



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 27, 2026 – 01:14 AM UTC

PDB ID : 4NQK / pdb_00004nqk
Title : Structure of an Ubiquitin complex
Authors : Peisley, A.; Wu, B.; Hur, S.
Deposited on : 2013-11-25
Resolution : 3.70 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

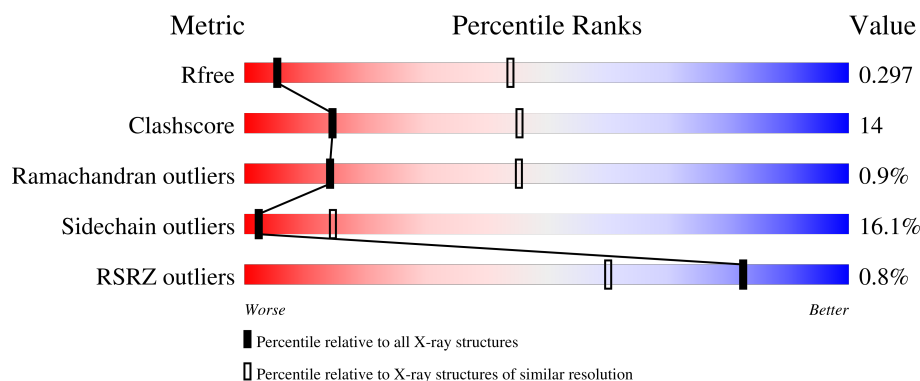
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.70 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	1131 (3.80-3.60)
Clashscore	190562	1171 (3.80-3.60)
Ramachandran outliers	187476	1129 (3.80-3.60)
Sidechain outliers	187428	1126 (3.80-3.60)
RSRZ outliers	180081	1130 (3.80-3.60)

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	203	
1	B	203	
1	C	203	
1	D	203	
2	E	79	

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Mol	Chain	Length	Quality of chain
2	F	79	% 58% 29% 5% 8%
2	G	79	70% 16% • 10%
2	H	79	% 71% 20% • 5%
2	I	79	4% 53% 22% 11% • 10%
2	J	79	% 70% 18% 5% 8%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 19139 atoms, of which 9584 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 Probable ATP-dependent RNA helicase DDX58.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	188	Total	C	H	N	O	S	0	0	0
			3041	990	1510	245	287	9			
1	B	188	Total	C	H	N	O	S	0	0	0
			3034	989	1505	243	288	9			
1	C	188	Total	C	H	N	O	S	0	0	0
			3000	981	1487	242	281	9			
1	D	188	Total	C	H	N	O	S	0	0	0
			3081	998	1535	248	291	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP O95786
A	-1	PRO	-	expression tag	UNP O95786
A	0	GLY	-	expression tag	UNP O95786
B	-2	GLY	-	expression tag	UNP O95786
B	-1	PRO	-	expression tag	UNP O95786
B	0	GLY	-	expression tag	UNP O95786
C	-2	GLY	-	expression tag	UNP O95786
C	-1	PRO	-	expression tag	UNP O95786
C	0	GLY	-	expression tag	UNP O95786
D	-2	GLY	-	expression tag	UNP O95786
D	-1	PRO	-	expression tag	UNP O95786
D	0	GLY	-	expression tag	UNP O95786

- Molecule 2 is a protein called Ubiquitin.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	E	72	Total	C	H	N	O	S	0	0	0
			1169	362	595	98	113	1			
2	F	73	Total	C	H	N	O	S	0	0	0
			1171	365	595	96	114	1			

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	G	71	Total	C	H	N	O	S	0	0	0
			1128	352	570	93	112	1			
2	H	75	Total	C	H	N	O	S	0	0	0
			1202	373	611	101	116	1			
2	I	71	Total	C	H	N	O	S	0	0	0
			1125	352	570	92	110	1			
2	J	73	Total	C	H	N	O	S	0	0	0
			1188	368	606	99	114	1			

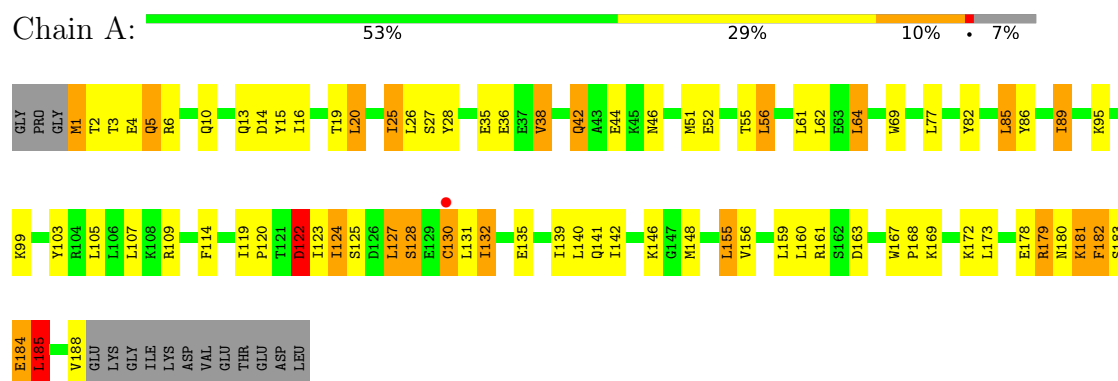
There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	-2	GLY	-	expression tag	UNP P0CG48
E	-1	PRO	-	expression tag	UNP P0CG48
F	-2	GLY	-	expression tag	UNP P0CG48
F	-1	PRO	-	expression tag	UNP P0CG48
G	-2	GLY	-	expression tag	UNP P0CG48
G	-1	PRO	-	expression tag	UNP P0CG48
H	-2	GLY	-	expression tag	UNP P0CG48
H	-1	PRO	-	expression tag	UNP P0CG48
I	-2	GLY	-	expression tag	UNP P0CG48
I	-1	PRO	-	expression tag	UNP P0CG48
J	-2	GLY	-	expression tag	UNP P0CG48
J	-1	PRO	-	expression tag	UNP P0CG48

