



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:31 AM UTC

PDB ID : 7CQJ / pdb_00007cqj
Title : Peroxiredoxin from Aeropyrum pernix K1 (ApPrx) C50S/K84A/C207S/C213
S mutant (ApPrx*K84A)
Authors : Himiyama, T.; Nakamura, T.
Deposited on : 2020-08-11
Resolution : 2.90 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

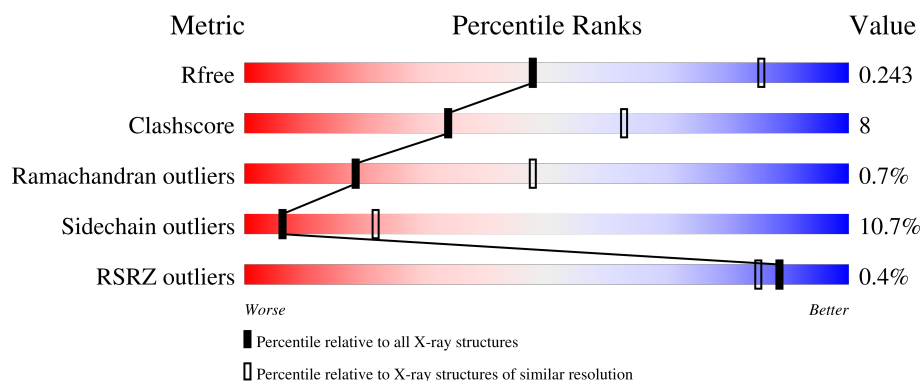
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	250	70% 24% ..
1	B	250	72% 23% ..
1	C	250	74% 21% ..
1	D	250	71% 25% ..
1	E	250	70% 24% ..

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Mol	Chain	Length	Quality of chain
1	F	250	68%25%
1	G	250	67%28%
1	H	250	74%20%
1	I	250	69%26%
1	J	250	73%22%
1	K	250	70%23%
1	L	250	76%19%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 23675 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 Peroxiredoxin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0
1	B	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0
1	C	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0
1	D	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0
1	E	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0
1	F	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0
1	G	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0
1	H	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0
1	I	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0
1	J	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0
1	K	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0
1	L	244	Total 1969	C 1265	N 346	O 354	S 4	0	0	0

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	50	SER	CYS	engineered mutation	UNP Q9Y9L0
A	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
A	207	SER	CYS	engineered mutation	UNP Q9Y9L0
A	213	SER	CYS	engineered mutation	UNP Q9Y9L0
B	50	SER	CYS	engineered mutation	UNP Q9Y9L0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
B	207	SER	CYS	engineered mutation	UNP Q9Y9L0
B	213	SER	CYS	engineered mutation	UNP Q9Y9L0
C	50	SER	CYS	engineered mutation	UNP Q9Y9L0
C	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
C	207	SER	CYS	engineered mutation	UNP Q9Y9L0
C	213	SER	CYS	engineered mutation	UNP Q9Y9L0
D	50	SER	CYS	engineered mutation	UNP Q9Y9L0
D	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
D	207	SER	CYS	engineered mutation	UNP Q9Y9L0
D	213	SER	CYS	engineered mutation	UNP Q9Y9L0
E	50	SER	CYS	engineered mutation	UNP Q9Y9L0
E	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
E	207	SER	CYS	engineered mutation	UNP Q9Y9L0
E	213	SER	CYS	engineered mutation	UNP Q9Y9L0
F	50	SER	CYS	engineered mutation	UNP Q9Y9L0
F	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
F	207	SER	CYS	engineered mutation	UNP Q9Y9L0
F	213	SER	CYS	engineered mutation	UNP Q9Y9L0
G	50	SER	CYS	engineered mutation	UNP Q9Y9L0
G	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
G	207	SER	CYS	engineered mutation	UNP Q9Y9L0
G	213	SER	CYS	engineered mutation	UNP Q9Y9L0
H	50	SER	CYS	engineered mutation	UNP Q9Y9L0
H	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
H	207	SER	CYS	engineered mutation	UNP Q9Y9L0
H	213	SER	CYS	engineered mutation	UNP Q9Y9L0
I	50	SER	CYS	engineered mutation	UNP Q9Y9L0
I	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
I	207	SER	CYS	engineered mutation	UNP Q9Y9L0
I	213	SER	CYS	engineered mutation	UNP Q9Y9L0
J	50	SER	CYS	engineered mutation	UNP Q9Y9L0
J	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
J	207	SER	CYS	engineered mutation	UNP Q9Y9L0
J	213	SER	CYS	engineered mutation	UNP Q9Y9L0
K	50	SER	CYS	engineered mutation	UNP Q9Y9L0
K	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
K	207	SER	CYS	engineered mutation	UNP Q9Y9L0
K	213	SER	CYS	engineered mutation	UNP Q9Y9L0
L	50	SER	CYS	engineered mutation	UNP Q9Y9L0
L	84	ALA	LYS	engineered mutation	UNP Q9Y9L0
L	207	SER	CYS	engineered mutation	UNP Q9Y9L0

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Chain	Residue	Modelled	Actual	Comment	Reference
L	213	SER	CYS	engineered mutation	UNP Q9Y9L0

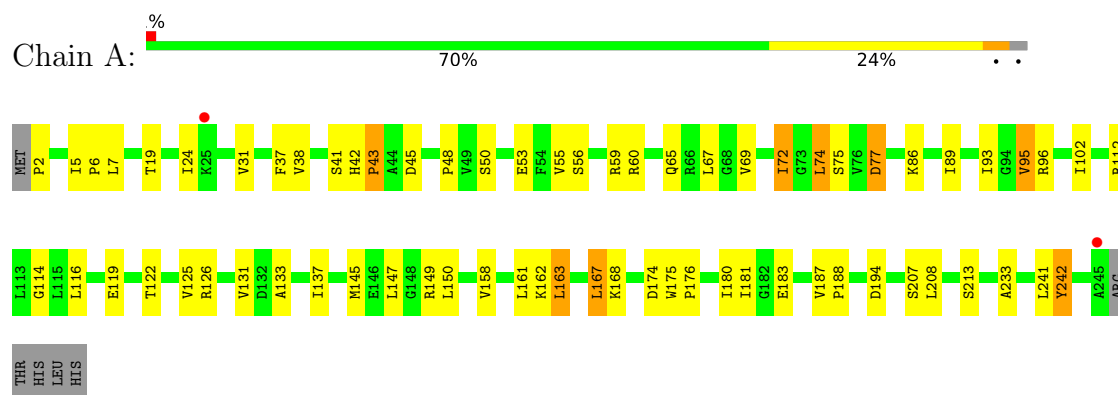
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	6	Total O 6 6	0	0
2	B	8	Total O 8 8	0	0
2	C	3	Total O 3 3	0	0
2	D	5	Total O 5 5	0	0
2	E	4	Total O 4 4	0	0
2	F	2	Total O 2 2	0	0
2	G	2	Total O 2 2	0	0
2	H	7	Total O 7 7	0	0
2	I	3	Total O 3 3	0	0
2	J	4	Total O 4 4	0	0
2	L	3	Total O 3 3	0	0

