



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 07:29 PM UTC

PDB ID : 4DDI / pdb_00004ddi
Title : Crystal structure of human OTUB1/UbcH5b Ub/Ub
Authors : Juang, Y.C.; Sanches, M.; Sicheri, F.
Deposited on : 2012-01-18
Resolution : 3.80 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

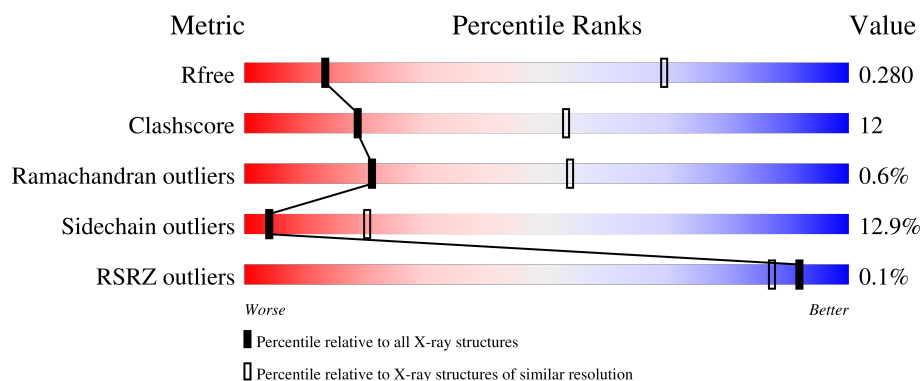
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.80 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	1065 (3.96-3.64)
Clashscore	190562	1012 (3.94-3.66)
Ramachandran outliers	187476	1048 (3.96-3.64)
Sidechain outliers	187428	1043 (3.96-3.64)
RSRZ outliers	180081	1064 (3.96-3.64)

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	399	 67% 28% 5%
1	B	399	 66% 30% .
1	C	399	 67% 29% .
2	D	76	 54% 36% 11%
2	E	76	 57% 33% 11%

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Mol	Chain	Length	Quality of chain
2	F	76	57% 33% 11%
2	G	76	66% 25% 9%
2	H	76	62% 29% 9%
2	I	76	64% 26% 9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 13284 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 Ubiquitin-conjugating enzyme E2 D2, Ubiquitin thioesterase OTUB1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	399	Total	C	N	O	S	0	0	0
			3226	2055	542	616	13			
1	B	399	Total	C	N	O	S	0	0	0
			3226	2055	542	616	13			
1	C	399	Total	C	N	O	S	0	0	0
			3226	2055	542	616	13			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP P62837
A	0	ALA	-	expression tag	UNP P62837
A	85	SER	CYS	conflict	UNP P62837
A	148	GLY	-	linker	UNP Q96FW1
A	149	GLY	-	linker	UNP Q96FW1
A	150	SER	-	linker	UNP Q96FW1
A	1091	SER	CYS	conflict	UNP Q96FW1
B	-1	GLY	-	expression tag	UNP P62837
B	0	ALA	-	expression tag	UNP P62837
B	85	SER	CYS	conflict	UNP P62837
B	148	GLY	-	linker	UNP Q96FW1
B	149	GLY	-	linker	UNP Q96FW1
B	150	SER	-	linker	UNP Q96FW1
B	1091	SER	CYS	conflict	UNP Q96FW1
C	-1	GLY	-	expression tag	UNP P62837
C	0	ALA	-	expression tag	UNP P62837
C	85	SER	CYS	conflict	UNP P62837
C	148	GLY	-	linker	UNP Q96FW1
C	149	GLY	-	linker	UNP Q96FW1
C	150	SER	-	linker	UNP Q96FW1
C	1091	SER	CYS	conflict	UNP Q96FW1

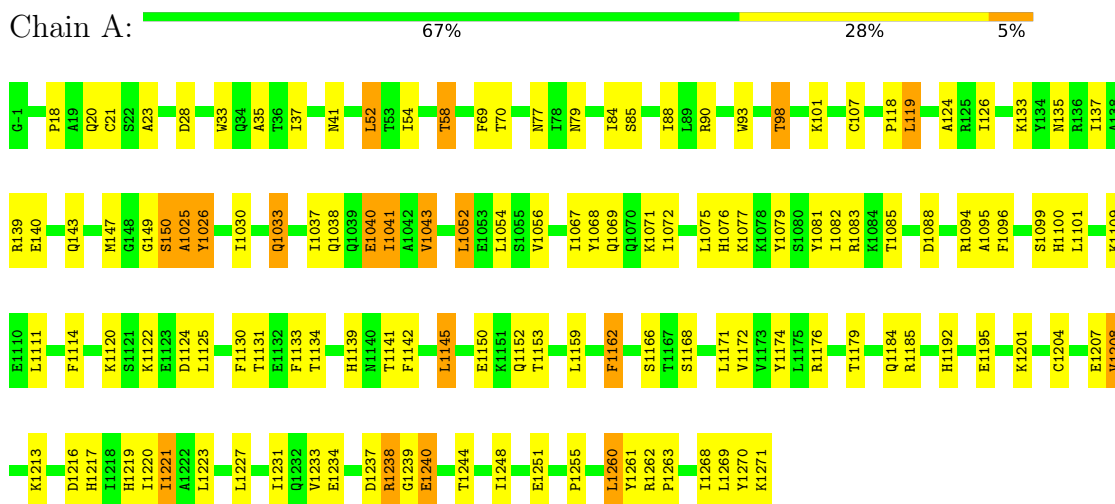
- Molecule 2 is a protein called Polyubiquitin-C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	76	Total 601	C 378	N 105	O 117	S 1	0	0	0
2	E	76	Total 601	C 378	N 105	O 117	S 1	0	0	0
2	F	76	Total 601	C 378	N 105	O 117	S 1	0	0	0
2	G	76	Total 601	C 378	N 105	O 117	S 1	0	0	0
2	H	76	Total 601	C 378	N 105	O 117	S 1	0	0	0
2	I	76	Total 601	C 378	N 105	O 117	S 1	0	0	0