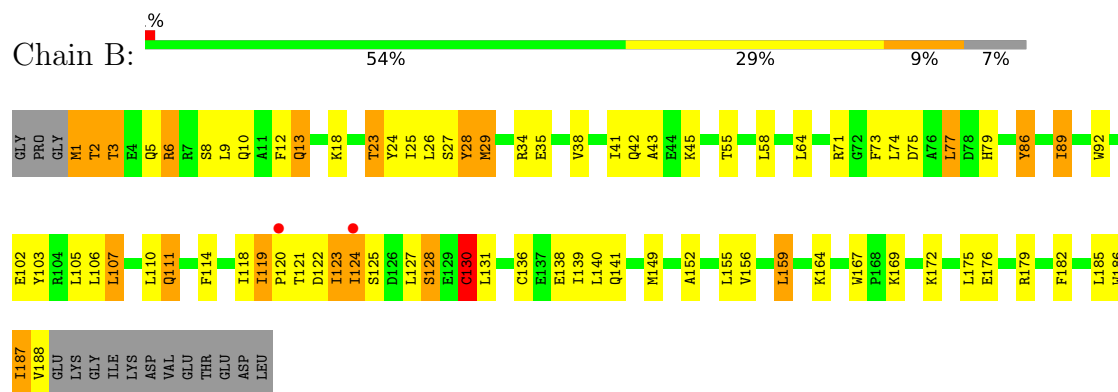
3 Residue-property plots

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.

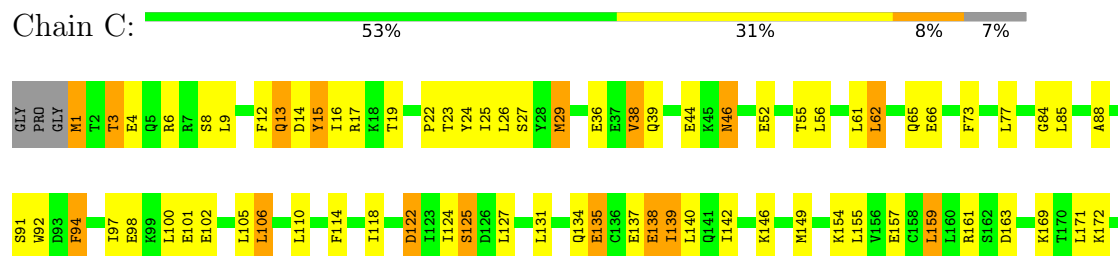
- Molecule 1: Probable ATP-dependent RNA helicase DDX58



- Molecule 1: Probable ATP-dependent RNA helicase DDX58



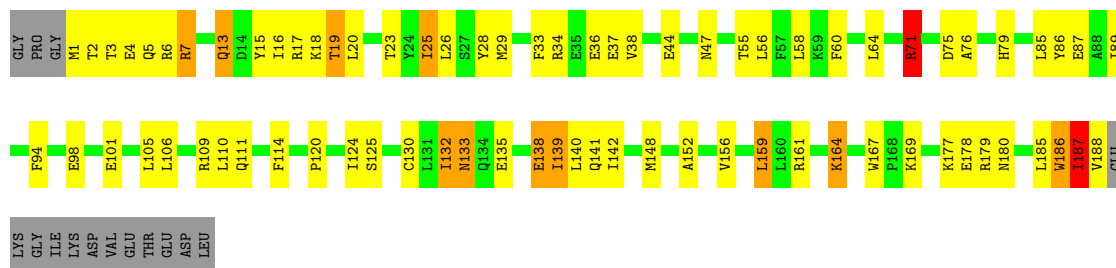
- Molecule 1: Probable ATP-dependent RNA helicase DDX58





- Molecule 1: Probable ATP-dependent RNA helicase DDX58

Chain D: 55% 31% 5% 7%



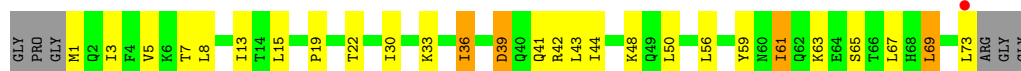
- Molecule 2: Ubiquitin

Chain E: 54% 32% 5% 9%



- Molecule 2: Ubiquitin

Chain F: 58% 29% 5% 8%



- Molecule 2: Ubiquitin

Chain G: 70% 16% 10% 4%



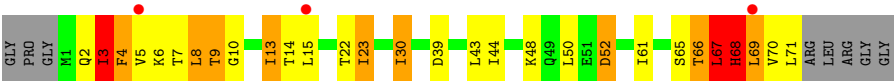
- Molecule 2: Ubiquitin

Chain H: 71% 20% 5% 4%

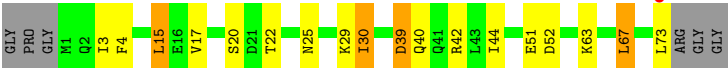


- Molecule 2: Ubiquitin

Chain I: 53% 22% 11% 10%



● Molecule 2: Ubiquitin



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	85.54Å 101.85Å 88.15Å 90.00° 106.88° 90.00°	Depositor
Resolution (Å)	46.00 – 3.70 46.00 – 3.70	Depositor EDS
% Data completeness (in resolution range)	98.5 (46.00-3.70) 98.5 (46.00-3.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.16	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.17 (at 3.48Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.222 , 0.285 0.235 , 0.297	Depositor DCC
R_{free} test set	903 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	79.2	Xtriage
Anisotropy	0.189	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 84.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.003 for l,k,h	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	19139	wwPDB-VP
Average B, all atoms (Å ²)	93.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.10% 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

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	0.64	0/1563	1.17	9/2110 (0.4%)
1	B	0.62	0/1561	1.17	9/2108 (0.4%)
1	C	0.55	0/1545	1.02	8/2088 (0.4%)
1	D	0.57	0/1578	1.07	6/2128 (0.3%)
2	E	0.44	0/580	0.90	1/781 (0.1%)
2	F	0.42	0/582	0.84	0/785
2	G	0.46	0/564	0.88	1/761 (0.1%)
2	H	0.45	0/597	0.85	0/804
2	I	1.03	7/561 (1.2%)	1.45	16/757 (2.1%)
2	J	0.42	0/588	0.87	0/792
All	All	0.59	7/9719 (0.1%)	1.07	50/13114 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	D	0	1
2	I	0	3
All	All	0	6