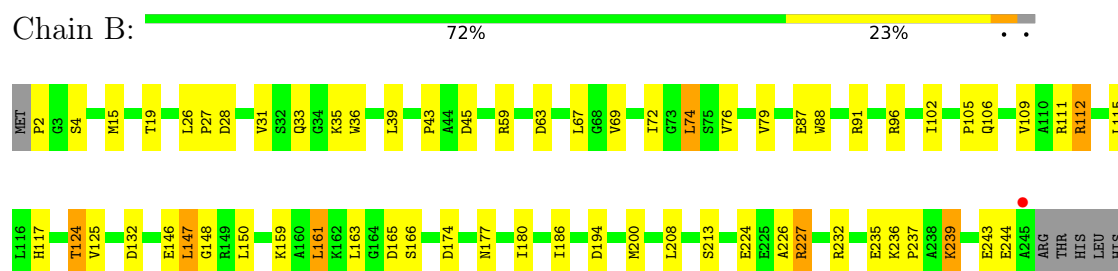
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

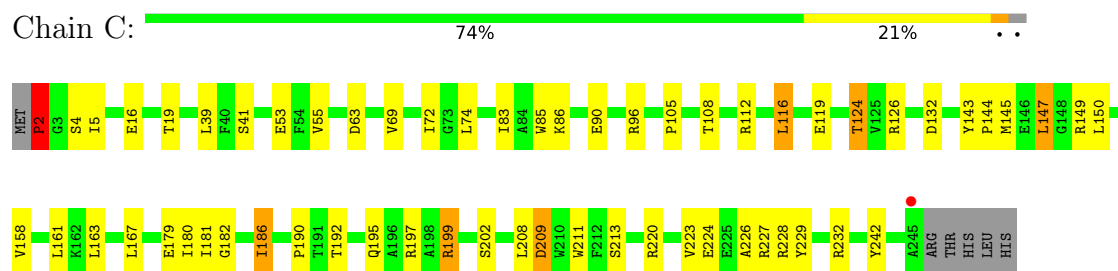
• Molecule 1: Peroxiredoxin



• Molecule 1: Peroxiredoxin

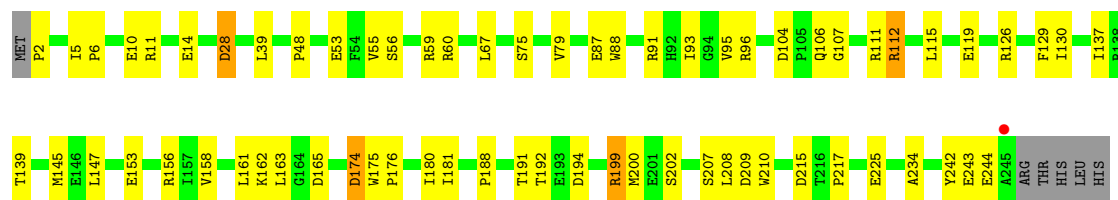


• Molecule 1: Peroxiredoxin



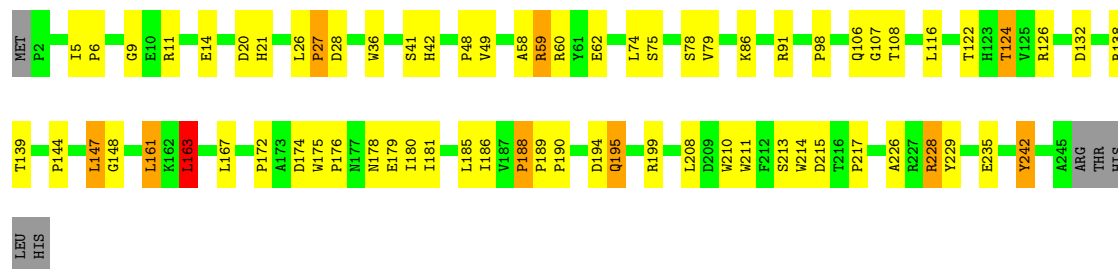
• Molecule 1: Peroxiredoxin





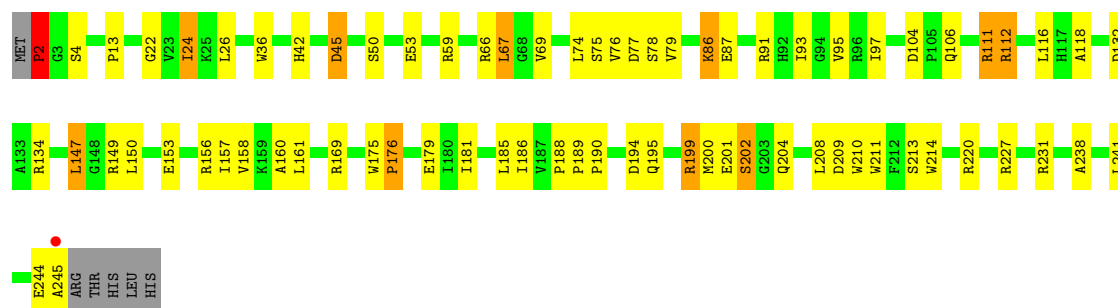
• Molecule 1: Peroxiredoxin

Chain E: 70% 24%



• Molecule 1: Peroxiredoxin

Chain F: 68% 25%



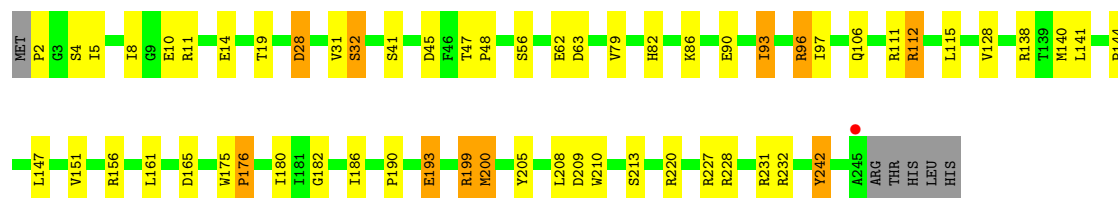
• Molecule 1: Peroxiredoxin

Chain G: 67% 28%



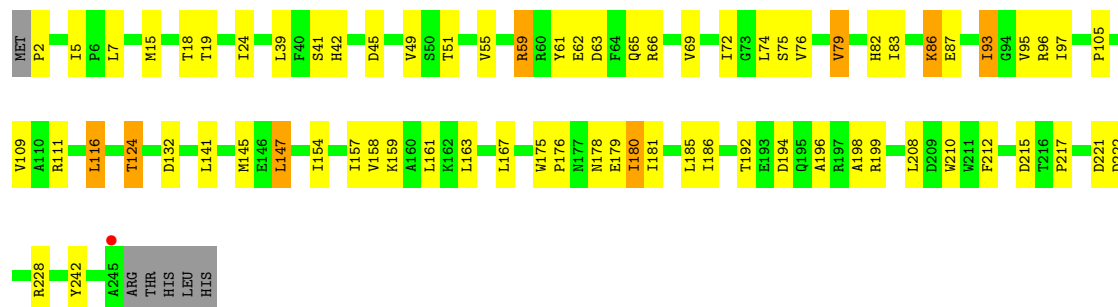
• Molecule 1: Peroxiredoxin

Chain H: 74% 20%



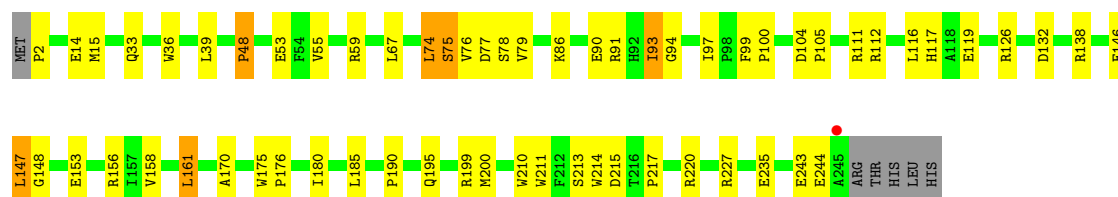
• Molecule 1: Peroxiredoxin

Chain I: 69% 26% ..