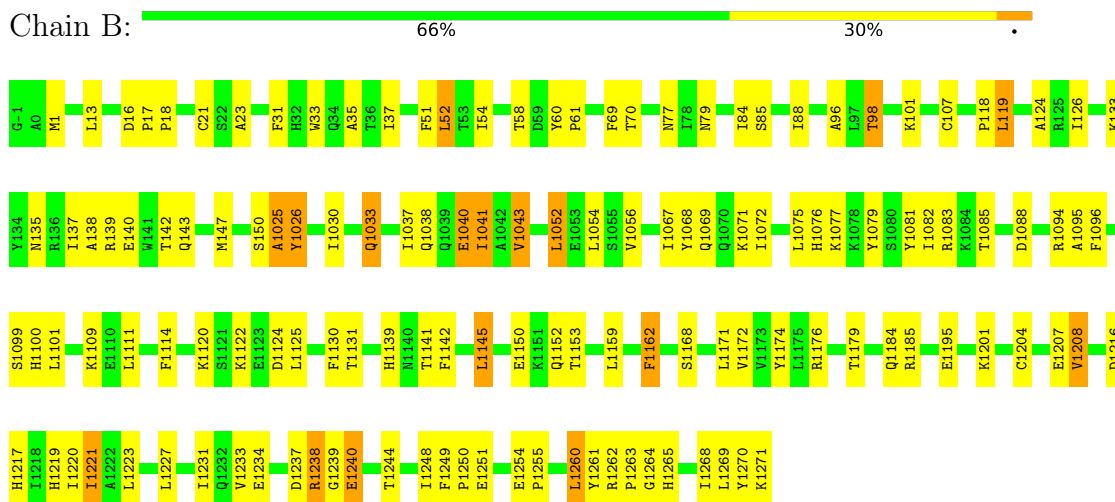
3 Residue-property plots

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: Ubiquitin-conjugating enzyme E2 D2, Ubiquitin thioesterase OTUB1

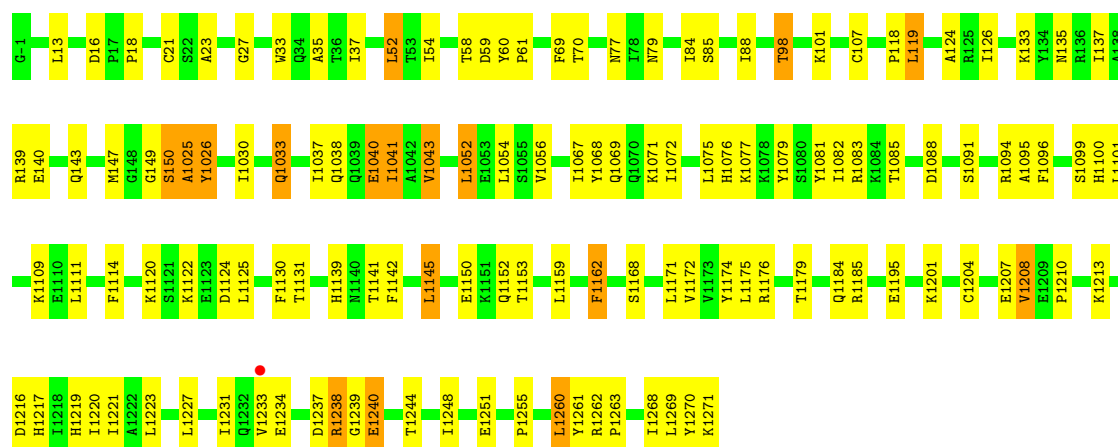


- Molecule 1: Ubiquitin-conjugating enzyme E2 D2, Ubiquitin thioesterase OTUB1

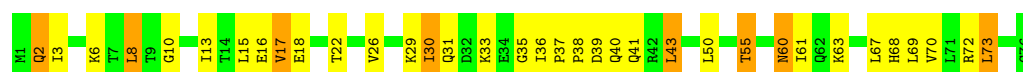


- Molecule 1: Ubiquitin-conjugating enzyme E2 D2, Ubiquitin thioesterase OTUB1





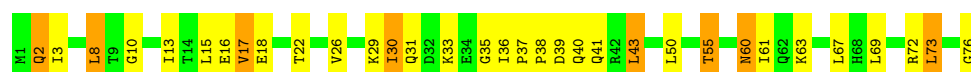
• Molecule 2: Polyubiquitin-C



• Molecule 2: Polyubiquitin-C



• Molecule 2: Polyubiquitin-C



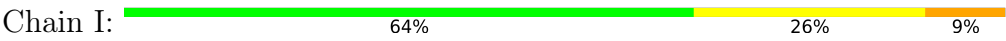
• Molecule 2: Polyubiquitin-C



• Molecule 2: Polyubiquitin-C



• Molecule 2: Polyubiquitin-C



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	173.74Å 106.22Å 134.69Å 90.00° 123.74° 90.00°	Depositor
Resolution (Å)	49.47 – 3.80 49.47 – 3.80	Depositor EDS
% Data completeness (in resolution range)	97.6 (49.47-3.80) 97.6 (49.47-3.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 3.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.1_743)	Depositor
R, R_{free}	0.218 , 0.273 0.234 , 0.280	Depositor DCC
R_{free} test set	987 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	126.9	Xtriage
Anisotropy	0.437	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 222.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13284	wwPDB-VP
Average B, all atoms (Å ²)	168.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.28% 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	0.33	0/3306	0.82	1/4476 (0.0%)
1	B	0.33	0/3306	0.82	1/4476 (0.0%)
1	C	0.33	0/3306	0.82	1/4476 (0.0%)
2	D	0.42	0/607	1.06	3/816 (0.4%)
2	E	0.43	0/607	1.05	3/816 (0.4%)
2	F	0.43	0/607	1.05	3/816 (0.4%)
2	G	0.34	0/607	0.90	1/816 (0.1%)
2	H	0.34	0/607	0.90	1/816 (0.1%)
2	I	0.34	0/607	0.90	1/816 (0.1%)
All	All	0.35	0/13560	0.86	15/18324 (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
2	D	0	1
2	E	0	1
2	F	0	1
All	All	0	3