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	6	LYS	CA-C	-6.19	1.45	1.52
2	I	2	GLN	CA-C	6.13	1.60	1.52
2	I	67	LEU	CA-C	-6.08	1.45	1.52
2	I	3	ILE	CA-C	5.90	1.60	1.52
2	I	67	LEU	C-O	5.79	1.30	1.23
2	I	4	PHE	CA-C	5.74	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	4	PHE	C-O	5.22	1.30	1.23

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	67	LEU	N-CA-C	-11.98	89.83	109.96
1	B	2	THR	N-CA-C	9.15	121.02	111.14
1	B	128	SER	N-CA-C	-8.26	102.36	111.36
1	D	138	GLU	N-CA-C	-8.19	98.96	110.50
1	A	130	CYS	N-CA-C	7.70	122.50	113.18
2	I	7	THR	N-CA-C	-7.46	100.69	110.53
2	I	67	LEU	O-C-N	7.33	131.66	123.16
1	A	128	SER	N-CA-C	-7.29	103.41	111.36
1	D	28	TYR	N-CA-C	7.20	119.21	111.36
1	D	140	LEU	N-CA-C	-7.13	103.44	111.14
1	A	38	VAL	N-CA-C	-6.99	103.66	110.72
1	C	15	TYR	N-CA-C	-6.77	103.98	111.36
1	B	130	CYS	N-CA-C	6.71	121.07	113.01
2	I	69	LEU	CA-C-N	6.36	132.73	122.68
2	I	69	LEU	C-N-CA	6.36	132.73	122.68
1	D	94	PHE	N-CA-C	6.30	120.67	112.92
2	I	6	LYS	N-CA-C	-6.28	96.84	108.02
1	B	43	ALA	N-CA-C	-6.24	104.01	111.69
2	I	2	GLN	N-CA-C	6.21	119.37	108.75
1	B	124	ILE	N-CA-CB	6.14	117.73	110.55
2	I	67	LEU	N-CA-CB	6.11	120.13	110.23
1	C	38	VAL	N-CA-C	-6.11	104.55	110.72
2	I	4	PHE	CA-C-O	5.91	127.79	121.11
1	A	28	TYR	N-CA-C	5.81	117.69	111.36
1	B	128	SER	N-CA-CB	5.71	118.62	110.16
1	C	65	GLN	N-CA-C	5.70	120.24	112.04
2	I	4	PHE	N-CA-C	5.69	118.83	109.72
1	A	35	GLU	N-CA-C	-5.67	105.19	111.71
1	B	28	TYR	N-CA-C	5.62	117.49	111.36
2	I	69	LEU	O-C-N	5.59	129.71	122.89
1	A	85	LEU	N-CA-C	-5.53	104.66	111.40
2	I	68	HIS	CA-C-N	-5.50	113.06	120.87
2	I	68	HIS	C-N-CA	-5.50	113.06	120.87
2	G	3	ILE	N-CA-C	5.42	115.15	108.53
1	D	85	LEU	N-CA-C	-5.42	105.02	111.69
2	I	68	HIS	N-CA-C	-5.42	99.89	108.73
1	A	5	GLN	N-CA-C	-5.35	105.49	112.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	25	ILE	N-CA-C	-5.33	106.08	113.00
1	C	179	ARG	CB-CA-C	-5.33	110.41	116.54
1	A	122	ASP	N-CA-C	-5.33	105.48	111.28
2	I	4	PHE	CA-C-N	-5.32	115.43	122.98
2	I	4	PHE	C-N-CA	-5.32	115.43	122.98
1	B	179	ARG	CB-CA-C	-5.29	110.46	116.54
1	C	118	ILE	N-CA-C	5.25	116.54	108.71
1	A	15	TYR	N-CA-C	-5.25	105.00	111.40
2	E	12	THR	N-CA-C	5.22	117.68	108.75
1	C	114	PHE	N-CA-C	-5.21	105.49	111.07
1	C	94	PHE	N-CA-C	5.16	119.27	112.92
1	D	71	ARG	CG-CD-NE	5.11	123.24	112.00
1	B	111	GLN	N-CA-C	5.09	121.05	109.81

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	179	ARG	Peptide
1	B	13	GLN	Peptide
1	D	186	TRP	Peptide
2	I	3	ILE	Peptide
2	I	66	THR	Peptide
2	I	68	HIS	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1531	1510	1510	50	0
1	B	1529	1505	1505	54	0
1	C	1513	1487	1487	54	0
1	D	1546	1535	1535	45	0
2	E	574	595	599	18	0
2	F	576	595	599	13	0
2	G	558	570	574	5	0
2	H	591	611	615	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	I	555	570	574	28	0
2	J	582	606	610	7	0
All	All	9555	9584	9608	269	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 14.