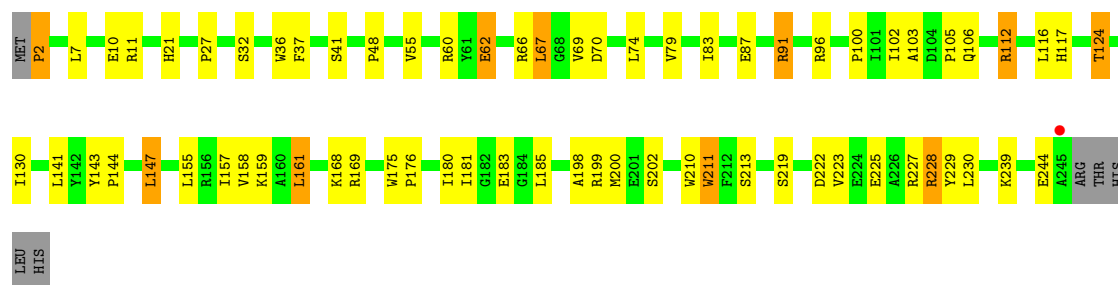
• Molecule 1: Peroxiredoxin

Chain J: 73% 22% ..



• Molecule 1: Peroxiredoxin

Chain K: 70% 23% ..



• Molecule 1: Peroxiredoxin

Chain L: 76% 19% ..





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.01Å 210.57Å 308.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.48 – 2.90 49.48 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.48-2.90) 99.8 (49.48-2.90)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.48 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.179 , 0.256 0.171 , 0.243	Depositor DCC
R_{free} test set	4355 reflections (5.19%)	wwPDB-VP
Wilson B-factor (Å ²)	63.2	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	23675	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 20.61 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 8.3704e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.04	0/2024	1.56	9/2753 (0.3%)
1	B	1.05	0/2024	1.53	5/2753 (0.2%)
1	C	1.07	0/2024	1.45	2/2753 (0.1%)
1	D	1.04	0/2024	1.50	6/2753 (0.2%)
1	E	1.06	0/2024	1.54	18/2753 (0.7%)
1	F	1.04	0/2024	1.53	8/2753 (0.3%)
1	G	1.03	0/2024	1.49	4/2753 (0.1%)
1	H	1.06	1/2024 (0.0%)	1.55	6/2753 (0.2%)
1	I	1.04	0/2024	1.46	5/2753 (0.2%)
1	J	1.04	0/2024	1.44	2/2753 (0.1%)
1	K	1.08	0/2024	1.51	4/2753 (0.1%)
1	L	1.07	0/2024	1.50	3/2753 (0.1%)
All	All	1.05	1/24288 (0.0%)	1.51	72/33036 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	93	ILE	N-CA	5.25	1.50	1.46

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	160	ALA	CA-C-N	7.40	130.52	120.38
1	F	160	ALA	C-N-CA	7.40	130.52	120.38
1	A	188	PRO	CB-CA-C	6.91	117.89	111.39
1	H	62	GLU	CB-CG-CD	6.72	124.03	112.60
1	I	198	ALA	CA-C-N	6.44	128.81	120.44
1	I	198	ALA	C-N-CA	6.44	128.81	120.44
1	E	106	GLN	CB-CA-C	6.40	120.04	111.14
1	I	228	ARG	CB-CA-C	6.30	120.77	110.88
1	I	2	PRO	CA-N-CD	-6.26	103.23	112.00
1	A	2	PRO	CA-N-CD	-6.15	103.39	112.00
1	A	45	ASP	CA-CB-CG	6.00	118.60	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	187	VAL	N-CA-CB	5.97	115.63	110.08
1	K	2	PRO	CA-N-CD	-5.94	103.68	112.00
1	F	2	PRO	CA-N-CD	-5.87	103.78	112.00
1	F	157	ILE	CA-C-O	-5.87	115.01	121.29
1	L	230	LEU	CA-C-N	5.78	128.30	120.44
1	L	230	LEU	C-N-CA	5.78	128.30	120.44
1	K	222	ASP	CA-CB-CG	5.71	118.31	112.60
1	B	132	ASP	CA-C-N	5.69	129.36	120.82
1	B	132	ASP	C-N-CA	5.69	129.36	120.82
1	A	188	PRO	N-CA-C	-5.68	105.02	110.47
1	A	242	TYR	CA-C-N	5.65	128.63	120.38
1	A	242	TYR	C-N-CA	5.65	128.63	120.38
1	F	45	ASP	CA-CB-CG	5.63	118.23	112.60
1	L	70	ASP	CB-CA-C	5.62	119.07	109.80
1	E	6	PRO	N-CA-C	-5.59	102.80	111.13
1	F	22	GLY	CA-C-O	-5.54	117.20	122.13
1	C	2	PRO	CA-N-CD	-5.53	104.25	112.00
1	I	228	ARG	N-CA-C	-5.47	105.21	111.07
1	J	77	ASP	CA-C-N	5.42	128.54	120.90
1	J	77	ASP	C-N-CA	5.42	128.54	120.90
1	G	195	GLN	CB-CA-C	-5.40	102.33	110.92
1	E	126	ARG	CA-C-N	5.38	126.04	120.60
1	E	126	ARG	C-N-CA	5.38	126.04	120.60
1	E	186	ILE	CA-C-N	5.37	126.46	122.59
1	E	186	ILE	C-N-CA	5.37	126.46	122.59
1	H	47	THR	CB-CA-C	5.36	117.17	108.87
1	E	49	VAL	N-CA-C	-5.27	105.36	110.42
1	B	45	ASP	CA-CB-CG	5.23	117.83	112.60
1	D	234	ALA	CA-C-N	5.23	129.58	121.26
1	D	234	ALA	C-N-CA	5.23	129.58	121.26
1	F	160	ALA	O-C-N	5.22	127.45	122.07
1	G	123	HIS	CA-C-N	5.20	128.10	120.71
1	G	123	HIS	C-N-CA	5.20	128.10	120.71
1	B	165	ASP	CA-CB-CG	5.17	117.77	112.60
1	E	178	ASN	CA-C-N	5.16	127.20	120.28
1	E	178	ASN	C-N-CA	5.16	127.20	120.28
1	B	2	PRO	CA-N-CD	-5.15	104.80	112.00
1	D	163	LEU	CA-C-N	5.13	125.54	119.94
1	D	163	LEU	C-N-CA	5.13	125.54	119.94
1	F	231	ARG	CB-CA-C	5.13	119.57	110.85
1	G	195	GLN	CB-CG-CD	-5.13	103.88	112.60
1	A	89	ILE	CA-C-N	5.13	127.46	120.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	89	ILE	C-N-CA	5.13	127.46	120.54
1	E	28	ASP	CA-C-N	5.13	127.10	120.44
1	E	28	ASP	C-N-CA	5.13	127.10	120.44
1	H	31	VAL	N-CA-C	-5.11	105.52	110.42
1	H	19	THR	CA-CB-OG1	-5.10	101.95	109.60
1	K	27	PRO	CA-C-N	5.09	127.41	120.54
1	K	27	PRO	C-N-CA	5.09	127.41	120.54
1	D	165	ASP	CA-C-N	5.07	127.33	120.38
1	D	165	ASP	C-N-CA	5.07	127.33	120.38
1	C	96	ARG	CB-CA-C	5.07	118.11	109.75
1	E	144	PRO	CA-C-N	5.07	129.25	121.19
1	E	144	PRO	C-N-CA	5.07	129.25	121.19
1	E	27	PRO	CA-C-N	5.06	127.31	120.38
1	E	27	PRO	C-N-CA	5.06	127.31	120.38
1	H	200	MET	CA-C-N	5.06	127.06	120.28
1	H	200	MET	C-N-CA	5.06	127.06	120.28
1	E	195	GLN	CB-CG-CD	-5.03	104.05	112.60
1	E	163	LEU	CA-C-N	5.03	125.67	119.99
1	E	163	LEU	C-N-CA	5.03	125.67	119.99

