



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:24 PM UTC

PDB ID : 4MYC / pdb_00004myc
Title : Structure of the mitochondrial ABC transporter, Atm1
Authors : Srinivasan, V.
Deposited on : 2013-09-27
Resolution : 3.06 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

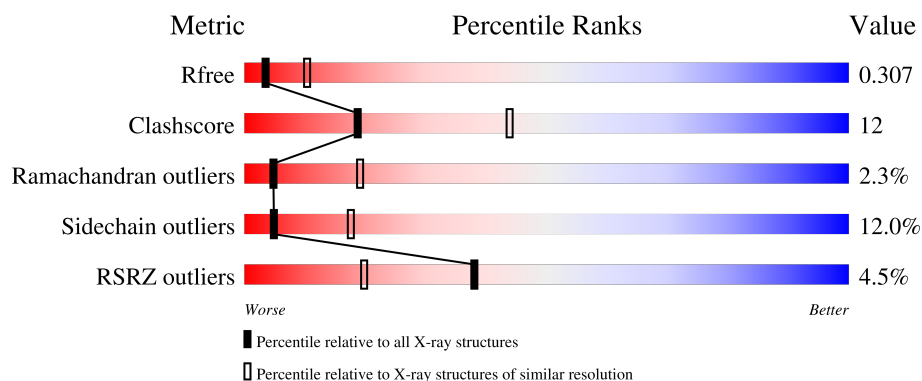
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.06 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	2469 (3.10-3.02)
Clashscore	190562	2569 (3.10-3.02)
Ramachandran outliers	187476	2424 (3.10-3.02)
Sidechain outliers	187428	2423 (3.10-3.02)
RSRZ outliers	180081	2469 (3.10-3.02)

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	598	5% 65% 29% 6%
1	B	598	4% 65% 30% 5%
1	C	598	4% 69% 27% • •

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 14028 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 Iron-sulfur clusters transporter ATM1, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	598	Total	C	N	O	S	0	0	0
			4676	3002	802	855	17			
1	C	598	Total	C	N	O	S	0	0	0
			4676	3002	802	855	17			
1	B	598	Total	C	N	O	S	0	0	0
			4676	3002	802	855	17			

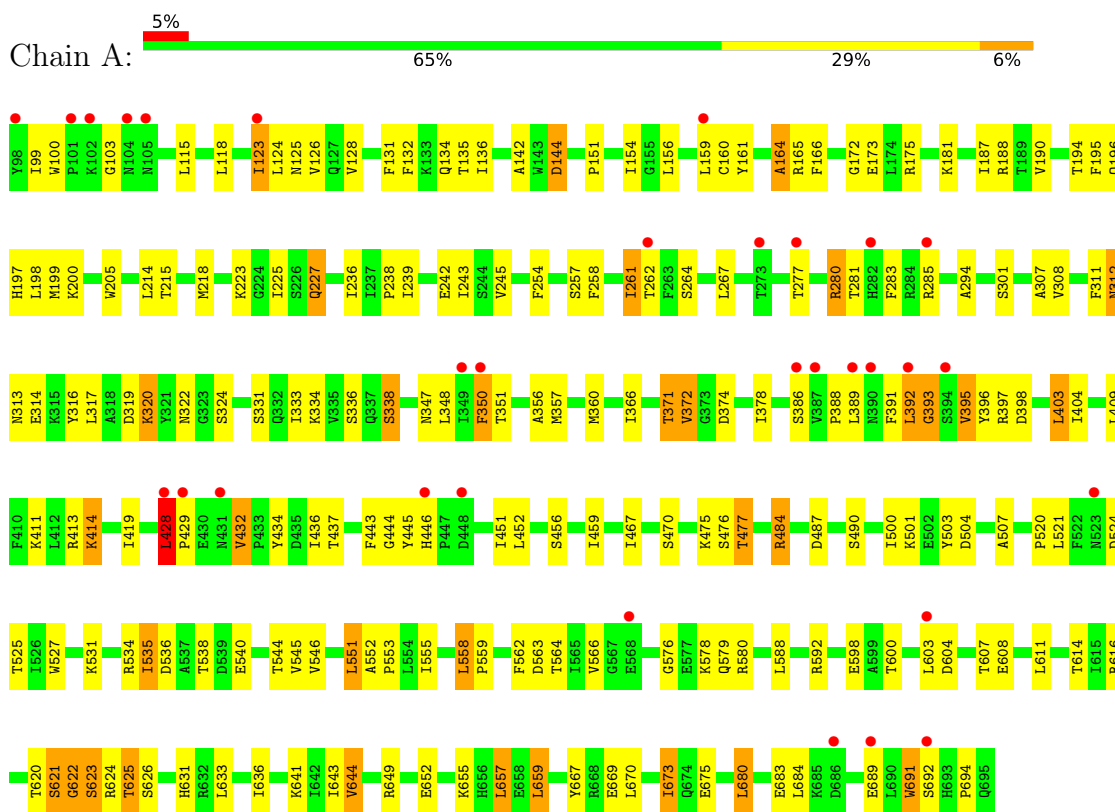
There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	691	TRP	-	expression tag	UNP P40416
A	692	SER	-	expression tag	UNP P40416
A	693	HIS	-	expression tag	UNP P40416
A	694	PRO	-	expression tag	UNP P40416
A	695	GLN	-	expression tag	UNP P40416
C	691	TRP	-	expression tag	UNP P40416
C	692	SER	-	expression tag	UNP P40416
C	693	HIS	-	expression tag	UNP P40416
C	694	PRO	-	expression tag	UNP P40416
C	695	GLN	-	expression tag	UNP P40416
B	691	TRP	-	expression tag	UNP P40416
B	692	SER	-	expression tag	UNP P40416
B	693	HIS	-	expression tag	UNP P40416
B	694	PRO	-	expression tag	UNP P40416
B	695	GLN	-	expression tag	UNP P40416