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	36	ILE	CA-C-N	6.18	124.11	119.66
2	D	36	ILE	C-N-CA	6.18	124.11	119.66
2	E	36	ILE	CA-C-N	6.14	124.08	119.66
2	E	36	ILE	C-N-CA	6.14	124.08	119.66
2	F	36	ILE	CA-C-N	6.05	124.02	119.66
2	F	36	ILE	C-N-CA	6.05	124.02	119.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	60	ASN	N-CA-C	5.51	122.10	113.72
2	G	73	LEU	N-CA-C	-5.36	101.23	109.76
2	I	73	LEU	N-CA-C	-5.33	101.28	109.76
2	F	60	ASN	N-CA-C	5.33	122.16	110.80
2	H	73	LEU	N-CA-C	-5.33	101.28	109.76
2	E	60	ASN	N-CA-C	5.30	122.08	110.80
1	C	1207	GLU	N-CA-C	5.13	119.69	113.38
1	B	1207	GLU	N-CA-C	5.08	119.62	113.38
1	A	1207	GLU	N-CA-C	5.07	119.61	113.38

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	60	ASN	Peptide
2	E	60	ASN	Peptide
2	F	60	ASN	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	3226	0	3141	79	0
1	B	3226	0	3141	76	0
1	C	3226	0	3141	75	0
2	D	601	0	629	19	0
2	E	601	0	629	20	0
2	F	601	0	629	21	0
2	G	601	0	629	16	0
2	H	601	0	629	18	0
2	I	601	0	629	21	0
All	All	13284	0	13197	315	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 (315) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1220:ILE:HD12	2:D:73:LEU:HD23	1.64	0.79
1:B:1111:LEU:HD21	1:B:1150:GLU:HA	1.66	0.78
1:C:1111:LEU:HD21	1:C:1150:GLU:HA	1.66	0.77
1:C:1204:CYS:HA	1:C:1208:VAL:HG13	1.67	0.77
2:F:31:GLN:HG3	2:F:37:PRO:HA	1.67	0.76
2:E:31:GLN:HG3	2:E:37:PRO:HA	1.68	0.76
2:D:31:GLN:HG3	2:D:37:PRO:HA	1.68	0.76
1:A:1111:LEU:HD21	1:A:1150:GLU:HA	1.66	0.75
1:B:1204:CYS:HA	1:B:1208:VAL:HG13	1.67	0.75
1:C:150:SER:O	1:C:1026:TYR:N	2.19	0.75
1:A:1204:CYS:HA	1:A:1208:VAL:HG13	1.67	0.74
1:B:1081:TYR:HB2	1:B:1270:TYR:HB2	1.71	0.72
1:A:1081:TYR:HB2	1:A:1270:TYR:HB2	1.71	0.72
1:C:1038:GLN:HA	1:C:1041:ILE:HB	1.73	0.71
1:C:1081:TYR:HB2	1:C:1270:TYR:HB2	1.71	0.71
2:G:5:VAL:HG13	2:G:13:ILE:HG23	1.73	0.70
2:I:5:VAL:HG13	2:I:13:ILE:HG23	1.74	0.70
1:A:1038:GLN:HA	1:A:1041:ILE:HB	1.73	0.69
1:B:1038:GLN:HA	1:B:1041:ILE:HB	1.73	0.69
1:C:150:SER:O	1:C:1025:ALA:C	2.35	0.68
2:H:5:VAL:HG13	2:H:13:ILE:HG23	1.73	0.68
1:A:1145:LEU:HD13	1:A:1171:LEU:HD12	1.78	0.66
1:C:149:GLY:O	1:C:150:SER:C	2.36	0.66
1:C:1145:LEU:HD13	1:C:1171:LEU:HD12	1.78	0.66
2:E:39:ASP:O	2:E:72:ARG:NH2	2.28	0.66
2:F:39:ASP:O	2:F:72:ARG:NH2	2.28	0.65
1:A:149:GLY:O	1:A:150:SER:C	2.39	0.65
1:B:1145:LEU:HD13	1:B:1171:LEU:HD12	1.78	0.65
2:D:39:ASP:O	2:D:72:ARG:NH2	2.28	0.64
1:A:1094:ARG:NH1	1:A:1168:SER:OG	2.31	0.64
1:C:1094:ARG:NH1	1:C:1168:SER:OG	2.31	0.64
1:A:143:GLN:HA	1:A:147:MET:HB2	1.80	0.64
1:B:1094:ARG:NH1	1:B:1168:SER:OG	2.31	0.64
1:A:1133:PHE:HZ	1:B:98:THR:HB	1.63	0.63
1:B:143:GLN:HA	1:B:147:MET:HB2	1.80	0.63
1:C:143:GLN:HA	1:C:147:MET:HB2	1.80	0.63
1:B:1265:HIS:HB2	2:H:48:LYS:HE3	1.80	0.62
2:H:23:ILE:HD13	2:H:50:LEU:HB3	1.81	0.62
2:I:23:ILE:HD13	2:I:50:LEU:HB3	1.81	0.62
1:C:52:LEU:HB3	1:C:69:PHE:HA	1.82	0.61
1:B:79:ASN:HB3	1:B:119:LEU:HD22	1.83	0.61
1:C:79:ASN:HB3	1:C:119:LEU:HD22	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:G:23:ILE:HD13	2:G:50:LEU:HB3	1.81	0.61
1:B:52:LEU:HB3	1:B:69:PHE:HA	1.82	0.61