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1969	0	1951	31	0
1	B	1969	0	1951	30	0
1	C	1969	0	1951	31	0
1	D	1969	0	1951	34	0
1	E	1969	0	1951	32	0
1	F	1969	0	1951	37	0
1	G	1969	0	1951	43	0
1	H	1969	0	1951	30	0
1	I	1969	0	1951	37	0
1	J	1969	0	1951	35	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	K	1969	0	1951	46	0
1	L	1969	0	1951	31	0
2	A	6	0	0	0	0
2	B	8	0	0	0	0
2	C	3	0	0	0	0
2	D	5	0	0	0	0
2	E	4	0	0	0	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	7	0	0	0	0
2	I	3	0	0	0	0
2	J	4	0	0	0	0
2	L	3	0	0	0	0
All	All	23675	0	23412	363	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 8.

All (363) 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:2:PRO:HB2	1:D:10:GLU:OE2	1.75	0.85
1:A:75:SER:OG	1:A:77:ASP:OD1	1.98	0.80
1:A:69:VAL:HG21	1:A:158:VAL:HG11	1.67	0.76
1:I:69:VAL:HG21	1:I:158:VAL:HG11	1.67	0.75
1:G:69:VAL:HG21	1:G:158:VAL:HG11	1.67	0.75
1:F:24:ILE:HD11	1:F:26:LEU:HD21	1.70	0.74
1:A:37:PHE:HD2	1:A:72:ILE:HD11	1.52	0.73
1:K:67:LEU:HD21	1:K:159:LYS:HD2	1.71	0.73
1:C:69:VAL:HG21	1:C:158:VAL:HG11	1.70	0.72
1:F:69:VAL:HG21	1:F:158:VAL:HG21	1.73	0.71
1:K:69:VAL:HG21	1:K:158:VAL:HG11	1.72	0.71
1:J:74:LEU:HD13	1:J:75:SER:N	2.07	0.69
1:A:31:VAL:HG13	1:A:133:ALA:O	1.91	0.69
1:I:215:ASP:OD1	1:I:217:PRO:HD3	1.93	0.69
1:C:199:ARG:HG3	1:C:199:ARG:HH11	1.57	0.68
1:L:216:THR:O	1:L:216:THR:OG1	2.11	0.68
1:K:2:PRO:O	1:L:7:LEU:HG	1.94	0.68
1:A:74:LEU:HD23	1:A:102:ILE:HB	1.75	0.67
1:G:144:PRO:HG3	1:H:138:ARG:O	1.95	0.66
1:D:158:VAL:O	1:D:162:LYS:HG3	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:106:GLN:O	1:D:111:ARG:NH2	2.29	0.64
1:F:156:ARG:HD2	1:F:175:TRP:O	1.97	0.63
1:D:188:PRO:O	1:D:199:ARG:NH2	2.31	0.63
1:J:147:LEU:HD22	1:J:148:GLY:O	1.98	0.63
1:F:200:MET:HE3	1:F:210:TRP:HA	1.80	0.62
1:G:2:PRO:HB2	1:H:10:GLU:OE2	2.00	0.62
1:D:39:LEU:C	1:D:39:LEU:HD23	2.25	0.62
1:L:67:LEU:HD13	1:L:158:VAL:HG23	1.80	0.62
1:E:20:ASP:OD2	1:E:86:LYS:NZ	2.26	0.61
1:C:55:VAL:HG21	1:D:180:ILE:HD11	1.83	0.61
1:G:49:VAL:O	1:G:52:THR:OG1	2.19	0.60
1:K:69:VAL:CG2	1:K:158:VAL:HG11	2.31	0.60
1:C:199:ARG:HH11	1:C:199:ARG:CG	2.15	0.60
1:D:67:LEU:HD13	1:D:158:VAL:HG23	1.84	0.60
1:F:93:ILE:HG22	1:F:95:VAL:HG23	1.84	0.59
1:I:93:ILE:CG2	1:I:95:VAL:HG23	2.33	0.59
1:J:55:VAL:O	1:J:59:ARG:HG3	2.03	0.59
1:K:112:ARG:HG3	1:K:112:ARG:HH11	1.66	0.59
1:F:244:GLU:O	1:F:245:ALA:C	2.46	0.59
1:A:150:LEU:HD12	1:B:174:ASP:HA	1.84	0.59
1:K:112:ARG:HH11	1:K:112:ARG:CG	2.16	0.58
1:G:41:SER:HB2	1:G:124:THR:CG2	2.34	0.58
1:G:93:ILE:HG22	1:G:95:VAL:HG23	1.85	0.58
1:L:146:GLU:OE2	1:L:146:GLU:N	2.36	0.58
1:E:11:ARG:NH1	1:E:11:ARG:HG2	2.18	0.58
1:A:175:TRP:CG	1:A:176:PRO:HA	2.39	0.58
1:I:49:VAL:HG21	1:J:170:ALA:HB1	1.86	0.58
1:B:147:LEU:HD22	1:B:148:GLY:O	2.02	0.57
1:H:86:LYS:HD3	1:H:97:ILE:HB	1.87	0.57
1:I:93:ILE:HG22	1:I:95:VAL:HG23	1.87	0.57
1:I:55:VAL:HG21	1:J:180:ILE:HD11	1.87	0.57
1:B:87:GLU:OE2	1:C:209:ASP:OD1	2.22	0.56
1:F:36:TRP:HB2	1:F:69:VAL:HG22	1.88	0.56
1:F:45:ASP:N	1:F:77:ASP:OD2	2.21	0.56
1:J:243:GLU:HB2	1:J:244:GLU:OE1	2.05	0.56
1:G:14:GLU:O	1:G:112:ARG:NH2	2.38	0.56
1:I:45:ASP:OD2	1:I:82:HIS:ND1	2.33	0.56
1:F:132:ASP:OD2	1:F:134:ARG:NH1	2.39	0.56
1:C:53:GLU:OE2	1:C:149:ARG:NE	2.34	0.55
1:K:211:TRP:H	1:K:211:TRP:CD1	2.24	0.55
1:J:74:LEU:HD13	1:J:74:LEU:C	2.30	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:41:SER:HB2	1:I:124:THR:HG21	1.89	0.55
1:D:175:TRP:CG	1:D:176:PRO:HA	2.41	0.55
1:A:72:ILE:H	1:A:72:ILE:HD12	1.72	0.55
1:I:157:ILE:O	1:I:161:LEU:HB2	2.06	0.54
1:K:36:TRP:HH2	1:K:161:LEU:HB3	1.71	0.54
1:A:233:ALA:O	1:B:177:ASN:ND2	2.35	0.54
1:D:180:ILE:HG22	1:D:181:ILE:HG23	1.88	0.54
1:G:39:LEU:C	1:G:39:LEU:HD23	2.32	0.54
1:A:72:ILE:HD12	1:A:72:ILE:N	2.23	0.54
1:C:112:ARG:HG3	1:C:112:ARG:HH11	1.72	0.54
1:G:132:ASP:OD2	1:G:134:ARG:NH1	2.41	0.54
1:D:153:GLU:OE1	1:D:153:GLU:HA	2.08	0.53
1:K:143:TYR:OH	1:L:153:GLU:OE2	2.23	0.53
1:B:15:MET:SD	1:B:109:VAL:HG13	2.47	0.53
1:C:179:GLU:OE2	1:D:59:ARG:HD3	2.08	0.53
1:F:202:SER:OG	1:F:204:GLN:NE2	2.42	0.53
1:A:41:SER:HB3	1:A:74:LEU:HD12	1.89	0.53
1:A:93:ILE:HG22	1:A:95:VAL:HG23	1.89	0.53
1:D:55:VAL:O	1:D:59:ARG:HG3	2.08	0.53
1:F:36:TRP:CD2	1:F:132:ASP:HA	2.44	0.53