All (269) 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)
2:I:5:VAL:HG13	2:I:69:LEU:HB3	1.45	0.98
1:A:135:GLU:OE1	1:A:161:ARG:NH1	2.17	0.77
1:B:107:LEU:HD23	1:B:156:VAL:HG13	1.69	0.74
1:C:154:LYS:NZ	1:C:157:GLU:OE2	2.19	0.74
2:I:5:VAL:HG13	2:I:69:LEU:CB	2.17	0.74
1:D:179:ARG:CB	1:D:180:ASN:HA	2.17	0.73
2:I:5:VAL:HB	2:I:13:ILE:HG23	1.70	0.73
1:C:6:ARG:NH1	1:C:62:LEU:O	2.22	0.72
1:D:15:TYR:O	1:D:19:THR:OG1	2.05	0.72
1:C:91:SER:O	2:H:42:ARG:NH2	2.25	0.70
1:B:74:LEU:HD21	1:B:89:ILE:HD11	1.74	0.69
2:E:39:ASP:O	2:E:42:ARG:NH1	2.27	0.68
1:B:124:ILE:O	1:B:128:SER:N	2.29	0.66
1:B:121:THR:O	1:B:124:ILE:HG22	1.96	0.65
1:C:105:LEU:HD12	1:C:106:LEU:N	2.11	0.65
1:C:139:ILE:O	1:C:142:ILE:N	2.31	0.64
1:A:125:SER:O	1:A:128:SER:OG	2.11	0.64
2:E:55:THR:OG1	2:E:58:ASP:OD1	2.15	0.63
1:A:26:LEU:HD13	1:A:38:VAL:HG13	1.81	0.63
2:J:39:ASP:OD1	2:J:40:GLN:N	2.32	0.63
1:B:42:GLN:HA	1:B:45:LYS:HB3	1.82	0.62
1:B:2:THR:HG22	1:B:6:ARG:HD2	1.81	0.62
2:I:61:ILE:HD13	2:I:67:LEU:HD22	1.81	0.61
1:D:25:ILE:HG22	1:D:76:ALA:HB1	1.83	0.60
1:A:184:GLU:O	1:A:185:LEU:HB2	2.00	0.60
1:C:105:LEU:HD13	2:I:8:LEU:HB2	1.83	0.60
2:I:3:ILE:HD12	2:I:67:LEU:HD21	1.84	0.59
2:I:5:VAL:HG21	2:I:30:ILE:HD11	1.84	0.59
1:A:124:ILE:O	1:A:128:SER:N	2.35	0.59
1:B:26:LEU:HD13	1:B:38:VAL:HG13	1.84	0.58
1:A:103:TYR:CE1	1:A:188:VAL:HG12	2.38	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:135:GLU:OE2	1:D:161:ARG:NH1	2.37	0.57
1:C:1:MET:HE2	1:C:1:MET:O	2.05	0.57
2:I:5:VAL:HG11	2:I:30:ILE:HD11	1.85	0.57
1:A:42:GLN:O	1:A:46:ASN:ND2	2.38	0.56
2:J:3:ILE:HD11	2:J:17:VAL:HG21	1.88	0.56
1:A:44:GLU:HG3	1:A:56:LEU:HD13	1.88	0.55
1:B:103:TYR:CE1	1:B:188:VAL:HG21	2.42	0.55
1:D:186:TRP:O	1:D:187:ILE:HG13	2.06	0.55
1:C:149:MET:HE1	1:D:133:ASN:HB3	1.88	0.55
2:I:9:THR:OG1	2:I:10:GLY:N	2.35	0.55
1:B:123:ILE:HD11	1:B:155:LEU:HD11	1.88	0.55
1:A:51:MET:HE1	1:B:35:GLU:HB2	1.88	0.55
1:D:179:ARG:CB	1:D:180:ASN:CA	2.85	0.55
1:C:122:ASP:OD1	1:C:122:ASP:N	2.41	0.54
1:C:149:MET:HE1	1:D:133:ASN:CB	2.38	0.54
2:E:18:GLU:N	2:E:21:ASP:OD2	2.38	0.54
1:B:182:PHE:O	1:B:186:TRP:N	2.37	0.54
1:B:152:ALA:O	1:B:156:VAL:HG23	2.09	0.53
1:A:20:LEU:HD12	1:A:82:TYR:CE1	2.43	0.53
1:C:26:LEU:HD22	1:C:38:VAL:HG13	1.90	0.53
1:C:88:ALA:O	1:C:92:TRP:N	2.42	0.53
1:D:44:GLU:HG3	1:D:56:LEU:HD13	1.91	0.53
1:A:179:ARG:O	1:A:181:LYS:N	2.39	0.53
1:A:155:LEU:HD22	1:A:159:LEU:HD13	1.90	0.53
2:I:5:VAL:CG1	2:I:69:LEU:HB3	2.29	0.52
1:A:82:TYR:HE2	1:A:132:ILE:HG12	1.73	0.52
2:H:39:ASP:OD1	2:H:40:GLN:N	2.42	0.52
2:I:13:ILE:HD11	2:I:15:LEU:HD13	1.91	0.52
1:A:4:GLU:OE1	1:A:4:GLU:N	2.31	0.52
2:I:4:PHE:O	2:I:67:LEU:O	2.27	0.52
2:I:69:LEU:HG	2:I:70:VAL:N	2.24	0.52
1:B:175:LEU:HD22	1:B:182:PHE:HB2	1.92	0.52
1:B:124:ILE:HD11	1:B:136:CYS:HA	1.92	0.52
2:G:23:ILE:HB	2:G:52:ASP:HA	1.91	0.52
1:A:2:THR:O	1:A:5:GLN:HG2	2.09	0.51
2:F:19:PRO:HA	2:F:56:LEU:HB2	1.91	0.51
1:B:73:PHE:CZ	1:B:77:LEU:HD11	2.45	0.51
1:D:138:GLU:O	1:D:139:ILE:HB	2.10	0.51
1:A:1:MET:SD	1:A:2:THR:OG1	2.58	0.51
2:E:27:LYS:O	2:E:41:GLN:NE2	2.37	0.51
1:C:105:LEU:CD1	2:I:8:LEU:HB2	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:177:LYS:HG3	1:D:178:GLU:N	2.27	0.50
1:D:2:THR:O	1:D:5:GLN:HG2	2.12	0.50
1:A:132:ILE:HD12	1:A:132:ILE:H	1.76	0.50
1:C:22:PRO:O	1:C:26:LEU:N	2.43	0.50
2:E:6:LYS:HG3	2:E:66:THR:CG2	2.42	0.50
1:D:139:ILE:HA	1:D:142:ILE:HB	1.94	0.50
1:B:127:LEU:HA	1:B:130:CYS:HB2	1.93	0.49
1:A:120:PRO:HD2	1:A:148:MET:HG3	1.95	0.49
2:E:61:ILE:HG23	2:E:65:SER:HB2	1.94	0.49
1:B:107:LEU:HD12	1:B:186:TRP:CZ3	2.47	0.49
2:I:23:ILE:HB	2:I:52:ASP:HA	1.93	0.49
1:D:152:ALA:O	1:D:156:VAL:HG23	2.12	0.49
1:A:109:ARG:HG3	2:E:8:LEU:CD1	2.43	0.49
1:A:16:ILE:CD1	1:A:85:LEU:HD21	2.43	0.48
1:B:28:TYR:CZ	1:B:79:HIS:HB3	2.49	0.48
1:C:106:LEU:HD13	1:C:106:LEU:C	2.38	0.48
1:A:131:LEU:HD21	1:A:159:LEU:HD12	1.95	0.48
1:B:125:SER:O	1:B:128:SER:OG	2.27	0.48
1:D:7:ARG:HG3	1:D:7:ARG:HH11	1.78	0.48
2:G:19:PRO:HA	2:G:56:LEU:HB2	1.96	0.48
1:C:29:MET:CE	1:C:73:PHE:HA	2.44	0.48
