	C	444	LEU
1	D	315	GLN

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	346/365 (95%)	331 (96%)	15 (4%)	29	50
1	B	345/365 (94%)	329 (95%)	16 (5%)	27	47
1	C	347/365 (95%)	329 (95%)	18 (5%)	23	41
1	D	345/365 (94%)	326 (94%)	19 (6%)	21	39
All	All	1383/1460 (95%)	1315 (95%)	68 (5%)	25	44

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	GLU
1	A	22	VAL
1	A	74	GLU
1	A	111	ASN
1	A	142	THR
1	A	183	SER
1	A	271	ARG

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Mol	Chain	Res	Type
1	A	329	LEU
1	A	357	GLU
1	A	379	VAL
1	A	382	LEU
1	A	399	ARG
1	A	414	SER
1	A	442	ARG
1	A	458	LEU
1	B	20	VAL
1	B	22	VAL
1	B	74	GLU
1	B	111	ASN
1	B	142	THR
1	B	183	SER
1	B	271	ARG
1	B	314	LEU
1	B	329	LEU
1	B	357	GLU
1	B	379	VAL
1	B	382	LEU
1	B	399	ARG
1	B	410	THR
1	B	450	SER
1	B	458	LEU
1	C	25	LYS
1	C	74	GLU
1	C	111	ASN
1	C	142	THR
1	C	183	SER
1	C	271	ARG
1	C	324	LYS
1	C	329	LEU
1	C	357	GLU
1	C	379	VAL
1	C	382	LEU
1	C	399	ARG
1	C	427	GLN
1	C	437	GLN
1	C	442	ARG
1	C	445	ILE
1	C	454	LEU
1	C	458	LEU

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Mol	Chain	Res	Type
1	D	22	VAL
1	D	74	GLU
1	D	111	ASN
1	D	142	THR
1	D	183	SER
1	D	271	ARG
1	D	329	LEU
1	D	357	GLU
1	D	379	VAL
1	D	382	LEU
1	D	399	ARG
1	D	407	SER
1	D	420	GLU
1	D	444	LEU
1	D	449	LEU
1	D	451	ILE
1	D	453	ASP
1	D	454	LEU
1	D	458	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	B	38	ASN
1	C	38	ASN
1	C	353	ASN
1	C	437	GLN
1	D	38	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

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	450/475 (94%)	0.76	63 (14%) 2 2	20, 48, 93, 125	17 (3%)
1	B	452/475 (95%)	0.37	23 (5%) 28 29	20, 46, 77, 113	24 (5%)
1	C	453/475 (95%)	0.78	61 (13%) 3 2	20, 49, 98, 122	18 (3%)
1	D	452/475 (95%)	0.39	24 (5%) 26 27	20, 46, 76, 128	24 (5%)
All	All	1807/1900 (95%)	0.58	171 (9%) 8 7	20, 47, 89, 128	83 (4%)

All (171) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	8	ALA	10.8
1	C	314	LEU	9.9
1	C	449	LEU	9.9
1	D	316	PRO	7.6
1	B	8	ALA	7.6
1	C	439	VAL	6.6
1	D	16	THR	6.5
1	C	447	ASP	6.4
1	A	439	VAL	6.3
1	C	409	VAL	5.8
1	C	20	VAL	5.5
1	C	412	LEU	5.3
1	D	323	GLY	5.3
1	A	447	ASP	5.3
1	C	124	GLY	5.2
1	A	453	ASP	5.1
1	C	8	ALA	5.1
1	C	425	ALA	5.1
1	D	314	LEU	5.1
1	D	10	TYR	5.0
1	A	449	LEU	5.0