1:I:69:VAL:CG2	1:I:158:VAL:HG11	2.37	0.53
1:B:27:PRO:O	1:B:31:VAL:HG23	2.08	0.53
1:G:20:ASP:OD2	1:G:86:LYS:HE3	2.09	0.53
1:D:93:ILE:HG22	1:D:95:VAL:HG23	1.90	0.53
1:K:21:HIS:ND1	1:K:100:PRO:HB3	2.23	0.53
1:H:56:SER:HB3	1:H:151:VAL:HG21	1.89	0.52
1:C:41:SER:HB2	1:C:124:THR:HG21	1.89	0.52
1:I:42:HIS:CE1	1:I:75:SER:HB2	2.44	0.52
1:L:5:ILE:HD12	1:L:140:MET:HE1	1.92	0.52
1:E:147:LEU:HD22	1:E:148:GLY:O	2.09	0.52
1:L:36:TRP:CD2	1:L:132:ASP:HA	2.45	0.52
1:H:90:GLU:OE2	1:H:96:ARG:NH2	2.43	0.52
1:B:39:LEU:HA	1:B:72:ILE:O	2.10	0.52
1:C:150:LEU:HD12	1:D:174:ASP:HA	1.91	0.52
1:E:161:LEU:HD13	1:F:147:LEU:HG	1.91	0.52
1:I:178:ASN:HB3	1:I:181:ILE:O	2.10	0.52
1:I:180:ILE:HD12	1:J:93:ILE:HD13	1.92	0.52
1:C:143:TYR:HD2	1:C:147:LEU:HD13	1.76	0.51
1:G:175:TRP:CG	1:G:176:PRO:HA	2.46	0.51
1:L:59:ARG:HH12	1:L:241:LEU:HD11	1.74	0.51
1:E:14:GLU:HA	1:E:26:LEU:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:ILE:HB	1:H:210:TRP:CZ2	2.45	0.51
1:L:190:PRO:HD2	1:L:211:TRP:HA	1.91	0.51
1:D:210:TRP:CE2	1:G:83:ILE:HG22	2.46	0.51
1:B:200:MET:HA	1:B:200:MET:HE2	1.91	0.51
1:D:56:SER:OG	1:D:60:ARG:NH1	2.45	0.50
1:F:53:GLU:CD	1:F:149:ARG:HH21	2.19	0.50
1:L:90:GLU:O	1:L:94:GLY:HA2	2.11	0.50
1:K:183:GLU:O	1:K:183:GLU:HG3	2.12	0.50
1:E:41:SER:HB2	1:E:124:THR:HG21	1.93	0.50
1:G:11:ARG:HG2	1:G:11:ARG:HH11	1.77	0.50
1:H:156:ARG:HD2	1:H:175:TRP:O	2.11	0.50
1:K:7:LEU:O	1:K:10:GLU:HB2	2.12	0.49
1:D:112:ARG:CG	1:D:112:ARG:HH11	2.25	0.49
1:J:99:PHE:HB2	1:J:100:PRO:HD2	1.93	0.49
1:D:53:GLU:OE1	1:D:126:ARG:NH2	2.45	0.49
1:I:74:LEU:HD13	1:I:74:LEU:C	2.37	0.49
1:J:175:TRP:CG	1:J:176:PRO:HA	2.48	0.49
1:K:200:MET:HE2	1:K:200:MET:N	2.27	0.49
1:L:6:PRO:HB2	1:L:137:ILE:HD13	1.94	0.49
1:B:239:LYS:NZ	1:B:243:GLU:HB2	2.28	0.49
1:I:154:ILE:O	1:I:158:VAL:HG23	2.12	0.49
1:J:67:LEU:HD13	1:J:158:VAL:HG23	1.94	0.49
1:J:153:GLU:OE1	1:J:153:GLU:HA	2.13	0.49
1:C:223:VAL:HG12	1:C:227:ARG:HH12	1.77	0.49
1:C:69:VAL:CG2	1:C:158:VAL:HG11	2.41	0.49
1:F:175:TRP:CG	1:F:176:PRO:HA	2.47	0.49
1:G:5:ILE:HD13	1:H:5:ILE:HD13	1.95	0.49
1:L:74:LEU:HD23	1:L:102:ILE:HB	1.95	0.48
1:L:243:GLU:HA	1:L:243:GLU:OE2	2.13	0.48
1:G:41:SER:HB2	1:G:124:THR:HG21	1.95	0.48
1:I:196:ALA:HB1	1:I:210:TRP:HB3	1.96	0.48
1:L:147:LEU:HD22	1:L:148:GLY:O	2.13	0.48
1:C:180:ILE:HG22	1:C:181:ILE:HG23	1.95	0.48
1:F:188:PRO:O	1:F:199:ARG:NH2	2.42	0.48
1:I:7:LEU:HD13	1:J:117:HIS:HA	1.95	0.48
1:J:86:LYS:HE3	1:J:97:ILE:HB	1.95	0.48
1:K:62:GLU:HA	1:K:62:GLU:OE2	2.12	0.48
1:K:74:LEU:HD23	1:K:102:ILE:HB	1.95	0.48
1:E:11:ARG:HG2	1:E:11:ARG:HH11	1.79	0.48
1:J:185:LEU:O	1:J:214:TRP:HB2	2.14	0.48
1:K:175:TRP:CG	1:K:176:PRO:HA	2.49	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:153:GLU:HA	1:F:153:GLU:OE1	2.14	0.48
1:G:142:TYR:O	1:H:8:ILE:HD11	2.14	0.48
1:G:196:ALA:HB1	1:G:210:TRP:HB3	1.95	0.48
1:H:45:ASP:OD2	1:H:82:HIS:ND1	2.40	0.48
1:K:103:ALA:C	1:K:105:PRO:HD3	2.39	0.48
1:F:111:ARG:HG3	1:F:116:LEU:HD12	1.95	0.48
1:H:199:ARG:HD2	1:H:205:TYR:CD2	2.49	0.48
1:G:141:LEU:HD22	1:H:141:LEU:HD22	1.95	0.47
1:H:200:MET:HE2	1:H:200:MET:HA	1.97	0.47
1:E:180:ILE:HA	1:F:241:LEU:HB2	1.96	0.47
1:F:13:PRO:HB3	1:F:112:ARG:NH2	2.29	0.47
1:K:74:LEU:HD23	1:K:102:ILE:CG2	2.45	0.47
1:A:180:ILE:HG22	1:A:181:ILE:HG23	1.97	0.47
1:C:190:PRO:HD2	1:C:211:TRP:HA	1.96	0.47
1:D:14:GLU:O	1:D:112:ARG:NH2	2.48	0.47
1:K:60:ARG:HG2	1:K:155:LEU:HD11	1.97	0.47
1:K:144:PRO:HD3	1:L:139:THR:OG1	2.14	0.47
1:K:175:TRP:CE2	1:K:176:PRO:HB3	2.50	0.47
1:L:226:ALA:HA	1:L:229:TYR:CD2	2.50	0.47
1:E:36:TRP:CD2	1:E:132:ASP:HA	2.50	0.47
1:I:175:TRP:CG	1:I:176:PRO:HA	2.50	0.47
1:K:112:ARG:CG	1:K:112:ARG:NH1	2.78	0.47
1:B:236:LYS:HG3	1:B:237:PRO:CD	2.45	0.47
1:D:14:GLU:OE1	1:D:28:ASP:OD1	2.32	0.47
1:E:215:ASP:OD1	1:E:217:PRO:HD3	2.15	0.47
1:I:79:VAL:O	1:I:82:HIS:HB2	2.15	0.47
1:A:59:ARG:HH22	1:A:241:LEU:HD21	1.80	0.47
1:K:37:PHE:HA	1:K:70:ASP:O	2.14	0.47
1:E:9:GLY:HA3	1:F:118:ALA:HB3	1.97	0.46
1:G:139:THR:OG1	1:H:144:PRO:HD3	2.15	0.46
1:G:156:ARG:HD2	1:G:175:TRP:O	2.15	0.46
1:J:15:MET:HB3	1:J:112:ARG:NH2	2.30	0.46
1:A:37:PHE:CD2	1:A:72:ILE:HD11	2.42	0.46
1:A:174:ASP:HA	1:B:150:LEU:HD12	1.96	0.46
1:H:112:ARG:HG3	1:H:112:ARG:HH11	1.80	0.46
1:A:53:GLU:OE1	1:A:126:ARG:NH2	2.49	0.46