1:C:46:ASN:O	1:D:47:ASN:ND2	2.44	0.48
2:H:59:TYR:HB2	2:H:61:ILE:CD1	2.43	0.48
1:C:24:TYR:CE1	1:C:134:GLN:HG2	2.49	0.48
2:J:3:ILE:CD1	2:J:17:VAL:HG21	2.43	0.48
1:A:19:THR:HG22	1:A:85:LEU:HD12	1.94	0.47
2:I:3:ILE:HA	2:I:4:PHE:CD1	2.48	0.47
1:C:135:GLU:OE1	1:C:161:ARG:NH1	2.47	0.47
1:A:36:GLU:C	1:A:38:VAL:H	2.22	0.47
2:I:23:ILE:HD13	2:I:23:ILE:N	2.29	0.47
1:B:2:THR:HA	1:B:5:GLN:HE21	1.80	0.47
1:B:149:MET:N	1:C:137:GLU:OE2	2.43	0.47
1:C:138:GLU:O	1:C:140:LEU:N	2.48	0.47
2:F:7:THR:HG22	2:F:69:LEU:HD12	1.97	0.47
2:F:69:LEU:C	2:F:69:LEU:HD13	2.40	0.47
2:I:5:VAL:HG11	2:I:30:ILE:CD1	2.45	0.47
1:C:105:LEU:HD22	2:I:8:LEU:O	2.15	0.47
1:C:155:LEU:HD23	1:C:159:LEU:HD13	1.97	0.47
1:D:2:THR:HG22	1:D:6:ARG:HD2	1.97	0.47
1:D:132:ILE:N	1:D:132:ILE:HD12	2.29	0.47
1:A:142:ILE:O	1:A:146:LYS:N	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:13:ILE:HG12	2:I:14:THR:N	2.30	0.46
1:A:52:GLU:O	1:A:55:THR:HG22	2.15	0.46
1:B:25:ILE:O	1:B:29:MET:N	2.47	0.46
1:C:44:GLU:HG3	1:C:56:LEU:HD22	1.96	0.46
2:E:61:ILE:HG23	2:E:65:SER:CB	2.44	0.46
1:B:1:MET:SD	1:B:3:THR:HG23	2.56	0.46
1:D:98:GLU:O	1:D:101:GLU:HB3	2.15	0.46
1:B:89:ILE:HD12	1:B:89:ILE:C	2.41	0.46
2:F:41:GLN:C	2:F:42:ARG:HD2	2.40	0.46
2:I:8:LEU:HD13	2:I:8:LEU:H	1.81	0.46
1:B:23:THR:CG2	1:B:42:GLN:HG2	2.46	0.46
2:F:43:LEU:C	2:F:44:ILE:HD12	2.41	0.46
2:E:15:LEU:HD21	2:E:30:ILE:HG13	1.98	0.46
2:G:60:ASN:O	2:G:60:ASN:ND2	2.41	0.46
1:A:127:LEU:HA	1:A:130:CYS:HB2	1.97	0.45
1:A:172:LYS:NZ	1:A:183:SER:OG	2.49	0.45
1:C:15:TYR:OH	1:C:98:GLU:OE1	2.29	0.45
1:D:20:LEU:HD21	1:D:25:ILE:HD11	1.98	0.45
2:F:1:MET:HE3	2:F:63:LYS:HA	1.97	0.45
2:E:7:THR:HG22	2:E:11:LYS:HB3	1.98	0.45
2:G:8:LEU:HD11	2:G:70:VAL:HG23	1.97	0.45
1:A:123:ILE:HG13	1:A:124:ILE:N	2.31	0.45
1:B:75:ASP:O	1:B:79:HIS:ND1	2.44	0.45
1:B:138:GLU:C	1:B:140:LEU:H	2.24	0.45
1:B:111:GLN:HG2	1:B:156:VAL:HG11	1.97	0.45
1:C:139:ILE:O	1:C:140:LEU:C	2.58	0.45
1:D:132:ILE:HB	1:D:135:GLU:HG3	1.97	0.45
2:E:6:LYS:HG3	2:E:66:THR:HG22	1.99	0.45
1:C:24:TYR:CD1	1:C:24:TYR:N	2.83	0.45
1:C:88:ALA:HB1	1:C:94:PHE:CD1	2.51	0.45
2:E:5:VAL:HG11	2:E:30:ILE:HD11	1.98	0.45
1:D:23:THR:O	1:D:26:LEU:HB2	2.17	0.45
1:B:127:LEU:C	1:B:130:CYS:H	2.24	0.45
1:C:12:PHE:CE2	1:C:94:PHE:CD1	3.05	0.45
1:C:142:ILE:O	1:C:146:LYS:N	2.48	0.45
1:C:14:ASP:HA	1:C:17:ARG:HG2	1.99	0.45
1:D:16:ILE:HD12	1:D:16:ILE:H	1.81	0.45
2:I:43:LEU:HD23	2:I:69:LEU:HA	1.98	0.45
1:B:24:TYR:O	1:B:27:SER:HB2	2.17	0.45
1:B:86:TYR:CD1	1:B:86:TYR:C	2.95	0.44
2:F:36:ILE:HD13	2:F:36:ILE:N	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:23:ILE:HB	2:E:52:ASP:HA	1.97	0.44
1:C:19:THR:CG2	1:C:163:ASP:HB3	2.48	0.44
1:B:119:ILE:N	1:B:119:ILE:HD12	2.33	0.44
2:F:61:ILE:HG21	2:F:67:LEU:HD21	1.99	0.44
1:D:86:TYR:CD2	1:D:86:TYR:C	2.95	0.44
2:I:67:LEU:HD23	2:I:67:LEU:H	1.83	0.44
1:A:16:ILE:HD11	1:A:89:ILE:HD11	2.00	0.44
1:B:120:PRO:O	1:B:124:ILE:HB	2.17	0.44
1:D:13:GLN:HA	1:D:16:ILE:HD13	1.99	0.44
1:D:56:LEU:HD21	1:D:60:PHE:CE2	2.52	0.44
1:D:177:LYS:CG	1:D:178:GLU:N	2.80	0.44
2:H:8:LEU:HD21	2:H:70:VAL:HG22	2.00	0.44
2:I:48:LYS:O	2:I:50:LEU:HD12	2.17	0.44
2:H:13:ILE:HG22	2:H:15:LEU:HD13	1.99	0.44
1:A:6:ARG:NH2	1:A:62:LEU:O	2.50	0.44
1:A:124:ILE:O	1:A:128:SER:HB3	2.17	0.44
1:A:169:LYS:O	1:A:173:LEU:HD22	2.18	0.44
1:C:106:LEU:HD12	1:C:186:TRP:CZ3	2.51	0.44
1:A:64:LEU:HD23	1:A:69:TRP:HB2	2.00	0.44
1:B:131:LEU:HD21	1:B:159:LEU:HD12	1.99	0.44
1:C:177:LYS:CG	1:C:178:GLU:N	2.81	0.44
1:D:33:PHE:CE1	1:D:64:LEU:HD11	2.53	0.44
1:B:124:ILE:O	1:B:128:SER:HB3	2.18	0.43
2:J:3:ILE:HG22	2:J:4:PHE:H	1.83	0.43
1:C:138:GLU:HG2	1:C:139:ILE:H	1.83	0.43
1:A:123:ILE:O	1:A:127:LEU:HG	2.18	0.43
1:B:12:PHE:HD1	1:B:12:PHE:O	2.01	0.43
1:A:114:PHE:C	1:A:114:PHE:CD2	2.96	0.43
1:A:167:TRP:N	1:A:168:PRO:HD2	2.33	0.43
1:C:84:GLY:N	1:C:163:ASP:OD2	2.52	0.43
2:I:66:THR:HG22	2:I:68:HIS:ND1	2.34	0.43
1:B:149:MET:H	1:C:137:GLU:CD	2.27	0.43
1:B:164:LYS:HB2	1:B:167:TRP:HB3	2.01	0.43
1:B:172:LYS:HD2	1:B:187:ILE:HA	1.99	0.43