1:A:79:ASN:HB3	1:A:119:LEU:HD22	1.83	0.60
2:E:30:ILE:HB	2:E:41:GLN:HE22	1.67	0.60
1:A:52:LEU:HB3	1:A:69:PHE:HA	1.82	0.59
1:C:1220:ILE:HD12	2:F:73:LEU:HD23	1.84	0.59
1:A:28:ASP:HB2	1:B:31:PHE:CD1	2.38	0.59
1:A:98:THR:HG23	1:A:101:LYS:HG2	1.85	0.59
2:F:30:ILE:HB	2:F:41:GLN:HE22	1.67	0.59
1:C:1071:LYS:NZ	1:C:1239:GLY:O	2.36	0.59
1:A:150:SER:O	1:A:1025:ALA:C	2.46	0.59
1:B:98:THR:HG23	1:B:101:LYS:HG2	1.84	0.58
1:C:98:THR:HG23	1:C:101:LYS:HG2	1.85	0.58
1:A:1260:LEU:HD21	1:A:1262:ARG:HG3	1.86	0.58
1:B:1071:LYS:NZ	1:B:1239:GLY:O	2.36	0.58
1:A:58:THR:HG23	1:C:27:GLY:HA2	1.86	0.58
2:D:30:ILE:HB	2:D:41:GLN:HE22	1.67	0.57
1:A:28:ASP:HB2	1:B:31:PHE:HD1	1.70	0.57
2:E:61:ILE:O	2:E:61:ILE:HG22	2.04	0.57
2:F:61:ILE:HG22	2:F:61:ILE:O	2.05	0.57
1:B:1083:ARG:HB3	1:B:1268:ILE:HG13	1.87	0.56
1:C:1083:ARG:HB3	1:C:1268:ILE:HG13	1.87	0.56
2:I:7:THR:OG1	2:I:8:LEU:N	2.39	0.56
2:E:22:THR:HA	2:E:55:THR:HA	1.87	0.56
1:C:133:LYS:O	1:C:137:ILE:HG12	2.05	0.56
1:C:1260:LEU:HD21	1:C:1262:ARG:HG3	1.86	0.56
1:B:1260:LEU:HD21	1:B:1262:ARG:HG3	1.86	0.56
1:A:133:LYS:O	1:A:137:ILE:HG12	2.05	0.56
1:B:1120:LYS:NZ	1:B:1124:ASP:OD1	2.33	0.56
1:A:1083:ARG:HB3	1:A:1268:ILE:HG13	1.88	0.55
1:B:133:LYS:O	1:B:137:ILE:HG12	2.05	0.55
1:B:1220:ILE:HD12	2:E:73:LEU:HD23	1.88	0.55
1:C:1082:ILE:HG13	1:C:1269:LEU:HD23	1.87	0.55
1:A:1082:ILE:HG13	1:A:1269:LEU:HD23	1.87	0.55
1:B:1082:ILE:HG13	1:B:1269:LEU:HD23	1.87	0.55
2:D:22:THR:HA	2:D:55:THR:HA	1.87	0.55
2:F:22:THR:HA	2:F:55:THR:HA	1.88	0.55
1:A:1125:LEU:HB3	1:A:1130:PHE:HD2	1.72	0.55
1:A:1166:SER:OG	1:B:1:MET:HE3	2.07	0.55
2:E:2:GLN:HB2	2:E:16:GLU:HA	1.89	0.55
1:A:79:ASN:H	1:A:119:LEU:HD13	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:ASN:H	1:C:119:LEU:HD13	1.72	0.55
1:A:1071:LYS:NZ	1:A:1239:GLY:O	2.36	0.54
2:F:2:GLN:HB2	2:F:16:GLU:HA	1.89	0.54
2:D:2:GLN:HB2	2:D:16:GLU:HA	1.89	0.54
1:B:79:ASN:H	1:B:119:LEU:HD13	1.72	0.54
1:B:1125:LEU:HB3	1:B:1130:PHE:HD2	1.72	0.54
1:C:1184:GLN:HE22	1:C:1201:LYS:HG3	1.73	0.54
1:C:1125:LEU:HB3	1:C:1130:PHE:HD2	1.72	0.54
1:A:150:SER:O	1:A:1026:TYR:N	2.41	0.54
2:G:1:MET:H3	2:G:63:LYS:HG2	1.72	0.54
1:C:1237:ASP:CG	1:C:1238:ARG:H	2.16	0.53
1:B:1179:THR:HG22	1:B:1223:LEU:HB2	1.90	0.53
1:C:1026:TYR:O	1:C:1030:ILE:HG13	2.09	0.53
1:B:1026:TYR:O	1:B:1030:ILE:HG13	2.09	0.53
2:E:17:VAL:HG22	2:E:18:GLU:H	1.74	0.53
1:A:1237:ASP:CG	1:A:1238:ARG:H	2.16	0.53
1:B:1237:ASP:CG	1:B:1238:ARG:H	2.16	0.53
1:C:1179:THR:HG22	1:C:1223:LEU:HB2	1.90	0.53
1:A:41:ASN:ND2	2:G:11:LYS:HD3	2.23	0.53
1:B:135:ASN:HB3	1:B:139:ARG:HH22	1.74	0.53
1:C:135:ASN:HB3	1:C:139:ARG:HH22	1.74	0.53
2:G:7:THR:OG1	2:G:8:LEU:N	2.38	0.53
1:A:1184:GLN:HE22	1:A:1201:LYS:HG3	1.73	0.53
2:D:17:VAL:HG22	2:D:18:GLU:H	1.74	0.52
1:A:1221:ILE:HG13	2:D:8:LEU:HD13	1.91	0.52
2:E:17:VAL:HG23	2:E:29:LYS:HE3	1.92	0.52
1:B:1052:LEU:HD12	1:B:1056:VAL:HG21	1.91	0.52
2:F:17:VAL:HG22	2:F:18:GLU:H	1.74	0.52
1:A:135:ASN:HB3	1:A:139:ARG:HH22	1.74	0.52
1:A:1026:TYR:O	1:A:1030:ILE:HG13	2.09	0.52
1:A:1179:THR:HG22	1:A:1223:LEU:HB2	1.90	0.52
1:B:1184:GLN:HE22	1:B:1201:LYS:HG3	1.73	0.52