1:G:31:VAL:HG13	1:G:133:ALA:O	2.15	0.46
1:K:180:ILE:HG22	1:K:181:ILE:HG23	1.97	0.46
1:A:6:PRO:HB2	1:A:137:ILE:CD1	2.46	0.46
1:F:150:LEU:HD23	1:F:153:GLU:HB2	1.97	0.46
1:J:215:ASP:C	1:J:217:PRO:HD3	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:175:TRP:CG	1:H:176:PRO:HA	2.51	0.46
1:H:228:ARG:HA	1:H:231:ARG:HG3	1.97	0.46
1:I:61:TYR:CE1	1:I:65:GLN:NE2	2.84	0.46
1:C:86:LYS:HD3	1:H:193:GLU:OE1	2.16	0.45
1:E:172:PRO:HG3	1:E:181:ILE:HD11	1.98	0.45
1:L:115:LEU:O	1:L:124:THR:HA	2.15	0.45
1:D:130:ILE:HD13	1:D:139:THR:HB	1.98	0.45
1:E:42:HIS:CE1	1:E:75:SER:HB2	2.51	0.45
1:E:175:TRP:CG	1:E:176:PRO:HA	2.52	0.45
1:E:185:LEU:O	1:E:214:TRP:HB2	2.16	0.45
1:H:41:SER:HB3	1:H:115:LEU:HD13	1.98	0.45
1:K:200:MET:HE2	1:K:200:MET:HA	1.97	0.45
1:L:153:GLU:OE1	1:L:153:GLU:HA	2.17	0.45
1:L:179:GLU:CD	1:L:179:GLU:H	2.24	0.45
1:A:7:LEU:HD13	1:B:117:HIS:HA	1.98	0.45
1:J:90:GLU:O	1:J:94:GLY:HA2	2.16	0.45
1:K:180:ILE:HG22	1:K:181:ILE:CG2	2.46	0.45
1:A:67:LEU:O	1:A:162:LYS:NZ	2.31	0.45
1:D:115:LEU:HD21	1:D:129:PHE:HE1	1.81	0.45
1:G:104:ASP:N	1:G:105:PRO:HD3	2.32	0.45
1:H:11:ARG:NH1	1:H:14:GLU:OE2	2.49	0.45
1:I:39:LEU:HA	1:I:72:ILE:O	2.17	0.45
1:I:132:ASP:OD1	1:I:132:ASP:C	2.59	0.45
1:J:200:MET:N	1:J:200:MET:HE2	2.31	0.45
1:L:126:ARG:HB3	1:L:149:ARG:CZ	2.47	0.45
1:A:55:VAL:HG21	1:B:180:ILE:HD11	1.98	0.45
1:C:53:GLU:CD	1:C:149:ARG:HH21	2.25	0.45
1:K:67:LEU:HD11	1:K:229:TYR:OH	2.17	0.45
1:D:175:TRP:CD1	1:D:176:PRO:HA	2.52	0.45
1:J:104:ASP:N	1:J:105:PRO:HD3	2.32	0.45
1:F:76:VAL:HA	1:F:104:ASP:O	2.17	0.45
1:E:48:PRO:HG2	1:F:186:ILE:HG21	1.98	0.44
1:K:200:MET:HE2	1:K:200:MET:CA	2.48	0.44
1:B:63:ASP:OD2	1:B:232:ARG:NH1	2.46	0.44
1:C:63:ASP:OD2	1:C:232:ARG:NH2	2.39	0.44
1:G:41:SER:HB2	1:G:124:THR:HG23	1.99	0.44
1:J:119:GLU:OE1	1:J:146:GLU:OE2	2.35	0.44
1:J:200:MET:HE3	1:J:210:TRP:HA	1.99	0.44
1:K:147:LEU:HG	1:L:161:LEU:HD13	1.98	0.44
1:E:188:PRO:HA	1:E:189:PRO:HD2	1.87	0.44
1:A:56:SER:OG	1:A:60:ARG:NH1	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:198:ALA:O	1:K:202:SER:HB3	2.18	0.44
1:E:59:ARG:NH1	1:F:179:GLU:OE1	2.50	0.44
1:G:147:LEU:HD22	1:G:148:GLY:O	2.17	0.44
1:I:42:HIS:O	1:I:124:THR:HG22	2.18	0.44
1:I:221:ASP:O	1:I:222:ASP:C	2.60	0.44
1:K:141:LEU:HD22	1:L:141:LEU:HD22	2.00	0.44
1:E:226:ALA:HA	1:E:229:TYR:CD2	2.52	0.44
1:F:190:PRO:HB3	1:F:195:GLN:HB3	2.00	0.44
1:J:39:LEU:HD23	1:J:39:LEU:C	2.43	0.44
1:D:112:ARG:CG	1:D:112:ARG:NH1	2.80	0.43
1:G:40:PHE:HA	1:G:127:GLY:O	2.17	0.43
1:I:15:MET:SD	1:I:109:VAL:HG13	2.58	0.43
1:J:111:ARG:HG3	1:J:116:LEU:HD12	1.98	0.43
1:K:225:GLU:O	1:K:228:ARG:HG2	2.18	0.43
1:F:87:GLU:OE2	1:G:209:ASP:OD1	2.36	0.43
1:G:163:LEU:HD12	1:G:163:LEU:HA	1.80	0.43
1:K:48:PRO:CG	1:L:186:ILE:HG21	2.48	0.43
1:C:143:TYR:CD2	1:C:147:LEU:HD13	2.54	0.43
1:F:42:HIS:HB2	1:F:50:SER:OG	2.19	0.43
1:H:112:ARG:HH11	1:H:112:ARG:CG	2.31	0.43
1:J:211:TRP:CD1	1:J:211:TRP:H	2.35	0.43
1:B:112:ARG:O	1:B:112:ARG:HG3	2.18	0.43
1:F:86:LYS:HG2	1:F:97:ILE:HB	2.00	0.43
1:H:242:TYR:CD1	1:H:242:TYR:C	2.96	0.43
1:L:158:VAL:O	1:L:162:LYS:HG3	2.19	0.43
1:A:175:TRP:CD1	1:A:176:PRO:HA	2.54	0.43
1:C:105:PRO:O	1:G:105:PRO:O	2.37	0.43
1:B:88:TRP:CE3	1:B:88:TRP:C	2.97	0.43
1:D:200:MET:HE1	1:D:207:SER:OG	2.18	0.43
1:E:190:PRO:HB3	1:E:195:GLN:HB3	1.99	0.43
1:G:51:THR:O	1:G:55:VAL:HG23	2.18	0.43
1:I:76:VAL:O	1:I:105:PRO:HA	2.19	0.43
1:J:36:TRP:CD2	1:J:132:ASP:HA	2.54	0.43
1:J:190:PRO:HA	1:J:195:GLN:NE2	2.34	0.43
1:C:132:ASP:OD1	1:C:132:ASP:C	2.62	0.42
1:G:242:TYR:O	1:G:243:GLU:C	2.61	0.42
1:B:67:LEU:HD21	1:B:159:LYS:HD2	2.01	0.42
1:E:228:ARG:HH11	1:E:228:ARG:CG	2.32	0.42
1:I:63:ASP:OD1	1:I:66:ARG:NH1	2.52	0.42
1:L:11:ARG:HG2	1:L:11:ARG:HH11	1.83	0.42
1:A:163:LEU:HG	1:A:167:LEU:CD2	2.48	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:26:LEU:HA	1:E:27:PRO:HA	1.84	0.42
1:E:242:TYR:CD1	1:F:214:TRP:HH2	2.37	0.42
1:G:74:LEU:HD13	1:G:74:LEU:C	2.45	0.42
1:I:93:ILE:CG2	1:I:95:VAL:CG2	2.97	0.42
1:I:51:THR:OG1	1:J:211:TRP:HZ3	2.02	0.42
1:B:74:LEU:HD13	1:B:74:LEU:C	2.44	0.42
1:B:106:GLN:OE1	1:B:106:GLN:HA	2.19	0.42
1:E:21:HIS:HE1	1:E:98:PRO:O	2.02	0.42
1:I:147:LEU:HG	1:J:161:LEU:HD13	2.01	0.42
1:J:210:TRP:CE2	1:J:211:TRP:HD1	2.38	0.42