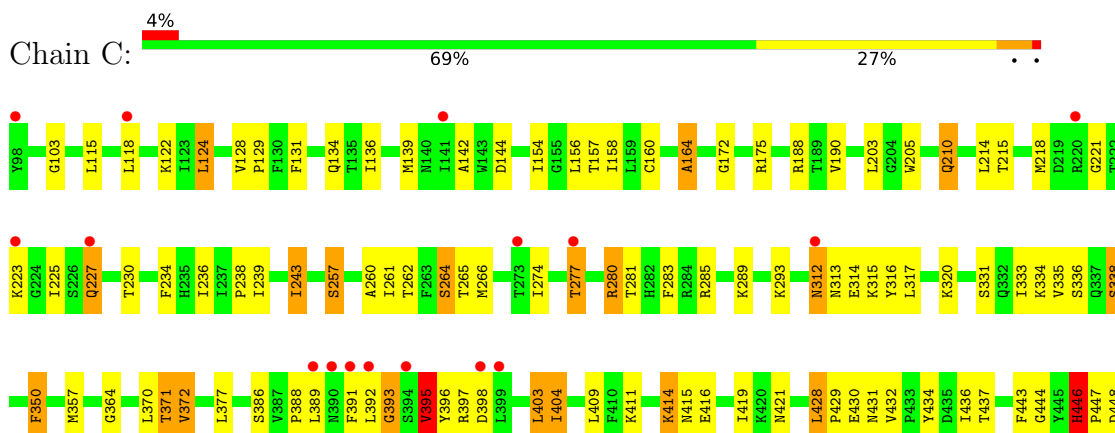
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry 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: Iron-sulfur clusters transporter ATM1, mitochondrial

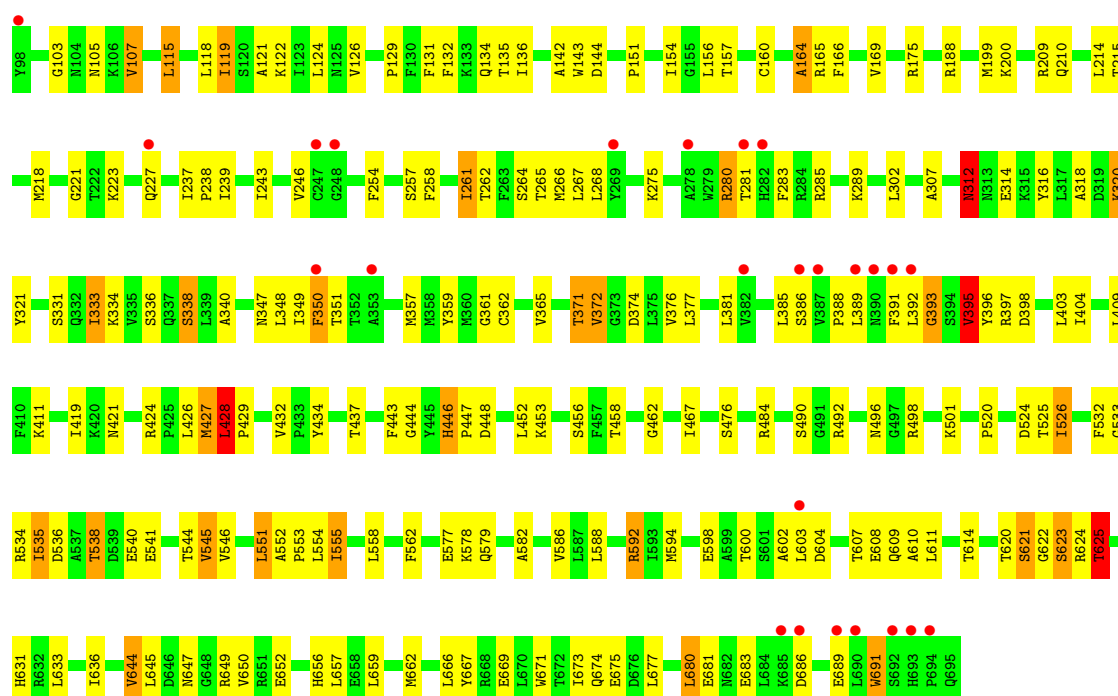


- Molecule 1: Iron-sulfur clusters transporter ATM1, mitochondrial





● Molecule 1: Iron-sulfur clusters transporter ATM1, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	156.62Å 156.62Å 520.56Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.60 – 3.06 48.60 – 3.06	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.60-3.06) 99.6 (48.60-3.06)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.251 , 0.298 0.263 , 0.307	Depositor DCC
R_{free} test set	3829 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	71.7	Xtriage
Anisotropy	0.115	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 11.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	14028	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.01% 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

5.1 Standard geometry

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	0.66	1/4763 (0.0%)	0.98	8/6451 (0.1%)
1	B	0.68	0/4763	1.01	11/6451 (0.2%)
1	C	0.68	0/4763	1.01	10/6451 (0.2%)
All	All	0.67	1/14289 (0.0%)	1.00	29/19353 (0.1%)

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
1	B	0	1
1	C	0	1
All	All	0	3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	128	VAL	CA-CB	5.38	1.57	1.53

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	428	LEU	CA-C-N	11.86	134.66	119.84
1	B	428	LEU	C-N-CA	11.86	134.66	119.84
1	B	427	MET	CA-C-O	7.16	124.76	119.68
1	A	100	TRP	CA-C-N	6.70	127.64	120.47
1	A	100	TRP	C-N-CA	6.70	127.64	120.47
1	C	446	HIS	CA-C-N	6.12	125.80	119.56
1	C	446	HIS	C-N-CA	6.12	125.80	119.56
1	B	200	LYS	N-CA-C	-6.08	105.86	113.28
1	A	558	LEU	CA-C-N	5.99	127.33	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	558	LEU	C-N-CA	5.99	127.33	119.84
1	B	526	ILE	CB-CA-C	-5.92	104.39	111.97
1	C	624	ARG	N-CA-C	-5.66	100.48	109.59
1	A	475	LYS	N-CA-C	-5.65	105.64	112.54
1	A	445	TYR	N-CA-C	-5.51	100.70	109.96
1	B	107	VAL	N-CA-C	-5.38	105.75	113.07
1	B	662	MET	CA-C-N	5.37	125.13	119.76
1	B	662	MET	C-N-CA	5.37	125.13	119.76
1	C	683	GLU	N-CA-C	-5.34	105.54	111.36
1	B	625	THR	CB-CA-C	5.34	118.84	110.19
1	B	320	LYS	N-CA-C	-5.31	105.66	111.82
1	A	320	LYS	N-CA-C	-5.25	105.73	111.82
1	C	625	THR	CB-CA-C	5.22	119.14	110.78
1	C	335	VAL	N-CA-C	-5.18	106.63	111.45
1	B	620	THR	N-CA-C	5.14	117.03	107.99
1	A	673	ILE	CB-CA-C	-5.11	105.43	111.97
1	C	643	ILE	CA-C-N	-5.05	116.57	122.93
1	C	643	ILE	C-N-CA	-5.05	116.57	122.93
1	C	128	VAL	CA-C-N	-5.02	113.87	119.19
1	C	128	VAL	C-N-CA	-5.02	113.87	119.19