1:D:34:ARG:O	1:D:38:VAL:HG23	2.19	0.43
1:D:111:GLN:HG2	1:D:156:VAL:HG21	2.00	0.43
1:D:75:ASP:O	1:D:79:HIS:ND1	2.51	0.43
2:F:5:VAL:HB	2:F:13:ILE:HG23	2.00	0.43
1:A:20:LEU:HD11	1:A:25:ILE:CD1	2.49	0.43
2:H:45:PHE:HB2	2:H:67:LEU:CD1	2.49	0.43
2:H:45:PHE:HB2	2:H:67:LEU:HD13	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:13:GLN:O	1:C:14:ASP:HB3	2.19	0.43
1:C:131:LEU:HD21	1:C:159:LEU:HD12	2.01	0.43
1:A:13:GLN:O	1:A:14:ASP:HB3	2.18	0.42
1:A:155:LEU:CD2	1:A:159:LEU:HD13	2.49	0.42
1:D:37:GLU:HB2	1:D:60:PHE:CD2	2.53	0.42
1:A:114:PHE:HZ	1:A:155:LEU:HD13	1.84	0.42
1:D:34:ARG:NH1	1:D:37:GLU:OE1	2.51	0.42
1:C:98:GLU:O	1:C:101:GLU:HB3	2.19	0.42
1:C:124:ILE:HG23	1:C:125:SER:N	2.33	0.42
1:D:186:TRP:CD1	1:D:187:ILE:H	2.36	0.42
1:A:107:LEU:HG	1:A:156:VAL:HG23	2.02	0.42
1:D:7:ARG:HH11	1:D:7:ARG:CG	2.33	0.42
1:A:52:GLU:HA	1:A:55:THR:HG22	2.01	0.42
1:B:102:GLU:O	1:B:105:LEU:HG	2.19	0.42
1:D:132:ILE:N	1:D:132:ILE:CD1	2.83	0.42
1:B:8:SER:HA	2:F:8:LEU:CD2	2.49	0.42
1:A:19:THR:CG2	1:A:163:ASP:HB3	2.49	0.42
1:B:41:ILE:O	1:B:42:GLN:HB2	2.20	0.42
1:B:127:LEU:HB2	1:B:130:CYS:HB3	2.01	0.42
1:D:71:ARG:HH11	1:D:71:ARG:HB2	1.85	0.42
2:E:8:LEU:N	2:E:8:LEU:HD22	2.35	0.42
2:E:71:LEU:HD12	2:E:71:LEU:N	2.34	0.42
1:A:86:TYR:CD2	1:A:86:TYR:C	2.97	0.42
1:B:3:THR:HA	1:B:6:ARG:CD	2.50	0.42
1:D:105:LEU:HD12	1:D:106:LEU:N	2.35	0.42
1:D:114:PHE:C	1:D:114:PHE:CD2	2.98	0.42
2:F:39:ASP:OD1	2:F:39:ASP:C	2.62	0.42
1:A:16:ILE:HD11	1:A:89:ILE:CD1	2.50	0.41
1:A:182:PHE:HB2	1:A:185:LEU:HB2	2.03	0.41
1:B:130:CYS:SG	1:B:167:TRP:HB2	2.59	0.41
1:B:175:LEU:HD12	1:B:176:GLU:N	2.35	0.41
1:C:135:GLU:O	1:C:139:ILE:CD1	2.68	0.41
1:D:139:ILE:H	1:D:142:ILE:HG12	1.85	0.41
2:J:67:LEU:N	2:J:67:LEU:HD23	2.34	0.41
1:A:135:GLU:O	1:A:139:ILE:HG13	2.21	0.41
1:C:106:LEU:C	1:C:106:LEU:CD1	2.94	0.41
1:C:124:ILE:CG2	1:C:125:SER:N	2.84	0.41
1:C:138:GLU:CG	1:C:139:ILE:H	2.33	0.41
2:I:43:LEU:HD23	2:I:69:LEU:CD1	2.50	0.41
1:A:178:GLU:O	1:A:181:LYS:HE2	2.20	0.41
1:B:114:PHE:N	1:B:182:PHE:HZ	2.17	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:8:SER:O	1:C:12:PHE:HB2	2.21	0.41
1:C:23:THR:HA	1:C:26:LEU:HD12	2.03	0.41
1:C:52:GLU:HA	1:C:55:THR:HG22	2.02	0.41
1:B:182:PHE:HA	1:B:185:LEU:HB3	2.02	0.41
1:C:97:ILE:HA	1:C:100:LEU:HB2	2.02	0.41
2:F:50:LEU:HD23	2:F:59:TYR:CD1	2.55	0.41
1:B:114:PHE:CD1	1:B:118:ILE:HD12	2.56	0.41
1:C:77:LEU:HD23	1:C:85:LEU:HB3	2.02	0.41
1:C:124:ILE:HD13	1:C:140:LEU:HD21	2.01	0.41
1:A:122:ASP:OD1	1:A:122:ASP:N	2.35	0.41
1:B:92:TRP:O	2:F:42:ARG:NH1	2.54	0.41
1:D:130:CYS:HA	1:D:164:LYS:HE3	2.03	0.41
1:B:23:THR:HG22	1:B:42:GLN:HG2	2.03	0.41
1:D:18:LYS:HB3	1:D:18:LYS:HE2	1.91	0.41
1:B:123:ILE:HD12	1:B:127:LEU:HG	2.03	0.41
2:E:7:THR:CG2	2:E:11:LYS:HB3	2.51	0.41
2:H:61:ILE:HG23	2:H:65:SER:HB2	2.03	0.41
2:J:15:LEU:HD11	2:J:30:ILE:HG12	2.03	0.41
2:J:25:ASN:O	2:J:29:LYS:HG3	2.20	0.41
1:C:36:GLU:C	1:C:38:VAL:H	2.29	0.41
2:I:61:ILE:HG23	2:I:65:SER:HB2	2.02	0.41
1:B:71:ARG:HB3	2:E:64:GLU:O	2.21	0.40
1:C:3:THR:HA	1:C:6:ARG:HG2	2.03	0.40
2:G:50:LEU:HD11	2:G:67:LEU:HD22	2.03	0.40
2:H:67:LEU:N	2:H:67:LEU:HD22	2.36	0.40
2:I:4:PHE:HB2	2:I:66:THR:HG23	2.03	0.40
1:A:2:THR:CG2	1:A:69:TRP:HE1	2.35	0.40
1:B:26:LEU:O	1:B:27:SER:C	2.64	0.40
1:D:120:PRO:HD2	1:D:148:MET:HG3	2.03	0.40
2:E:8:LEU:HD21	2:E:70:VAL:HG22	2.03	0.40
1:C:172:LYS:HA	1:C:175:LEU:HD12	2.04	0.40
1:D:124:ILE:HG23	1:D:125:SER:N	2.36	0.40
1:D:159:LEU:HD23	1:D:167:TRP:CZ3	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	186/203 (92%)	172 (92%)	11 (6%)	3 (2%)	7	35
1	B	186/203 (92%)	171 (92%)	14 (8%)	1 (0%)	24	56
1	C	186/203 (92%)	170 (91%)	13 (7%)	3 (2%)	7	35
1	D	186/203 (92%)	176 (95%)	8 (4%)	2 (1%)	11	42
2	E	70/79 (89%)	69 (99%)	1 (1%)	0	100	100
2	F	71/79 (90%)	69 (97%)	2 (3%)	0	100	100
2	G	69/79 (87%)	68 (99%)	1 (1%)	0	100	100
2	H	73/79 (92%)	71 (97%)	2 (3%)	0	100	100
2	I	69/79 (87%)	65 (94%)	3 (4%)	1 (1%)	9	37
2	J	71/79 (90%)	69 (97%)	2 (3%)	0	100	100
All	All	1167/1286 (91%)	1100 (94%)	57 (5%)	10 (1%)	14	45