1:B:115:LEU:HB3	1:B:124:THR:HG22	2.01	0.42
1:G:76:VAL:O	1:G:105:PRO:HA	2.20	0.42
1:I:111:ARG:HG3	1:I:116:LEU:HD23	2.00	0.42
1:A:5:ILE:HG22	1:A:114:GLY:HA3	2.00	0.42
1:F:210:TRP:CZ3	1:F:211:TRP:HB3	2.54	0.42
1:D:156:ARG:HD2	1:D:175:TRP:O	2.19	0.42
1:E:210:TRP:CE2	1:E:211:TRP:HD1	2.38	0.42
1:J:53:GLU:OE1	1:J:126:ARG:NH2	2.53	0.42
1:L:240:LEU:N	1:L:240:LEU:HD23	2.34	0.42
1:B:19:THR:HG22	1:B:102:ILE:HG12	2.02	0.42
1:C:226:ALA:HA	1:C:229:TYR:CD2	2.55	0.42
1:F:169:ARG:HB3	1:F:185:LEU:HB3	2.01	0.42
1:G:36:TRP:CD2	1:G:132:ASP:HA	2.54	0.42
1:H:28:ASP:O	1:H:32:SER:OG	2.35	0.42
1:C:53:GLU:OE1	1:C:126:ARG:NH2	2.54	0.41
1:D:87:GLU:O	1:D:88:TRP:C	2.63	0.41
1:D:215:ASP:C	1:D:217:PRO:HD3	2.44	0.41
1:J:156:ARG:HD2	1:J:175:TRP:O	2.20	0.41
1:B:76:VAL:O	1:B:105:PRO:HA	2.20	0.41
1:C:5:ILE:HD11	1:D:5:ILE:HG21	2.02	0.41
1:D:104:ASP:OD2	1:D:107:GLY:HA2	2.19	0.41
1:K:41:SER:HB2	1:K:124:THR:HG21	2.01	0.41
1:A:42:HIS:CD2	1:A:50:SER:HB3	2.56	0.41
1:D:119:GLU:OE2	1:D:145:MET:HG2	2.20	0.41
1:G:161:LEU:HD12	1:G:161:LEU:HA	1.90	0.41
1:G:171:VAL:HG22	1:G:185:LEU:HD22	2.01	0.41
1:K:117:HIS:HA	1:L:7:LEU:HD13	2.02	0.41
1:A:48:PRO:HG2	1:B:186:ILE:HG21	2.02	0.41
1:A:65:GLN:HA	1:A:65:GLN:NE2	2.34	0.41
1:B:72:ILE:HD11	1:B:102:ILE:CG1	2.51	0.41
1:C:85:TRP:O	1:C:86:LYS:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:106:GLN:O	1:H:111:ARG:NH2	2.53	0.41
1:K:228:ARG:O	1:K:229:TYR:C	2.62	0.41
1:L:132:ASP:OD2	1:L:134:ARG:NH1	2.54	0.41
1:F:67:LEU:HD13	1:F:158:VAL:CG2	2.51	0.41
1:I:41:SER:HB2	1:I:124:THR:CG2	2.49	0.41
1:K:87:GLU:O	1:K:91:ARG:HB2	2.21	0.41
1:B:159:LYS:HE3	1:B:163:LEU:HD21	2.03	0.41
1:E:5:ILE:O	1:F:2:PRO:HB3	2.21	0.41
1:E:58:ALA:C	1:E:60:ARG:H	2.28	0.41
1:F:189:PRO:HA	1:F:190:PRO:HD3	1.94	0.41
1:K:55:VAL:HG11	1:L:180:ILE:HD11	2.02	0.41
1:F:210:TRP:CZ2	1:K:83:ILE:CG2	3.04	0.41
1:G:88:TRP:HE3	1:G:89:ILE:N	2.17	0.41
1:I:83:ILE:HG22	1:K:210:TRP:CE2	2.56	0.41
1:A:38:VAL:O	1:A:72:ILE:HD12	2.21	0.41
1:A:43:PRO:HD2	1:A:126:ARG:NH2	2.36	0.41
1:B:26:LEU:HA	1:B:27:PRO:HA	1.85	0.41
1:D:6:PRO:HB2	1:D:137:ILE:CD1	2.51	0.41
1:F:200:MET:HA	1:F:200:MET:HE2	2.02	0.41
1:I:186:ILE:HG21	1:J:48:PRO:HG2	2.03	0.41
1:J:138:ARG:HA	1:J:138:ARG:HD2	1.88	0.41
1:K:48:PRO:HG2	1:L:186:ILE:HG21	2.03	0.41
1:L:161:LEU:HD12	1:L:161:LEU:HA	1.78	0.41
1:C:39:LEU:HA	1:C:72:ILE:HG23	2.01	0.41
1:E:138:ARG:C	1:E:139:THR:HG1	2.28	0.41
1:G:63:ASP:OD2	1:G:232:ARG:NH1	2.51	0.41
1:G:157:ILE:O	1:G:161:LEU:HB2	2.21	0.41
1:G:186:ILE:HG21	1:H:48:PRO:HG2	2.03	0.41
1:H:112:ARG:CG	1:H:112:ARG:NH1	2.83	0.40
1:H:128:VAL:O	1:H:140:MET:HA	2.21	0.40
1:I:49:VAL:CG2	1:J:170:ALA:HB1	2.49	0.40
1:C:116:LEU:HD12	1:C:116:LEU:HA	1.92	0.40
1:E:163:LEU:HD12	1:E:163:LEU:HA	1.82	0.40
1:E:190:PRO:HB3	1:E:195:GLN:CB	2.51	0.40
1:I:180:ILE:HD11	1:I:212:PHE:CZ	2.57	0.40
1:B:36:TRP:HB2	1:B:69:VAL:HG22	2.02	0.40
1:C:186:ILE:HG21	1:D:48:PRO:HG2	2.04	0.40
1:E:242:TYR:CD1	1:F:214:TRP:CH2	3.09	0.40
1:G:60:ARG:HD3	1:G:151:VAL:HG12	2.03	0.40
1:G:60:ARG:O	1:G:61:TYR:C	2.65	0.40
1:K:130:ILE:HD13	1:K:157:ILE:HG21	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:169:ARG:CB	1:K:185:LEU:HB3	2.51	0.40
1:K:223:VAL:HG12	1:K:227:ARG:NH1	2.36	0.40
1:B:226:ALA:O	1:B:227:ARG:C	2.64	0.40
1:C:41:SER:HB2	1:C:124:THR:CG2	2.51	0.40
1:D:39:LEU:C	1:D:39:LEU:CD2	2.94	0.40
1:G:208:LEU:O	1:G:209:ASP:HB2	2.21	0.40
1:G:241:LEU:HB2	1:H:180:ILE:HA	2.02	0.40
1:H:63:ASP:OD2	1:H:232:ARG:NH2	2.50	0.40
1:H:190:PRO:HG2	1:H:210:TRP:HE3	1.86	0.40
1:A:53:GLU:OE2	1:A:149:ARG:NE	2.44	0.40
1:B:33:GLN:HB3	1:B:35:LYS:HG3	2.04	0.40
1:B:161:LEU:HD12	1:B:161:LEU:HA	1.96	0.40
1:E:122:THR:HG22	1:K:106:GLN:HG3	2.04	0.40
1:I:86:LYS:HG2	1:I:97:ILE:HB	2.04	0.40
1:K:41:SER:HB2	1:K:124:THR:CG2	2.52	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	242/250 (97%)	224 (93%)	16 (7%)	2 (1%)	16	44
1	B	242/250 (97%)	227 (94%)	13 (5%)	2 (1%)	16	44
1	C	242/250 (97%)	225 (93%)	15 (6%)	2 (1%)	16	44
1	D	242/250 (97%)	224 (93%)	17 (7%)	1 (0%)	30	59
1	E	242/250 (97%)	227 (94%)	14 (6%)	1 (0%)	30	59
1	F	242/250 (97%)	227 (94%)	11 (4%)	4 (2%)	7	26
1	G	242/250 (97%)	227 (94%)	14 (6%)	1 (0%)	30	59
1	H	242/250 (97%)	221 (91%)	19 (8%)	2 (1%)	16	44