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	428	LEU	Peptide
1	B	428	LEU	Peptide
1	C	428	LEU	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	4676	0	4738	136	0
1	B	4676	0	4738	131	0
1	C	4676	0	4738	103	0
All	All	14028	0	14214	335	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 (335) 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:453:LYS:NZ	1:C:647:ASN:OD1	2.04	0.90
1:A:546:VAL:HG13	1:A:551:LEU:HB3	1.57	0.86
1:B:526:ILE:HD11	1:B:558:LEU:HD12	1.57	0.86
1:A:631:HIS:HB2	1:B:691:TRP:HH2	1.47	0.80
1:C:261:ILE:HD11	1:C:357:MET:HG3	1.64	0.78
1:A:691:TRP:CH2	1:B:631:HIS:HB2	2.21	0.76
1:C:218:MET:HA	1:C:409:LEU:HD13	1.68	0.75
1:B:426:LEU:HB3	1:B:428:LEU:HD23	1.67	0.75
1:B:546:VAL:HG13	1:B:551:LEU:HB3	1.68	0.74
1:A:691:TRP:HH2	1:B:631:HIS:HB2	1.54	0.73
1:A:181:LYS:HA	1:B:333:ILE:HD11	1.70	0.72
1:C:436:ILE:HD11	1:C:593:ILE:HD13	1.71	0.72
1:A:261:ILE:HD11	1:A:357:MET:HG3	1.72	0.72
1:A:350:PHE:CE1	1:A:386:SER:HB3	2.27	0.69
1:C:124:LEU:HD13	1:C:164:ALA:HA	1.75	0.68
1:A:531:LYS:NZ	1:A:535:ILE:O	2.26	0.68
1:A:428:LEU:HB3	1:A:429:PRO:HD3	1.74	0.68
1:B:391:PHE:O	1:B:393:GLY:N	2.26	0.68
1:B:427:MET:H	1:B:428:LEU:HD22	1.59	0.67
1:A:484:ARG:NH1	1:A:501:LYS:HG2	2.11	0.66
1:A:280:ARG:NH1	1:A:338:SER:OG	2.29	0.66
1:B:124:LEU:HD22	1:B:164:ALA:HA	1.77	0.65
1:A:262:THR:OG1	1:A:357:MET:HE1	1.96	0.65
1:B:621:SER:O	1:B:623:SER:N	2.29	0.65
1:C:280:ARG:NH1	1:C:338:SER:OG	2.28	0.64
1:B:218:MET:HA	1:B:409:LEU:HD13	1.79	0.64
1:C:350:PHE:CE1	1:C:386:SER:HB3	2.31	0.64
1:B:434:TYR:O	1:B:625:THR:HG21	1.97	0.64
1:B:600:THR:HB	1:B:608:GLU:HG3	1.80	0.63
1:A:218:MET:HA	1:A:409:LEU:HD13	1.81	0.63
1:A:633:LEU:HD23	1:A:667:TYR:CE1	2.33	0.63
1:A:118:LEU:HD11	1:A:238:PRO:HG3	1.81	0.63
1:A:125:ASN:ND2	1:A:242:GLU:OE2	2.30	0.62
1:C:668:ARG:NH2	1:C:669:GLU:OE2	2.32	0.62
1:A:631:HIS:HB2	1:B:691:TRP:CH2	2.34	0.62
1:C:632:ARG:HD3	1:C:634:ARG:HH21	1.64	0.62
1:A:616:ARG:O	1:A:620:THR:OG1	2.17	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:453:LYS:NZ	1:B:647:ASN:OD1	2.32	0.62
1:B:280:ARG:NH1	1:B:338:SER:OG	2.27	0.61
1:B:633:LEU:HD23	1:B:667:TYR:CE1	2.35	0.61
1:A:118:LEU:HD13	1:A:175:ARG:HD3	1.80	0.61
1:A:484:ARG:NH1	1:A:487:ASP:OD2	2.33	0.61
1:A:683:GLU:HG2	1:B:680:LEU:HD21	1.82	0.61
1:A:118:LEU:HD13	1:A:175:ARG:HH11	1.66	0.60
1:A:391:PHE:O	1:A:393:GLY:N	2.33	0.60
1:C:131:PHE:CE2	1:C:160:CYS:HB3	2.36	0.60
1:A:124:LEU:HD12	1:A:164:ALA:HA	1.83	0.60
1:A:411:LYS:HE2	1:A:414:LYS:HZ1	1.66	0.60
1:C:633:LEU:HD23	1:C:667:TYR:CE1	2.37	0.59
1:C:391:PHE:O	1:C:393:GLY:N	2.34	0.59
1:B:266:MET:HE3	1:B:385:LEU:HD12	1.83	0.59
1:B:239:ILE:HD13	1:B:388:PRO:HG2	1.84	0.59
1:A:197:HIS:CD2	1:A:413:ARG:HA	2.39	0.58
1:A:680:LEU:HD21	1:B:683:GLU:HG2	1.84	0.58
1:A:644:VAL:HG13	1:A:652:GLU:HB2	1.85	0.58
1:A:131:PHE:CE2	1:A:160:CYS:HB3	2.39	0.58
1:B:261:ILE:HD11	1:B:357:MET:HG3	1.85	0.58
1:A:535:ILE:HG13	1:B:316:TYR:CD1	2.38	0.58
1:B:258:PHE:HD1	1:B:357:MET:HG2	1.68	0.58
1:B:645:LEU:HD23	1:B:650:VAL:HA	1.84	0.57
1:C:262:THR:OG1	1:C:357:MET:HE1	2.05	0.57
1:B:395:VAL:O	1:B:397:ARG:N	2.38	0.57
1:C:280:ARG:HH11	1:C:338:SER:HG	1.52	0.57
1:C:431:ASN:H	1:B:492:ARG:HH22	1.53	0.56
1:A:434:TYR:O	1:A:625:THR:HG21	2.05	0.56
1:A:205:TRP:HH2	1:A:214:LEU:HD21	1.71	0.56
1:A:395:VAL:O	1:A:397:ARG:N	2.38	0.56
1:C:462:GLY:H	1:C:625:THR:HB	1.69	0.56
1:A:536:ASP:OD2	1:A:536:ASP:N	2.35	0.56
1:B:524:ASP:OD1	1:B:525:THR:N	2.34	0.55
1:A:205:TRP:CH2	1:A:214:LEU:HD21	2.41	0.55
1:A:527:TRP:HD1	1:A:562:PHE:CG	2.24	0.55
1:C:395:VAL:O	1:C:397:ARG:N	2.40	0.55
1:C:536:ASP:OD2	1:C:536:ASP:N	2.37	0.55
1:B:421:ASN:OD1	1:B:421:ASN:N	2.40	0.54
1:A:411:LYS:HG3	1:A:414:LYS:HZ3	1.70	0.54
1:A:166:PHE:CE1	1:B:348:LEU:HD13	2.42	0.54
1:A:411:LYS:HG3	1:A:414:LYS:HD3	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:656:HIS:HE1	1:C:671:TRP:CZ3	2.26	0.54
1:B:350:PHE:CE1	1:B:386:SER:HB3	2.42	0.54
1:A:316:TYR:CD1	1:B:535:ILE:HG13	2.43	0.53
1:B:446:HIS:ND1	1:B:447:PRO:HD2	2.23	0.53
1:A:283:PHE:CE1	1:A:334:LYS:HB3	2.44	0.53
1:A:428:LEU:HB2	1:A:503:TYR:HE1	1.74	0.53
1:B:552:ALA:HB3	1:B:553:PRO:HD3	1.91	0.53
1:B:602:ALA:O	1:B:604:ASP:N	2.33	0.53
1:B:105:ASN:C	1:B:107:VAL:H	2.16	0.53
1:B:536:ASP:OD2	1:B:536:ASP:N	2.31	0.53
1:A:552:ALA:HB3	1:A:553:PRO:HD3	1.90	0.53
1:C:411:LYS:O	1:C:414:LYS:HB3	2.09	0.53
1:C:527:TRP:CZ3	1:C:531:LYS:HD2	2.43	0.53
1:A:527:TRP:CZ3	1:A:531:LYS:HD2	2.44	0.53
1:B:371:THR:OG1	1:B:372:VAL:N	2.41	0.52
1:A:411:LYS:HE2	1:A:414:LYS:NZ	2.24	0.52
1:C:609:GLN:O	1:C:613:ARG:HG3	2.09	0.52
1:A:131:PHE:HD2	1:A:161:TYR:HA	1.75	0.52
1:A:351:THR:HG23	1:B:165:ARG:HD3	1.91	0.52
1:C:283:PHE:HE1	1:C:334:LYS:HB3	1.75	0.52
1:C:446:HIS:ND1	1:C:447:PRO:HD2	2.24	0.51
1:A:428:LEU:HB3	1:A:429:PRO:CD	2.40	0.51
1:A:254:PHE:HE1	1:A:374:ASP:HB3	1.74	0.51
1:C:280:ARG:NH1	1:C:338:SER:HG	2.05	0.51
1:C:134:GLN:HB3	1:C:157:THR:HG21	1.92	0.51
1:B:677:LEU:O	1:B:681:GLU:HG2	2.11	0.51
1:C:118:LEU:HD13	1:C:175:ARG:HD3	1.91	0.51
1:B:594:MET:HE1	1:B:624:ARG:HH21	1.75	0.51
1:A:397:ARG:NH1	1:A:398:ASP:OD1	2.43	0.51
1:A:540:GLU:O	1:A:544:THR:HG23	2.10	0.51