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	187	ILE
1	C	66	GLU
1	C	139	ILE
1	D	139	ILE
1	A	180	ASN
1	A	185	LEU
1	A	182	PHE
2	I	9	THR
1	C	138	GLU
1	B	139	ILE

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	162/182 (89%)	135 (83%)	27 (17%)	2	14
1	B	162/182 (89%)	135 (83%)	27 (17%)	2	14
1	C	158/182 (87%)	133 (84%)	25 (16%)	2	15
1	D	166/182 (91%)	140 (84%)	26 (16%)	2	15
2	E	66/69 (96%)	57 (86%)	9 (14%)	3	19
2	F	66/69 (96%)	54 (82%)	12 (18%)	2	11
2	G	64/69 (93%)	55 (86%)	9 (14%)	3	18
2	H	67/69 (97%)	57 (85%)	10 (15%)	3	16
2	I	63/69 (91%)	52 (82%)	11 (18%)	2	12
2	J	67/69 (97%)	55 (82%)	12 (18%)	2	12
All	All	1041/1142 (91%)	873 (84%)	168 (16%)	2	14

All (168) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	3	THR
1	A	10	GLN
1	A	20	LEU
1	A	25	ILE
1	A	27	SER
1	A	42	GLN
1	A	56	LEU
1	A	61	LEU
1	A	64	LEU
1	A	77	LEU
1	A	89	ILE
1	A	95	LYS
1	A	99	LYS
1	A	105	LEU
1	A	119	ILE