2:I:4:PHE:HE1	2:I:64:GLU:HA	1.74	0.52
1:C:1248:ILE:HG12	1:C:1255:PRO:HG3	1.91	0.52
2:H:7:THR:OG1	2:H:8:LEU:N	2.38	0.52
1:A:1133:PHE:CZ	1:B:98:THR:HB	2.42	0.52
1:A:1248:ILE:HG12	1:A:1255:PRO:HG3	1.91	0.52
1:C:1052:LEU:HD12	1:C:1056:VAL:HG21	1.92	0.52
2:H:4:PHE:HE1	2:H:64:GLU:HA	1.75	0.51
1:C:1120:LYS:NZ	1:C:1124:ASP:OD1	2.33	0.51
2:G:4:PHE:HE1	2:G:64:GLU:HA	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:31:GLN:HG3	2:F:38:PRO:HD3	1.93	0.51
1:A:33:TRP:HB2	1:A:54:ILE:HB	1.93	0.51
1:B:1217:HIS:CE1	2:E:8:LEU:HD11	2.46	0.51
1:A:1120:LYS:NZ	1:A:1124:ASP:OD1	2.33	0.51
1:B:1248:ILE:HG12	1:B:1255:PRO:HG3	1.91	0.51
2:H:34:GLU:HB3	2:H:36:ILE:HG13	1.93	0.51
1:B:1264:GLY:HA2	2:E:75:GLY:HA2	1.93	0.50
1:C:33:TRP:HB2	1:C:54:ILE:HB	1.92	0.50
1:A:1052:LEU:HD12	1:A:1056:VAL:HG21	1.92	0.50
1:B:33:TRP:HB2	1:B:54:ILE:HB	1.92	0.50
2:D:17:VAL:HG23	2:D:29:LYS:HE3	1.92	0.50
1:A:1134:THR:HG21	1:B:96:ALA:O	2.10	0.50
1:C:1095:ALA:HB1	1:C:1268:ILE:HG23	1.94	0.50
2:G:1:MET:N	2:G:63:LYS:HG2	2.27	0.50
1:A:90:ARG:NH2	2:I:75:GLY:HA3	2.26	0.50
2:H:1:MET:N	2:H:63:LYS:HG2	2.26	0.50
1:A:1139:HIS:HA	1:A:1174:TYR:CE2	2.47	0.50
1:C:1122:LYS:HB2	1:C:1139:HIS:CE1	2.47	0.50
2:I:22:THR:HB	2:I:25:ASN:H	1.77	0.50
1:C:1139:HIS:HA	1:C:1174:TYR:CE2	2.47	0.50
2:I:1:MET:N	2:I:63:LYS:HG2	2.27	0.49
1:A:1122:LYS:HB2	1:A:1139:HIS:CE1	2.48	0.49
2:E:31:GLN:HG3	2:E:38:PRO:HD3	1.93	0.49
2:F:17:VAL:HG23	2:F:29:LYS:HE3	1.92	0.49
2:I:34:GLU:HB3	2:I:36:ILE:HG13	1.93	0.49
2:D:31:GLN:HG3	2:D:38:PRO:HD3	1.93	0.49
1:B:1221:ILE:HG13	2:E:8:LEU:HD13	1.94	0.49
2:G:34:GLU:HB3	2:G:36:ILE:HG13	1.93	0.49
1:A:1095:ALA:HB1	1:A:1268:ILE:HG23	1.94	0.49
2:H:22:THR:HB	2:H:25:ASN:H	1.77	0.49
2:G:22:THR:HB	2:G:25:ASN:H	1.77	0.49
1:A:118:PRO:HB3	1:A:124:ALA:HB2	1.94	0.48
1:B:1139:HIS:HA	1:B:1174:TYR:CE2	2.47	0.48
1:A:90:ARG:HH22	2:I:75:GLY:HA3	1.77	0.48
1:B:1095:ALA:HB1	1:B:1268:ILE:HG23	1.94	0.48
1:A:1260:LEU:HB2	1:A:1269:LEU:HD11	1.95	0.48
1:B:118:PRO:HB3	1:B:124:ALA:HB2	1.94	0.48
1:B:1122:LYS:HB2	1:B:1139:HIS:CE1	2.47	0.48
2:I:61:ILE:H	2:I:61:ILE:HG12	1.29	0.48
1:C:135:ASN:HB3	1:C:139:ARG:NH2	2.29	0.48
1:C:1030:ILE:HG12	2:I:8:LEU:HD21	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1217:HIS:HB3	2:D:70:VAL:HG13	1.96	0.48
1:B:135:ASN:HB3	1:B:139:ARG:NH2	2.29	0.48
2:I:1:MET:H1	2:I:63:LYS:HG2	1.78	0.48
1:C:118:PRO:HB3	1:C:124:ALA:HB2	1.94	0.48
1:B:1088:ASP:OD2	1:B:1094:ARG:NH2	2.47	0.48
1:C:1261:TYR:CE1	1:C:1263:PRO:HA	2.49	0.47
1:A:135:ASN:HB3	1:A:139:ARG:NH2	2.29	0.47
1:A:1261:TYR:CE1	1:A:1263:PRO:HA	2.50	0.47
1:C:1088:ASP:OD2	1:C:1094:ARG:NH2	2.47	0.47
1:C:1260:LEU:HB2	1:C:1269:LEU:HD11	1.95	0.47
1:B:1261:TYR:CE1	1:B:1263:PRO:HA	2.50	0.47
2:F:43:LEU:HD13	2:F:50:LEU:HD12	1.97	0.47
1:A:1071:LYS:HD2	1:A:1244:THR:HG22	1.97	0.47
1:B:1260:LEU:HB2	1:B:1269:LEU:HD11	1.95	0.47
1:A:1033:GLN:HG3	2:G:8:LEU:HD13	1.97	0.47
2:I:3:ILE:HG22	2:I:65:SER:H	1.80	0.47
1:A:1088:ASP:OD2	1:A:1094:ARG:NH2	2.47	0.46
2:D:43:LEU:HD13	2:D:50:LEU:HD12	1.97	0.46
1:B:1040:GLU:O	1:B:1043:VAL:HG12	2.15	0.46
1:A:20:GLN:NE2	2:G:9:THR:O	2.48	0.46