1:B:540:GLU:O	1:B:544:THR:HG23	2.11	0.51
1:C:645:LEU:HD23	1:C:650:VAL:HA	1.93	0.50
1:B:462:GLY:H	1:B:625:THR:HG22	1.76	0.50
1:A:166:PHE:HE1	1:B:348:LEU:HB2	1.77	0.50
1:C:118:LEU:HD11	1:C:238:PRO:HB3	1.92	0.50
1:A:366:ILE:HG12	1:B:143:TRP:CE3	2.47	0.50
1:A:684:LEU:HA	1:B:680:LEU:HD23	1.92	0.50
1:A:198:LEU:HD13	1:B:302:LEU:HD21	1.94	0.50
1:C:632:ARG:HD3	1:C:634:ARG:NH2	2.26	0.50
1:B:496:ASN:O	1:B:498:ARG:HD2	2.12	0.50
1:B:644:VAL:HG13	1:B:652:GLU:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:600:THR:HB	1:C:608:GLU:HG3	1.94	0.50
1:A:451:ILE:HG21	1:A:477:THR:HG21	1.94	0.49
1:B:669:GLU:O	1:B:673:ILE:HG13	2.12	0.49
1:C:432:VAL:O	1:C:592:ARG:NH1	2.44	0.49
1:C:546:VAL:HG13	1:C:551:LEU:HB3	1.93	0.49
1:A:236:ILE:HD13	1:A:392:LEU:HD11	1.94	0.49
1:C:277:THR:HA	1:C:280:ARG:HG2	1.95	0.49
1:C:411:LYS:HG3	1:C:414:LYS:HD3	1.94	0.49
1:B:448:ASP:OD1	1:B:448:ASP:N	2.41	0.49
1:A:294:ALA:HB2	1:A:324:SER:HB3	1.95	0.48
1:C:154:ILE:O	1:C:158:ILE:HG23	2.13	0.48
1:C:239:ILE:HD13	1:C:388:PRO:HG2	1.94	0.48
1:A:195:PHE:CE2	1:B:318:ALA:HA	2.48	0.48
1:A:348:LEU:HD13	1:B:166:PHE:CD1	2.48	0.48
1:C:397:ARG:NH1	1:C:398:ASP:OD1	2.46	0.48
1:C:313:ASN:HA	1:C:316:TYR:HB3	1.95	0.48
1:C:350:PHE:CD1	1:C:386:SER:HB3	2.48	0.48
1:A:411:LYS:O	1:A:414:LYS:HB3	2.14	0.48
1:A:181:LYS:HG2	1:B:333:ILE:HD13	1.95	0.47
1:A:165:ARG:HD3	1:B:351:THR:HG23	1.97	0.47
1:C:540:GLU:O	1:C:544:THR:HG23	2.15	0.47
1:B:538:THR:HG23	1:B:541:GLU:CD	2.40	0.47
1:A:520:PRO:HG3	1:B:307:ALA:HB2	1.95	0.47
1:B:359:TYR:O	1:B:362:CYS:N	2.48	0.47
1:B:582:ALA:O	1:B:586:VAL:HG13	2.14	0.47
1:C:129:PRO:HB2	1:C:377:LEU:HD12	1.97	0.47
1:C:434:TYR:O	1:C:625:THR:HG21	2.14	0.47
1:A:307:ALA:HB2	1:B:520:PRO:HG3	1.97	0.47
1:B:246:VAL:HG22	1:B:381:LEU:HG	1.96	0.47
1:C:371:THR:OG1	1:C:372:VAL:N	2.48	0.47
1:C:448:ASP:N	1:C:448:ASP:OD1	2.48	0.47
1:C:552:ALA:HB3	1:C:553:PRO:HD3	1.96	0.47
1:C:656:HIS:CE1	1:C:671:TRP:CZ3	3.03	0.47
1:B:361:GLY:O	1:B:365:VAL:HG23	2.14	0.47
1:A:576:GLY:O	1:A:580:ARG:HB2	2.15	0.47
1:B:122:LYS:HE3	1:B:238:PRO:O	2.15	0.47
1:C:274:ILE:O	1:C:277:THR:HG22	2.15	0.46
1:B:652:GLU:OE2	1:B:666:LEU:N	2.26	0.46
1:A:669:GLU:O	1:A:673:ILE:HG13	2.15	0.46
1:C:421:ASN:OD1	1:C:421:ASN:N	2.44	0.46
1:B:131:PHE:CD2	1:B:160:CYS:HB3	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:173:GLU:HB3	1:B:340:ALA:HB1	1.96	0.46
1:A:521:LEU:HD13	1:A:566:VAL:HB	1.97	0.46
1:A:633:LEU:O	1:A:636:ILE:HG13	2.15	0.46
1:B:214:LEU:O	1:B:218:MET:HG2	2.15	0.46
1:A:604:ASP:OD2	1:A:607:THR:OG1	2.27	0.46
1:A:301:SER:OG	1:A:317:LEU:HD13	2.15	0.46
1:A:500:ILE:HA	1:A:503:TYR:HD2	1.80	0.46
1:C:555:ILE:HG21	1:C:562:PHE:CE1	2.51	0.46
1:C:524:ASP:OD1	1:C:525:THR:N	2.43	0.46
1:A:620:THR:O	1:A:622:GLY:N	2.49	0.46
1:B:656:HIS:HE1	1:B:671:TRP:CZ3	2.33	0.46
1:A:283:PHE:HE1	1:A:334:LYS:HD3	1.80	0.46
1:B:267:LEU:HD12	1:B:267:LEU:HA	1.76	0.46
1:A:239:ILE:HD13	1:A:388:PRO:HG2	1.98	0.45
1:C:443:PHE:CG	1:C:444:GLY:N	2.84	0.45
1:C:462:GLY:H	1:C:625:THR:CB	2.28	0.45
1:C:644:VAL:HG13	1:C:652:GLU:HB2	1.97	0.45
1:A:196:GLN:O	1:A:200:LYS:HG3	2.17	0.45
1:A:436:ILE:HB	1:A:459:ILE:HB	1.99	0.45
1:A:558:LEU:HA	1:A:559:PRO:HD2	1.71	0.45
1:C:429:PRO:HD2	1:C:434:TYR:HE1	1.81	0.45
1:A:126:VAL:HG11	1:A:245:VAL:HG12	1.98	0.45
1:A:347:ASN:CG	1:B:169:VAL:HG21	2.42	0.45
1:C:260:ALA:O	1:C:264:SER:HB2	2.17	0.45
1:C:504:ASP:O	1:C:507:ALA:N	2.49	0.45
1:C:548:LYS:HB3	1:C:618:ASN:OD1	2.17	0.45
1:B:389:LEU:HA	1:B:389:LEU:HD12	1.73	0.45
1:B:462:GLY:H	1:B:625:THR:CG2	2.29	0.45
1:C:403:LEU:HD22	1:C:403:LEU:HA	1.69	0.45
1:B:427:MET:H	1:B:428:LEU:CD2	2.29	0.45
1:A:308:VAL:HG11	1:A:317:LEU:HD12	1.98	0.45
1:A:371:THR:OG1	1:A:372:VAL:N	2.49	0.45
1:A:680:LEU:HD12	1:A:680:LEU:HA	1.67	0.45
1:A:313:ASN:HA	1:A:316:TYR:HB3	1.97	0.45
1:A:670:LEU:HA	1:A:670:LEU:HD23	1.75	0.45
1:C:686:ASP:O	1:C:689:GLU:HB3	2.16	0.45
1:B:118:LEU:HD13	1:B:175:ARG:HH11	1.82	0.45
1:B:126:VAL:O	1:B:129:PRO:HD2	2.17	0.45
1:A:504:ASP:O	1:A:507:ALA:N	2.49	0.45
1:B:209:ARG:HH12	1:B:411:LYS:HZ1	1.63	0.45
1:B:258:PHE:HB3	1:B:357:MET:HE2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:594:MET:HE2	1:B:594:MET:HB2	1.83	0.44
1:A:524:ASP:OD1	1:A:525:THR:N	2.40	0.44
1:A:225:ILE:C	1:A:227:GLN:H	2.25	0.44
1:A:283:PHE:HE1	1:A:334:LYS:HB3	1.82	0.44
1:C:317:LEU:HD23	1:C:317:LEU:HA	1.67	0.44
1:B:428:LEU:HB3	1:B:498:ARG:NH1	2.32	0.44
1:A:600:THR:HB	1:A:608:GLU:HG3	1.99	0.44
1:C:484:ARG:NH1	1:C:501:LYS:HG2	2.32	0.44
1:A:356:ALA:O	1:A:360:MET:HG3	2.17	0.44
1:A:621:SER:O	1:A:623:SER:N	2.47	0.44
1:B:131:PHE:CE2	1:B:160:CYS:HB3	2.53	0.44
1:A:432:VAL:O	1:A:592:ARG:NH1	2.51	0.44
1:A:655:LYS:O	1:A:659:LEU:HB2	2.18	0.44
1:A:657:LEU:HD12	1:A:657:LEU:HA	1.81	0.44
1:C:257:SER:O	1:C:261:ILE:HG23	2.18	0.44
1:C:266:MET:HE1	1:C:350:PHE:CZ	2.53	0.44
1:A:172:GLY:O	1:A:175:ARG:HB3	2.18	0.44
1:C:558:LEU:HD13	1:C:564:THR:HG21	2.00	0.44
1:B:268:LEU:HD23	1:B:349:ILE:HG12	2.00	0.44
1:B:554:LEU:HD13	1:B:577:GLU:CD	2.43	0.44
1:C:122:LYS:HE3	1:C:238:PRO:O	2.17	0.43
1:C:604:ASP:OD1	1:C:607:THR:OG1	2.26	0.43
1:C:680:LEU:HD12	1:C:680:LEU:HA	1.78	0.43
1:B:432:VAL:O	1:B:592:ARG:NH1	2.50	0.43