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Mol	Chain	Res	Type	RSRZ
1	C	418	TYR	5.0
1	A	422	ALA	4.9
1	D	12	ILE	4.7
1	C	446	GLY	4.6
1	C	440	ILE	4.6
1	A	444	LEU	4.6
1	C	450	SER	4.6
1	C	121	ALA	4.4
1	B	453	ASP	4.4
1	A	19	GLU	4.3
1	A	119	ILE	4.3
1	D	315	GLN	4.3
1	D	313	ASP	4.2
1	A	429	LEU	4.2
1	B	449	LEU	4.2
1	A	418	TYR	4.2
1	C	413	ASN	4.1
1	A	125	VAL	4.0
1	C	453	ASP	4.0
1	A	451	ILE	4.0
1	B	25	LYS	3.9
1	C	441	ASP	3.9
1	B	10	TYR	3.9
1	A	10	TYR	3.6
1	C	315	GLN	3.6
1	C	29	ARG	3.6
1	B	440	ILE	3.6
1	C	445	ILE	3.5
1	A	92	ASP	3.5
1	C	416	ILE	3.5
1	D	9	ASN	3.5
1	A	455	ASP	3.4
1	B	16	THR	3.4
1	D	15	ASP	3.4
1	C	451	ILE	3.4
1	A	175	LEU	3.4
1	A	415	ALA	3.3
1	A	21	ARG	3.3
1	C	9	ASN	3.2
1	D	19	GLU	3.2
1	A	174	ALA	3.2
1	A	127	LEU	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	445	ILE	3.2
1	A	446	GLY	3.2
1	B	19	GLU	3.2
1	C	415	ALA	3.2
1	A	76	ALA	3.1
1	C	466	ALA	3.1
1	A	452	GLU	3.1
1	D	447	ASP	3.1
1	A	424	VAL	3.1
1	C	83	ALA	3.1
1	B	323	GLY	3.1
1	C	400	ARG	3.1
1	D	342	GLY	3.1
1	C	452	GLU	3.0
1	B	12	ILE	3.0
1	C	10	TYR	3.0
1	D	453	ASP	3.0
1	A	409	VAL	3.0
1	A	30	ALA	3.0
1	C	55	GLY	3.0
1	A	121	ALA	2.9
1	C	438	THR	2.9
1	D	83	ALA	2.9
1	C	323	GLY	2.8
1	B	23	PRO	2.8
1	A	8	ALA	2.8
1	C	22	VAL	2.8
1	C	125	VAL	2.8
1	A	412	LEU	2.7
1	B	314	LEU	2.7
1	A	342	GLY	2.7
1	A	440	ILE	2.7
1	A	420	GLU	2.7
1	C	455	ASP	2.7
1	B	341	ILE	2.7
1	C	414	SER	2.7
1	A	126	THR	2.6
1	C	120	ALA	2.6
1	A	20	VAL	2.6
1	B	275	VAL	2.6
1	C	180	VAL	2.6
1	A	28	TRP	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	438	THR	2.6
1	A	26	ALA	2.5
1	A	345	ALA	2.5
1	A	23	PRO	2.5
1	A	414	SER	2.5
1	A	454	LEU	2.5
1	D	449	LEU	2.5
1	A	425	ALA	2.4
1	C	442	ARG	2.4
1	C	345	ALA	2.4
1	A	278	SER	2.4
1	C	92	ASP	2.4
1	A	25	LYS	2.4
1	A	401	LEU	2.4
1	D	21	ARG	2.3
1	D	444	LEU	2.3
1	B	15	ASP	2.3
1	A	87	ALA	2.3
1	C	348	ALA	2.3
1	A	123	GLY	2.3
1	C	57	LEU	2.3
1	D	341	ILE	2.3
1	A	300	SER	2.3
1	A	115	VAL	2.3
1	A	83	ALA	2.3
1	B	447	ASP	2.2
1	C	435	ILE	2.2
1	C	288	LEU	2.2
1	C	313	ASP	2.2
1	C	420	GLU	2.2
1	A	17	MET	2.2
1	B	20	VAL	2.2
1	D	20	VAL	2.2
1	A	80	ILE	2.2
1	C	253	THR	2.2
1	D	88	ASP	2.2
1	C	456	ARG	2.2
1	C	444	LEU	2.2
1	D	76	ALA	2.2
1	A	15	ASP	2.2
1	A	450	SER	2.1
1	C	344	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	117	ALA	2.1
1	C	424	VAL	2.1
1	B	452	GLU	2.1
1	B	454	LEU	2.1
1	A	419	GLU	2.1
1	A	443	GLY	2.1
1	C	316	PRO	2.1
1	A	293	ASN	2.1
1	C	246	VAL	2.1
1	A	457	ARG	2.1
1	B	9	ASN	2.1
1	A	22	VAL	2.1
1	C	175	LEU	2.1
1	B	197	GLY	2.1
1	A	171	ALA	2.0
1	C	283	THR	2.0
1	A	400	ARG	2.0
1	C	457	ARG	2.0
1	A	280	ALA	2.0
1	C	30	ALA	2.0
1	D	345	ALA	2.0
1	C	115	VAL	2.0
1	B	81	ALA	2.0
1	B	119	ILE	2.0

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.