1:B:1114:PHE:CE1	1:B:1227:LEU:HD23	2.50	0.46
1:C:126:ILE:HD12	1:C:133:LYS:HD2	1.97	0.46
1:C:1040:GLU:O	1:C:1043:VAL:HG12	2.16	0.46
1:B:23:ALA:HB2	1:B:35:ALA:HB2	1.98	0.46
1:C:1114:PHE:CE1	1:C:1227:LEU:HD23	2.50	0.46
1:A:1114:PHE:CE1	1:A:1227:LEU:HD23	2.50	0.46
1:B:1071:LYS:HD2	1:B:1244:THR:HG22	1.97	0.46
1:C:1071:LYS:HD2	1:C:1244:THR:HG22	1.97	0.46
1:A:126:ILE:HD12	1:A:133:LYS:HD2	1.97	0.46
1:A:1040:GLU:O	1:A:1043:VAL:HG12	2.16	0.46
2:G:3:ILE:HG22	2:G:65:SER:H	1.81	0.46
2:G:7:THR:HG22	2:G:11:LYS:O	2.16	0.46
1:B:126:ILE:HD12	1:B:133:LYS:HD2	1.97	0.45
1:B:1237:ASP:OD1	1:B:1237:ASP:N	2.50	0.45
2:G:61:ILE:H	2:G:61:ILE:HG12	1.29	0.45
2:H:7:THR:HG22	2:H:11:LYS:O	2.16	0.45
2:I:7:THR:HG22	2:I:11:LYS:O	2.16	0.45
1:A:77:ASN:HB3	1:A:85:SER:HB3	1.98	0.45
1:C:1091:SER:HB2	2:F:76:GLY:HA2	1.97	0.45
2:E:43:LEU:HD13	2:E:50:LEU:HD12	1.97	0.45
2:H:3:ILE:HG22	2:H:65:SER:H	1.80	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:59:ASP:OD1	1:C:59:ASP:N	2.47	0.45
1:A:1237:ASP:OD1	1:A:1237:ASP:N	2.50	0.45
1:C:77:ASN:HB3	1:C:85:SER:HB3	1.99	0.45
1:C:1033:GLN:HG3	2:I:8:LEU:HD13	1.98	0.45
1:A:1068:TYR:O	1:A:1072:ILE:HG13	2.17	0.45
1:B:77:ASN:HB3	1:B:85:SER:HB3	1.99	0.45
1:B:1068:TYR:O	1:B:1072:ILE:HG13	2.17	0.44
1:B:1101:LEU:HD23	1:B:1101:LEU:HA	1.86	0.44
2:H:37:PRO:HA	2:H:38:PRO:HD3	1.91	0.44
1:C:23:ALA:HB2	1:C:35:ALA:HB2	1.98	0.44
2:F:26:VAL:O	2:F:30:ILE:HG13	2.17	0.44
1:C:1025:ALA:O	1:C:1026:TYR:C	2.61	0.44
1:C:1068:TYR:O	1:C:1072:ILE:HG13	2.17	0.44
1:C:1240:GLU:CD	1:C:1240:GLU:H	2.26	0.44
1:C:37:ILE:HD12	1:C:52:LEU:HD21	2.00	0.44
1:C:1216:ASP:H	1:C:1219:HIS:CD2	2.36	0.44
1:A:23:ALA:HB2	1:A:35:ALA:HB2	1.98	0.44
1:B:37:ILE:HD12	1:B:52:LEU:HD21	2.00	0.44
1:A:1033:GLN:O	1:A:1037:ILE:HG13	2.18	0.43
1:C:1091:SER:CB	2:F:76:GLY:HA2	2.48	0.43
2:D:13:ILE:HD12	2:D:33:LYS:HE3	2.00	0.43
2:F:13:ILE:HD12	2:F:33:LYS:HE3	2.00	0.43
1:A:1216:ASP:H	1:A:1219:HIS:CD2	2.36	0.43
2:H:5:VAL:HA	2:H:67:LEU:O	2.19	0.43
2:E:26:VAL:O	2:E:30:ILE:HG13	2.18	0.43
2:G:5:VAL:HA	2:G:67:LEU:O	2.19	0.43
1:C:1213:LYS:HD3	1:C:1213:LYS:HA	1.84	0.43
1:A:1240:GLU:CD	1:A:1240:GLU:H	2.26	0.43
1:B:1216:ASP:H	1:B:1219:HIS:CD2	2.36	0.43
1:B:1033:GLN:O	1:B:1037:ILE:HG13	2.18	0.43
1:B:1176:ARG:HD3	1:B:1208:VAL:O	2.19	0.43
2:I:5:VAL:HA	2:I:67:LEU:O	2.19	0.43
1:B:17:PRO:HA	1:B:18:PRO:HD3	1.82	0.43
1:A:1168:SER:O	1:A:1172:VAL:HG23	2.19	0.43
1:B:138:ALA:O	1:B:142:THR:OG1	2.32	0.43
2:D:26:VAL:O	2:D:30:ILE:HG13	2.18	0.43
1:C:1168:SER:O	1:C:1172:VAL:HG23	2.19	0.43
1:B:1033:GLN:HG3	2:H:8:LEU:HD13	1.99	0.42
1:B:1168:SER:O	1:B:1172:VAL:HG23	2.19	0.42
1:C:149:GLY:O	1:C:150:SER:O	2.37	0.42
1:C:1033:GLN:O	1:C:1037:ILE:HG13	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:37:PRO:HG2	2:F:40:GLN:HG3	2.01	0.42
1:A:1176:ARG:HD3	1:A:1208:VAL:O	2.19	0.42
2:E:13:ILE:HD12	2:E:33:LYS:HE3	2.00	0.42
1:A:1096:PHE:CE2	1:A:1223:LEU:HD22	2.55	0.42
1:C:1096:PHE:CE2	1:C:1223:LEU:HD22	2.55	0.42
2:H:61:ILE:H	2:H:61:ILE:HG12	1.29	0.42
1:A:37:ILE:HD12	1:A:52:LEU:HD21	2.00	0.42
1:B:1083:ARG:HD2	1:B:1268:ILE:HD11	2.02	0.42
2:D:37:PRO:HG2	2:D:40:GLN:HG3	2.01	0.42