1:A:181:LYS:HG2	1:B:333:ILE:CD1	2.48	0.43
1:A:622:GLY:O	1:A:624:ARG:HG3	2.17	0.43
1:A:694:PRO:HB2	1:B:647:ASN:HD22	1.84	0.43
1:C:205:TRP:CH2	1:C:214:LEU:HD21	2.53	0.43
1:C:230:THR:O	1:C:234:PHE:HB2	2.19	0.43
1:C:502:GLU:OE2	1:B:424:ARG:HD3	2.18	0.43
1:C:689:GLU:O	1:C:692:SER:HB3	2.19	0.43
1:B:283:PHE:CE1	1:B:334:LYS:HB3	2.53	0.43
1:A:132:PHE:HA	1:A:135:THR:HG22	2.01	0.43
1:A:258:PHE:HE1	1:A:378:ILE:HD12	1.84	0.43
1:B:554:LEU:HD22	1:B:577:GLU:HB3	2.00	0.43
1:B:555:ILE:HG21	1:B:562:PHE:CE1	2.53	0.43
1:A:280:ARG:O	1:A:283:PHE:HB2	2.19	0.43
1:A:452:LEU:HD12	1:A:452:LEU:HA	1.88	0.43
1:A:199:MET:HE1	1:B:314:GLU:HA	2.01	0.43
1:C:243:ILE:HG12	1:C:266:MET:HG3	2.00	0.43
1:B:118:LEU:HD11	1:B:238:PRO:HB3	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:554:LEU:HD13	1:C:577:GLU:CD	2.43	0.43
1:C:556:LYS:HA	1:C:556:LYS:HD3	1.77	0.43
1:B:686:ASP:O	1:B:689:GLU:HB3	2.18	0.43
1:C:554:LEU:HD22	1:C:577:GLU:HB3	2.00	0.43
1:B:209:ARG:NH1	1:B:411:LYS:HZ1	2.17	0.43
1:B:644:VAL:CG1	1:B:652:GLU:HB2	2.49	0.43
1:A:311:PHE:CD1	1:B:533:GLY:HA2	2.54	0.42
1:B:262:THR:OG1	1:B:357:MET:HE1	2.18	0.42
1:B:283:PHE:HE1	1:B:334:LYS:HB3	1.84	0.42
1:A:579:GLN:HB2	1:A:611:LEU:HD11	2.01	0.42
1:A:644:VAL:CG1	1:A:652:GLU:HB2	2.48	0.42
1:C:139:MET:HB3	1:C:139:MET:HE3	1.72	0.42
1:C:210:GLN:CD	1:C:210:GLN:H	2.26	0.42
1:C:214:LEU:O	1:C:218:MET:HG2	2.19	0.42
1:A:123:ILE:HD13	1:A:123:ILE:HA	1.75	0.42
1:A:319:ASP:HA	1:A:322:ASN:HB2	2.01	0.42
1:A:308:VAL:HG12	1:B:532:PHE:HZ	1.84	0.42
1:A:689:GLU:O	1:A:692:SER:HB3	2.19	0.42
1:C:313:ASN:O	1:C:315:LYS:N	2.52	0.42
1:A:403:LEU:HD22	1:A:403:LEU:HA	1.82	0.42
1:C:280:ARG:O	1:C:283:PHE:HB2	2.19	0.42
1:A:267:LEU:HD12	1:A:267:LEU:HA	1.67	0.42
1:A:389:LEU:HD12	1:A:389:LEU:HA	1.80	0.42
1:C:221:GLY:O	1:C:225:ILE:HG13	2.20	0.42
1:C:549:ALA:O	1:C:580:ARG:HG3	2.20	0.42
1:B:397:ARG:NH1	1:B:398:ASP:OD1	2.52	0.42
1:A:151:PRO:HB2	1:A:154:ILE:HG12	2.02	0.42
1:B:254:PHE:HB3	1:B:258:PHE:CD2	2.54	0.42
1:B:428:LEU:HB3	1:B:498:ARG:HH12	1.85	0.42
1:B:579:GLN:HB2	1:B:611:LEU:HD11	2.02	0.42
1:C:389:LEU:HD12	1:C:389:LEU:HA	1.79	0.42
1:B:452:LEU:HD12	1:B:452:LEU:HA	1.86	0.42
1:A:277:THR:HA	1:A:280:ARG:HG2	2.02	0.42
1:B:132:PHE:HA	1:B:135:THR:HG22	2.02	0.42
1:B:281:THR:C	1:B:283:PHE:H	2.27	0.42
1:A:258:PHE:CE1	1:A:378:ILE:HD12	2.55	0.41
1:C:462:GLY:H	1:C:625:THR:CG2	2.33	0.41
1:B:115:LEU:O	1:B:118:LEU:HB3	2.20	0.41
1:B:680:LEU:HD12	1:B:680:LEU:HA	1.66	0.41
1:A:641:LYS:HE2	1:A:643:ILE:HD11	2.02	0.41
1:C:283:PHE:CE1	1:C:334:LYS:HB3	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:545:VAL:HG11	1:B:588:LEU:HG	2.02	0.41
1:B:604:ASP:OD2	1:B:607:THR:OG1	2.25	0.41
1:A:470:SER:HA	1:B:691:TRP:CH2	2.55	0.41
1:A:525:THR:HA	1:A:564:THR:O	2.20	0.41
1:A:694:PRO:CB	1:B:647:ASN:HD22	2.34	0.41
1:C:225:ILE:C	1:C:227:GLN:H	2.27	0.41
1:C:587:LEU:HD23	1:C:587:LEU:HA	1.91	0.41
1:B:151:PRO:O	1:B:154:ILE:HG12	2.20	0.41
1:A:525:THR:HG22	1:A:563:ASP:O	2.21	0.41
1:C:293:LYS:HD3	1:C:293:LYS:HA	1.73	0.41
1:B:134:GLN:HB3	1:B:157:THR:HG21	2.02	0.41
1:B:446:HIS:ND1	1:B:448:ASP:OD1	2.53	0.41
1:C:364:GLY:O	1:C:370:LEU:N	2.45	0.41
1:B:280:ARG:O	1:B:283:PHE:HB2	2.20	0.41
1:B:443:PHE:CG	1:B:444:GLY:N	2.88	0.41
1:B:633:LEU:O	1:B:636:ILE:HG13	2.20	0.41
1:A:527:TRP:HZ3	1:A:531:LYS:HD2	1.84	0.41
1:C:234:PHE:N	1:C:234:PHE:CD1	2.88	0.41
1:C:414:LYS:HE2	1:C:415:ASN:ND2	2.35	0.41
1:C:561:GLY:O	1:C:564:THR:OG1	2.30	0.41
1:C:633:LEU:HD12	1:C:633:LEU:H	1.85	0.41
1:A:239:ILE:HG12	1:A:389:LEU:HD13	2.02	0.41
1:B:119:ILE:C	1:B:121:ALA:H	2.29	0.41
1:B:610:ALA:O	1:B:614:THR:HG23	2.20	0.41
1:C:576:GLY:O	1:C:580:ARG:HB2	2.20	0.41
1:B:221:GLY:HA3	1:B:409:LEU:HD22	2.02	0.41
1:A:261:ILE:HG12	1:A:357:MET:HE3	2.03	0.41
1:A:314:GLU:HA	1:B:199:MET:HE1	2.03	0.41
1:A:545:VAL:HG11	1:A:588:LEU:HG	2.03	0.41
1:C:404:ILE:H	1:C:404:ILE:HG12	1.72	0.41
1:C:429:PRO:HD2	1:C:434:TYR:CE1	2.56	0.41
1:C:608:GLU:OE2	1:C:632:ARG:HD2	2.21	0.41
1:B:254:PHE:HE1	1:B:374:ASP:HB3	1.85	0.41
1:B:312:ASN:OD1	1:B:312:ASN:N	2.42	0.41
1:B:318:ALA:O	1:B:321:TYR:HB3	2.21	0.41
1:A:443:PHE:CG	1:A:444:GLY:N	2.88	0.41
1:C:203:LEU:HD22	1:C:485:PHE:CD2	2.55	0.41
1:C:281:THR:C	1:C:283:PHE:H	2.29	0.41
1:A:281:THR:C	1:A:283:PHE:H	2.28	0.40
1:A:546:VAL:HG13	1:A:551:LEU:CB	2.38	0.40
1:C:218:MET:HE3	1:C:218:MET:HB2	1.80	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:PHE:HA	1:B:347:ASN:OD1	2.22	0.40
1:C:172:GLY:O	1:C:175:ARG:HB3	2.21	0.40
1:B:254:PHE:CE1	1:B:374:ASP:HB3	2.56	0.40
1:A:173:GLU:HB3	1:B:340:ALA:CB	2.52	0.40
1:A:239:ILE:HD12	1:A:239:ILE:HA	1.88	0.40
1:A:254:PHE:CE1	1:A:374:ASP:HB3	2.56	0.40
1:B:280:ARG:HH11	1:B:338:SER:HG	1.63	0.40
1:A:144:ASP:OD2	1:A:144:ASP:N	2.47	0.40
1:B:129:PRO:HB2	1:B:377:LEU:CD1	2.51	0.40
1:B:484:ARG:NH1	1:B:501:LYS:HG2	2.36	0.40
1:C:158:ILE:HD13	1:C:158:ILE:HG21	1.87	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	596/598 (100%)	540 (91%)	42 (7%)	14 (2%)	5	19
1	B	596/598 (100%)	540 (91%)	42 (7%)	14 (2%)	5	19
1	C	596/598 (100%)	544 (91%)	38 (6%)	14 (2%)	5	19
All	All	1788/1794 (100%)	1624 (91%)	122 (7%)	42 (2%)	5	19