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Mol	Chain	Res	Type
1	A	122	ASP
1	A	124	ILE
1	A	127	LEU
1	A	132	ILE
1	A	140	LEU
1	A	141	GLN
1	A	155	LEU
1	A	160	LEU
1	A	181	LYS
1	A	184	GLU
1	A	185	LEU
1	B	1	MET
1	B	3	THR
1	B	6	ARG
1	B	9	LEU
1	B	10	GLN
1	B	13	GLN
1	B	18	LYS
1	B	23	THR
1	B	29	MET
1	B	34	ARG
1	B	55	THR
1	B	58	LEU
1	B	64	LEU
1	B	77	LEU
1	B	86	TYR
1	B	89	ILE
1	B	106	LEU
1	B	107	LEU
1	B	110	LEU
1	B	119	ILE
1	B	122	ASP
1	B	123	ILE
1	B	130	CYS
1	B	141	GLN
1	B	159	LEU
1	B	169	LYS
1	B	187	ILE
1	C	1	MET
1	C	3	THR
1	C	4	GLU
1	C	9	LEU

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Mol	Chain	Res	Type
1	C	13	GLN
1	C	16	ILE
1	C	27	SER
1	C	29	MET
1	C	39	GLN
1	C	46	ASN
1	C	61	LEU
1	C	62	LEU
1	C	102	GLU
1	C	106	LEU
1	C	110	LEU
1	C	122	ASP
1	C	125	SER
1	C	127	LEU
1	C	135	GLU
1	C	159	LEU
1	C	169	LYS
1	C	171	LEU
1	C	175	LEU
1	C	185	LEU
1	C	187	ILE
1	D	1	MET
1	D	3	THR
1	D	4	GLU
1	D	7	ARG
1	D	13	GLN
1	D	17	ARG
1	D	19	THR
1	D	25	ILE
1	D	29	MET
1	D	36	GLU
1	D	55	THR
1	D	58	LEU
1	D	71	ARG
1	D	87	GLU
1	D	89	ILE
1	D	109	ARG
1	D	110	LEU
1	D	132	ILE
1	D	133	ASN
1	D	141	GLN
1	D	159	LEU

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Mol	Chain	Res	Type
1	D	164	LYS
1	D	169	LYS
1	D	185	LEU
1	D	187	ILE
1	D	188	VAL
2	E	1	MET
2	E	6	LYS
2	E	8	LEU
2	E	22	THR
2	E	30	ILE
2	E	32	ASP
2	E	50	LEU
2	E	51	GLU
2	E	66	THR
2	F	3	ILE
2	F	15	LEU
2	F	22	THR
2	F	30	ILE
2	F	33	LYS
2	F	36	ILE
2	F	39	ASP
2	F	48	LYS
2	F	61	ILE
2	F	65	SER
2	F	69	LEU
2	F	73	LEU
2	G	1	MET
2	G	8	LEU
2	G	12	THR
2	G	22	THR
2	G	30	ILE
2	G	58	ASP
2	G	60	ASN
2	G	70	VAL
2	G	71	LEU
2	H	16	GLU
2	H	18	GLU
2	H	22	THR
2	H	24	GLU
2	H	39	ASP
2	H	42	ARG
2	H	43	LEU

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Mol	Chain	Res	Type
2	H	50	LEU
2	H	67	LEU
2	H	74	ARG
2	I	8	LEU
2	I	13	ILE
2	I	22	THR
2	I	23	ILE
2	I	30	ILE
2	I	39	ASP
2	I	44	ILE
2	I	52	ASP
2	I	67	LEU
2	I	68	HIS
2	I	71	LEU
2	J	15	LEU
2	J	20	SER
2	J	22	THR
2	J	30	ILE
2	J	39	ASP
2	J	42	ARG
2	J	44	ILE
2	J	51	GLU
2	J	52	ASP
2	J	63	LYS
2	J	67	LEU
2	J	73	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	46	ASN
1	B	5	GLN
1	B	111	GLN
2	E	40	GLN
2	E	68	HIS
2	F	25	ASN
2	G	41	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	188/203 (92%)	-0.22	1 (0%) 87 69	29, 66, 118, 165	0
1	B	188/203 (92%)	-0.12	2 (1%) 78 53	40, 80, 127, 166	0
1	C	188/203 (92%)	-0.21	0 100 100	46, 90, 141, 167	0
1	D	188/203 (92%)	-0.31	0 100 100	36, 69, 112, 126	0
2	E	72/79 (91%)	-0.34	0 100 100	44, 76, 117, 141	0
2	F	73/79 (92%)	-0.25	1 (1%) 73 48	44, 148, 194, 214	1 (1%)
2	G	71/79 (89%)	-0.18	0 100 100	109, 174, 214, 236	0
2	H	75/79 (94%)	-0.32	1 (1%) 75 50	38, 131, 170, 197	1 (1%)
2	I	71/79 (89%)	-0.05	3 (4%) 40 25	68, 102, 152, 166	0
2	J	73/79 (92%)	-0.21	1 (1%) 73 48	25, 74, 110, 141	1 (1%)
All	All	1187/1286 (92%)	-0.22	9 (0%) 82 60	25, 85, 172, 236	3 (0%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	J	73	LEU	6.2
2	F	73	LEU	3.5
2	H	73	LEU	2.7
2	I	15	LEU	2.5
1	A	130	CYS	2.5
1	B	124	ILE	2.4
2	I	5	VAL	2.3
2	I	69	LEU	2.2
1	B	120	PRO	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.