2:E:37:PRO:HG2	2:E:40:GLN:HG3	2.01	0.42
1:A:1192:HIS:HD2	2:D:68:HIS:NE2	2.16	0.42
1:A:1213:LYS:HD3	1:A:1213:LYS:HA	1.83	0.42
1:B:1240:GLU:H	1:B:1240:GLU:CD	2.26	0.42
1:C:1079:TYR:CE2	1:C:1234:GLU:HG3	2.55	0.42
2:I:4:PHE:CE1	2:I:64:GLU:HA	2.54	0.42
1:A:1079:TYR:CE2	1:A:1234:GLU:HG3	2.55	0.42
1:B:1096:PHE:CE2	1:B:1223:LEU:HD22	2.55	0.42
1:C:1083:ARG:HD2	1:C:1268:ILE:HD11	2.02	0.42
1:A:1054:LEU:HD12	1:A:1075:LEU:HD23	2.02	0.42
2:D:6:LYS:HB3	2:D:6:LYS:HE2	1.88	0.42
2:G:4:PHE:CE1	2:G:64:GLU:HA	2.54	0.42
1:B:51:PHE:CE2	1:B:1025:ALA:HB1	2.56	0.41
1:C:1175:LEU:O	1:C:1179:THR:HG23	2.20	0.41
2:D:31:GLN:O	2:D:35:GLY:HA2	2.21	0.41
2:E:31:GLN:O	2:E:35:GLY:HA2	2.21	0.41
1:A:1101:LEU:HD23	1:A:1101:LEU:HA	1.86	0.41
1:B:1079:TYR:CE2	1:B:1234:GLU:HG3	2.55	0.41
1:C:60:TYR:CD1	1:C:61:PRO:HA	2.56	0.41
1:C:1079:TYR:CE1	1:C:1271:LYS:HE3	2.55	0.41
1:C:1101:LEU:HD23	1:C:1101:LEU:HA	1.86	0.41
1:C:1217:HIS:CE1	2:F:8:LEU:HD11	2.56	0.41
2:F:8:LEU:C	2:F:10:GLY:H	2.28	0.41
1:B:1142:PHE:HD2	1:B:1174:TYR:CD2	2.38	0.41
1:C:1030:ILE:HG12	2:I:8:LEU:CD2	2.50	0.41
1:A:1083:ARG:HD2	1:A:1268:ILE:HD11	2.02	0.41
1:A:1142:PHE:HD2	1:A:1174:TYR:CD2	2.39	0.41
1:A:1162:PHE:O	1:A:1168:SER:HB2	2.21	0.41
1:B:1079:TYR:CE1	1:B:1271:LYS:HE3	2.55	0.41
1:C:1054:LEU:HD12	1:C:1075:LEU:HD23	2.02	0.41
2:F:31:GLN:O	2:F:35:GLY:HA2	2.20	0.41
2:I:23:ILE:CD1	2:I:50:LEU:HB3	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:1099:SER:HB2	1:C:1268:ILE:HG21	2.03	0.41
2:D:8:LEU:C	2:D:10:GLY:H	2.28	0.41
2:H:4:PHE:CD2	2:H:14:THR:HG22	2.56	0.41
1:A:1099:SER:HB2	1:A:1268:ILE:HG21	2.03	0.41
1:B:1054:LEU:HD12	1:B:1075:LEU:HD23	2.01	0.41
1:B:1254:GLU:HA	1:B:1255:PRO:HD3	1.89	0.41
1:C:18:PRO:HB2	1:C:21:CYS:O	2.21	0.41
1:C:1176:ARG:HD3	1:C:1208:VAL:O	2.19	0.41
2:E:8:LEU:C	2:E:10:GLY:H	2.28	0.41
1:B:60:TYR:CD1	1:B:61:PRO:HA	2.56	0.41
1:C:1142:PHE:HD2	1:C:1174:TYR:CD2	2.38	0.41
2:E:13:ILE:HD13	2:E:13:ILE:HA	1.96	0.41
1:B:1249:PHE:HA	1:B:1250:PRO:HA	1.88	0.40
2:F:37:PRO:HA	2:F:38:PRO:HD3	1.89	0.40
1:A:18:PRO:HB2	1:A:21:CYS:O	2.21	0.40
2:H:1:MET:HB2	2:H:63:LYS:HE2	2.04	0.40
2:H:4:PHE:CE1	2:H:64:GLU:HA	2.54	0.40
1:A:1079:TYR:CE1	1:A:1271:LYS:HE3	2.55	0.40
1:B:18:PRO:HB2	1:B:21:CYS:O	2.21	0.40
1:B:1162:PHE:O	1:B:1168:SER:HB2	2.21	0.40
1:A:93:TRP:O	1:C:1210:PRO:HB3	2.22	0.40
1:B:1099:SER:HB2	1:B:1268:ILE:HG21	2.03	0.40
1:C:1162:PHE:O	1:C:1168:SER:HB2	2.20	0.40
2:F:8:LEU:C	2:F:10:GLY:N	2.80	0.40
2:H:23:ILE:CD1	2:H:50:LEU:HB3	2.50	0.40
2:I:1:MET:HB2	2:I:63:LYS:HE2	2.04	0.40
2:I:4:PHE:CD2	2:I:14:THR:HG22	2.56	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	397/399 (100%)	372 (94%)	23 (6%)	2 (0%)	24	57
1	B	397/399 (100%)	372 (94%)	23 (6%)	2 (0%)	24	57
1	C	397/399 (100%)	371 (94%)	24 (6%)	2 (0%)	24	57
2	D	74/76 (97%)	66 (89%)	6 (8%)	2 (3%)	4	27
2	E	74/76 (97%)	66 (89%)	7 (10%)	1 (1%)	9	37
2	F	74/76 (97%)	66 (89%)	7 (10%)	1 (1%)	9	37
2	G	74/76 (97%)	73 (99%)	1 (1%)	0	100	100
2	H	74/76 (97%)	73 (99%)	1 (1%)	0	100	100
2	I	74/76 (97%)	72 (97%)	2 (3%)	0	100	100
All	All	1635/1653 (99%)	1531 (94%)	94 (6%)	10 (1%)	21	54