All (42) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	312	ASN
1	A	392	LEU
1	A	395	VAL
1	A	396	TYR
1	A	603	LEU
1	A	621	SER

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Mol	Chain	Res	Type
1	C	312	ASN
1	C	392	LEU
1	C	395	VAL
1	C	396	TYR
1	C	603	LEU
1	C	621	SER
1	B	312	ASN
1	B	392	LEU
1	B	395	VAL
1	B	396	TYR
1	B	603	LEU
1	A	164	ALA
1	C	164	ALA
1	B	164	ALA
1	B	393	GLY
1	B	429	PRO
1	B	621	SER
1	A	393	GLY
1	C	393	GLY
1	B	142	ALA
1	A	99	ILE
1	A	142	ALA
1	A	446	HIS
1	C	142	ALA
1	C	446	HIS
1	B	428	LEU
1	B	446	HIS
1	B	622	GLY
1	A	428	LEU
1	A	622	GLY
1	C	314	GLU
1	A	103	GLY
1	C	428	LEU
1	C	430	GLU
1	C	103	GLY
1	B	103	GLY

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	501/514 (98%)	441 (88%)	60 (12%)	5	17
1	B	501/514 (98%)	440 (88%)	61 (12%)	5	17
1	C	501/514 (98%)	441 (88%)	60 (12%)	5	17
All	All	1503/1542 (98%)	1322 (88%)	181 (12%)	5	17

All (181) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	115	LEU
1	A	123	ILE
1	A	134	GLN
1	A	136	ILE
1	A	144	ASP
1	A	156	LEU
1	A	159	LEU
1	A	187	ILE
1	A	188	ARG
1	A	190	VAL
1	A	194	THR
1	A	215	THR
1	A	223	LYS
1	A	227	GLN
1	A	243	ILE
1	A	257	SER
1	A	261	ILE
1	A	264	SER
1	A	280	ARG
1	A	285	ARG
1	A	312	ASN
1	A	320	LYS
1	A	331	SER
1	A	333	ILE
1	A	336	SER
1	A	338	SER
1	A	350	PHE
1	A	371	THR
1	A	372	VAL
1	A	403	LEU
1	A	404	ILE

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Mol	Chain	Res	Type
1	A	414	LYS
1	A	419	ILE
1	A	428	LEU
1	A	432	VAL
1	A	437	THR
1	A	456	SER
1	A	467	ILE
1	A	476	SER
1	A	477	THR
1	A	484	ARG
1	A	490	SER
1	A	534	ARG
1	A	535	ILE
1	A	538	THR
1	A	551	LEU
1	A	555	ILE
1	A	578	LYS
1	A	598	GLU
1	A	614	THR
1	A	623	SER
1	A	625	THR
1	A	626	SER
1	A	644	VAL
1	A	649	ARG
1	A	657	LEU
1	A	659	LEU
1	A	675	GLU
1	A	680	LEU
1	A	691	TRP
1	C	115	LEU
1	C	124	LEU
1	C	136	ILE
1	C	144	ASP
1	C	156	LEU
1	C	188	ARG
1	C	190	VAL
1	C	210	GLN
1	C	215	THR
1	C	223	LYS
1	C	227	GLN
1	C	236	ILE
1	C	243	ILE