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1025	ALA
1	B	150	SER
1	B	1025	ALA
1	C	150	SER
1	C	1025	ALA
1	A	150	SER
2	D	61	ILE
2	D	17	VAL
2	E	17	VAL
2	F	17	VAL

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	354/354 (100%)	315 (89%)	39 (11%)	6	24
1	B	354/354 (100%)	313 (88%)	41 (12%)	5	22
1	C	354/354 (100%)	313 (88%)	41 (12%)	5	22

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	68/68 (100%)	57 (84%)	11 (16%)	2	14
2	E	68/68 (100%)	57 (84%)	11 (16%)	2	14
2	F	68/68 (100%)	57 (84%)	11 (16%)	2	14
2	G	68/68 (100%)	56 (82%)	12 (18%)	2	12
2	H	68/68 (100%)	56 (82%)	12 (18%)	2	12
2	I	68/68 (100%)	56 (82%)	12 (18%)	2	12
All	All	1470/1470 (100%)	1280 (87%)	190 (13%)	4	20

All (190) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	52	LEU
1	A	58	THR
1	A	70	THR
1	A	84	ILE
1	A	88	ILE
1	A	98	THR
1	A	107	CYS
1	A	119	LEU
1	A	140	GLU
1	A	1026	TYR
1	A	1033	GLN
1	A	1040	GLU
1	A	1041	ILE
1	A	1043	VAL
1	A	1052	LEU
1	A	1067	ILE
1	A	1069	GLN
1	A	1076	HIS
1	A	1077	LYS
1	A	1085	THR
1	A	1100	HIS
1	A	1109	LYS
1	A	1131	THR
1	A	1141	THR
1	A	1145	LEU
1	A	1152	GLN
1	A	1153	THR
1	A	1159	LEU
1	A	1162	PHE

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Mol	Chain	Res	Type
1	A	1185	ARG
1	A	1195	GLU
1	A	1208	VAL
1	A	1221	ILE
1	A	1231	ILE
1	A	1233	VAL
1	A	1238	ARG
1	A	1240	GLU
1	A	1251	GLU
1	A	1260	LEU
1	B	13	LEU
1	B	16	ASP
1	B	52	LEU
1	B	58	THR
1	B	70	THR
1	B	84	ILE
1	B	88	ILE
1	B	98	THR
1	B	107	CYS
1	B	119	LEU
1	B	140	GLU
1	B	1026	TYR
1	B	1033	GLN
1	B	1040	GLU
1	B	1041	ILE
1	B	1043	VAL
1	B	1052	LEU
1	B	1067	ILE
1	B	1069	GLN
1	B	1076	HIS
1	B	1077	LYS
1	B	1085	THR
1	B	1100	HIS
1	B	1109	LYS
1	B	1131	THR
1	B	1141	THR
1	B	1145	LEU
1	B	1152	GLN
1	B	1153	THR
1	B	1159	LEU
1	B	1162	PHE
1	B	1185	ARG

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Mol	Chain	Res	Type
1	B	1195	GLU
1	B	1208	VAL
1	B	1221	ILE
1	B	1231	ILE
1	B	1233	VAL
1	B	1238	ARG
1	B	1240	GLU
1	B	1251	GLU
1	B	1260	LEU
1	C	13	LEU
1	C	16	ASP
1	C	52	LEU
1	C	58	THR
1	C	70	THR
1	C	84	ILE
1	C	88	ILE
1	C	98	THR
1	C	107	CYS
1	C	119	LEU
1	C	140	GLU
1	C	1026	TYR
1	C	1033	GLN
1	C	1040	GLU
1	C	1041	ILE
1	C	1043	VAL
1	C	1052	LEU
1	C	1067	ILE
1	C	1069	GLN
1	C	1076	HIS
1	C	1077	LYS
1	C	1085	THR
1	C	1100	HIS
1	C	1109	LYS
1	C	1131	THR
1	C	1141	THR
1	C	1145	LEU
1	C	1152	GLN
1	C	1153	THR
1	C	1159	LEU
1	C	1162	PHE
1	C	1185	ARG
1	C	1195	GLU

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Mol	Chain	Res	Type
1	C	1208	VAL
1	C	1221	ILE
1	C	1231	ILE
1	C	1233	VAL
1	C	1238	ARG
1	C	1240	GLU
1	C	1251	GLU
1	C	1260	LEU
2	D	2	GLN
2	D	3	ILE
2	D	8	LEU
2	D	15	LEU
2	D	30	ILE
2	D	43	LEU
2	D	55	THR
2	D	63	LYS
2	D	67	LEU
2	D	69	LEU
2	D	73	LEU
2	E	2	GLN
2	E	3	ILE
2	E	8	LEU
2	E	15	LEU
2	E	30	ILE
2	E	43	LEU
2	E	55	THR
2	E	63	LYS
2	E	67	LEU
2	E	69	LEU
2	E	73	LEU
2	F	2	GLN
2	F	3	ILE
2	F	8	LEU
2	F	15	LEU
2	F	30	ILE
2	F	43	LEU
2	F	55	THR
2	F	63	LYS
2	F	67	LEU
2	F	69	LEU
2	F	73	LEU
2	G	3	ILE

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Mol	Chain	Res	Type
2	G	5	VAL
2	G	16	GLU
2	G	22	THR
2	G	23	ILE
2	G	52	ASP
2	G	55	THR
2	G	61	ILE
2	G	67	LEU
2	G	72	ARG
2	G	73	LEU
2	G	74	ARG
2	H	3	ILE
2	H	5	VAL
2	H	16	GLU
2	H	22	THR
2	H	23	ILE
2	H	52	ASP
2	H	55	THR
2	H	61	ILE
2	H	67	LEU
2	H	72	ARG
2	H	73	LEU
2	H	74	ARG
2	I	3	ILE
2	I	5	VAL
2	I	16	GLU
2	I	22	THR
2	I	23	ILE
2	I	52	ASP
2	I	55	THR
2	I	61	ILE
2	I	67	LEU
2	I	72	ARG
2	I	73	LEU
2	I	74	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	HIS
1	A	1066	ASN
1	A	1192	HIS

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Mol	Chain	Res	Type
1	A	1247	HIS
1	B	55	HIS
1	B	1066	ASN
1	B	1192	HIS
1	C	55	HIS
1	C	135	ASN
1	C	1066	ASN
1	C	1192	HIS
1	C	1217	HIS
2	D	40	GLN
2	D	49	GLN
2	E	49	GLN
2	F	49	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	399/399 (100%)	-0.29	0 100 100	90, 159, 226, 269	0
1	B	399/399 (100%)	-0.45	0 100 100	105, 162, 227, 269	0
1	C	399/399 (100%)	-0.15	1 (0%) 90 78	88, 160, 226, 269	0
2	D	76/76 (100%)	-0.41	0 100 100	135, 201, 241, 262	0
2	E	76/76 (100%)	-0.42	0 100 100	137, 200, 241, 262	0
2	F	76/76 (100%)	0.02	0 100 100	135, 197, 242, 261	0
2	G	76/76 (100%)	-0.31	0 100 100	106, 153, 210, 237	0
2	H	76/76 (100%)	-0.32	0 100 100	126, 161, 212, 238	0
2	I	76/76 (100%)	-0.26	0 100 100	103, 152, 209, 238	0
All	All	1653/1653 (100%)	-0.29	1 (0%) 92 87	88, 164, 230, 269	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1233	VAL	2.8

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.