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Mol	Chain	Res	Type
1	C	257	SER
1	C	264	SER
1	C	265	THR
1	C	277	THR
1	C	280	ARG
1	C	285	ARG
1	C	289	LYS
1	C	312	ASN
1	C	320	LYS
1	C	331	SER
1	C	333	ILE
1	C	336	SER
1	C	338	SER
1	C	350	PHE
1	C	371	THR
1	C	372	VAL
1	C	395	VAL
1	C	403	LEU
1	C	404	ILE
1	C	414	LYS
1	C	416	GLU
1	C	419	ILE
1	C	437	THR
1	C	456	SER
1	C	467	ILE
1	C	476	SER
1	C	477	THR
1	C	490	SER
1	C	534	ARG
1	C	535	ILE
1	C	538	THR
1	C	551	LEU
1	C	555	ILE
1	C	578	LYS
1	C	592	ARG
1	C	598	GLU
1	C	609	GLN
1	C	614	THR
1	C	623	SER
1	C	625	THR
1	C	644	VAL
1	C	649	ARG

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Mol	Chain	Res	Type
1	C	657	LEU
1	C	659	LEU
1	C	675	GLU
1	C	680	LEU
1	C	690	LEU
1	B	115	LEU
1	B	119	ILE
1	B	136	ILE
1	B	144	ASP
1	B	156	LEU
1	B	188	ARG
1	B	210	GLN
1	B	215	THR
1	B	223	LYS
1	B	227	GLN
1	B	237	ILE
1	B	243	ILE
1	B	257	SER
1	B	261	ILE
1	B	264	SER
1	B	265	THR
1	B	275	LYS
1	B	280	ARG
1	B	285	ARG
1	B	289	LYS
1	B	312	ASN
1	B	320	LYS
1	B	331	SER
1	B	333	ILE
1	B	336	SER
1	B	338	SER
1	B	350	PHE
1	B	371	THR
1	B	372	VAL
1	B	376	VAL
1	B	395	VAL
1	B	403	LEU
1	B	404	ILE
1	B	419	ILE
1	B	428	LEU
1	B	437	THR
1	B	456	SER

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Mol	Chain	Res	Type
1	B	458	THR
1	B	467	ILE
1	B	476	SER
1	B	490	SER
1	B	534	ARG
1	B	535	ILE
1	B	538	THR
1	B	545	VAL
1	B	551	LEU
1	B	555	ILE
1	B	578	LYS
1	B	592	ARG
1	B	598	GLU
1	B	609	GLN
1	B	623	SER
1	B	625	THR
1	B	644	VAL
1	B	649	ARG
1	B	657	LEU
1	B	659	LEU
1	B	674	GLN
1	B	675	GLU
1	B	680	LEU
1	B	691	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	134	GLN
1	A	590	ASN
1	C	196	GLN
1	C	227	GLN
1	C	682	ASN
1	B	134	GLN
1	B	210	GLN
1	B	227	GLN
1	B	253	GLN
1	B	379	ASN
1	B	674	GLN
1	B	679	HIS
1	B	682	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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	598/598 (100%)	0.25	31 (5%)	33 16	5, 41, 61, 87	0
1	B	598/598 (100%)	0.06	25 (4%)	40 21	4, 37, 57, 77	0
1	C	598/598 (100%)	-0.08	25 (4%)	40 21	4, 28, 49, 69	0
All	All	1794/1794 (100%)	0.08	81 (4%)	38 20	4, 33, 57, 87	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	692	SER	9.3
1	B	689	GLU	6.8
1	C	389	LEU	5.3
1	A	390	ASN	5.3
1	C	394	SER	4.9
1	A	104	ASN	4.7
1	A	568	GLU	4.6
1	C	689	GLU	4.6
1	A	448	ASP	4.6
1	A	389	LEU	4.4
1	A	603	LEU	4.3
1	C	390	ASN	4.1
1	C	603	LEU	4.1
1	C	227	GLN	4.1
1	C	98	TYR	4.0
1	A	98	TYR	3.8
1	C	399	LEU	3.8
1	A	689	GLU	3.7
1	A	394	SER	3.7
1	C	392	LEU	3.6
1	B	389	LEU	3.6
1	A	431	ASN	3.5
1	B	386	SER	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	387	VAL	3.4
1	A	101	PRO	3.3
1	C	391	PHE	3.3
1	C	647	ASN	3.3
1	A	446	HIS	3.2
1	B	392	LEU	3.2
1	B	282	HIS	3.1
1	A	692	SER	3.1
1	A	350	PHE	3.0
1	B	382	VAL	2.9
1	B	686	ASP	2.9
1	A	273	THR	2.9
1	C	694	PRO	2.9
1	B	390	ASN	2.9
1	C	692	SER	2.9
1	B	690	LEU	2.9
1	A	386	SER	2.8
1	A	262	THR	2.8
1	C	398	ASP	2.8
1	B	98	TYR	2.8
1	A	392	LEU	2.8
1	B	603	LEU	2.8
1	B	350	PHE	2.7
1	A	686	ASP	2.7
1	B	694	PRO	2.6
1	B	685	LYS	2.6
1	A	428	LEU	2.6
1	C	118	LEU	2.6
1	C	312	ASN	2.5
1	A	387	VAL	2.5
1	A	285	ARG	2.5
1	C	648	GLY	2.5
1	C	141	ILE	2.4
1	B	391	PHE	2.4
1	A	105	ASN	2.4
1	A	277	THR	2.3
1	B	248	GLY	2.3
1	C	273	THR	2.3
1	A	123	ILE	2.2
1	A	159	LEU	2.2
1	C	649	ARG	2.2
1	C	690	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	353	ALA	2.2
1	B	693	HIS	2.2
1	C	223	LYS	2.2
1	A	282	HIS	2.2
1	C	277	THR	2.1
1	B	281	THR	2.1
1	A	523	ASN	2.1
1	A	102	LYS	2.1
1	C	220	ARG	2.1
1	B	227	GLN	2.1
1	B	278	ALA	2.1
1	A	349	ILE	2.1
1	C	693	HIS	2.0
1	B	269	TYR	2.0
1	B	247	CYS	2.0
1	A	429	PRO	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 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.