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	I	242/250 (97%)	224 (93%)	16 (7%)	2 (1%)	16	44
1	J	242/250 (97%)	221 (91%)	20 (8%)	1 (0%)	30	59
1	K	242/250 (97%)	225 (93%)	17 (7%)	0	100	100
1	L	242/250 (97%)	221 (91%)	20 (8%)	1 (0%)	30	59
All	All	2904/3000 (97%)	2693 (93%)	192 (7%)	19 (1%)	18	48

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	195	GLN
1	I	59	ARG
1	I	194	ASP
1	D	209	ASP
1	E	107	GLY
1	F	201	GLU
1	F	202	SER
1	F	209	ASP
1	F	238	ALA
1	H	209	ASP
1	H	182	GLY
1	L	125	VAL
1	A	43	PRO
1	A	125	VAL
1	G	48	PRO
1	C	182	GLY
1	B	43	PRO
1	B	125	VAL
1	J	48	PRO

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	209/215 (97%)	184 (88%)	25 (12%)	5	16

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	209/215 (97%)	187 (90%)	22 (10%)	6	22
1	C	209/215 (97%)	181 (87%)	28 (13%)	4	12
1	D	209/215 (97%)	188 (90%)	21 (10%)	7	24
1	E	209/215 (97%)	186 (89%)	23 (11%)	6	20
1	F	209/215 (97%)	184 (88%)	25 (12%)	5	16
1	G	209/215 (97%)	187 (90%)	22 (10%)	6	22
1	H	209/215 (97%)	189 (90%)	20 (10%)	8	26
1	I	209/215 (97%)	183 (88%)	26 (12%)	4	15
1	J	209/215 (97%)	192 (92%)	17 (8%)	11	33
1	K	209/215 (97%)	187 (90%)	22 (10%)	6	22
1	L	209/215 (97%)	192 (92%)	17 (8%)	11	33
All	All	2508/2580 (97%)	2240 (89%)	268 (11%)	6	22

All (268) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	THR
1	A	24	ILE
1	A	72	ILE
1	A	74	LEU
1	A	77	ASP
1	A	86	LYS
1	A	95	VAL
1	A	96	ARG
1	A	112	ARG
1	A	116	LEU
1	A	119	GLU
1	A	122	THR
1	A	131	VAL
1	A	145	MET
1	A	147	LEU
1	A	161	LEU
1	A	163	LEU
1	A	167	LEU
1	A	168	LYS
1	A	183	GLU
1	A	194	ASP
1	A	207	SER

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Mol	Chain	Res	Type
1	A	208	LEU
1	A	213	SER
1	A	242	TYR
1	B	4	SER
1	B	28	ASP
1	B	59	ARG
1	B	74	LEU
1	B	79	VAL
1	B	91	ARG
1	B	96	ARG
1	B	111	ARG
1	B	112	ARG
1	B	124	THR
1	B	146	GLU
1	B	147	LEU
1	B	161	LEU
1	B	166	SER
1	B	194	ASP
1	B	208	LEU
1	B	213	SER
1	B	224	GLU
1	B	227	ARG
1	B	235	GLU
1	B	239	LYS
1	B	244	GLU
1	C	2	PRO
1	C	4	SER
1	C	16	GLU
1	C	19	THR
1	C	74	LEU
1	C	90	GLU
1	C	108	THR
1	C	116	LEU
1	C	119	GLU
1	C	124	THR
1	C	144	PRO
1	C	145	MET
1	C	147	LEU
1	C	161	LEU
1	C	163	LEU
1	C	167	LEU
1	C	186	ILE

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Mol	Chain	Res	Type
1	C	192	THR
1	C	197	ARG
1	C	199	ARG
1	C	202	SER
1	C	208	LEU
1	C	209	ASP
1	C	213	SER
1	C	220	ARG
1	C	224	GLU
1	C	228	ARG
1	C	242	TYR
1	D	2	PRO
1	D	11	ARG
1	D	28	ASP
1	D	75	SER
1	D	79	VAL
1	D	91	ARG
1	D	96	ARG
1	D	112	ARG
1	D	147	LEU
1	D	161	LEU
1	D	174	ASP
1	D	191	THR
1	D	192	THR
1	D	194	ASP
1	D	199	ARG
1	D	202	SER
1	D	208	LEU
1	D	225	GLU
1	D	242	TYR
1	D	243	GLU
1	D	244	GLU
1	E	59	ARG
1	E	62	GLU
1	E	74	LEU
1	E	78	SER
1	E	79	VAL
1	E	91	ARG
1	E	108	THR
1	E	116	LEU
1	E	124	THR
1	E	147	LEU

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Mol	Chain	Res	Type
1	E	161	LEU
1	E	163	LEU
1	E	167	LEU
1	E	174	ASP
1	E	179	GLU
1	E	188	PRO
1	E	194	ASP
1	E	199	ARG
1	E	208	LEU
1	E	213	SER
1	E	228	ARG
1	E	235	GLU
1	E	242	TYR
1	F	2	PRO
1	F	4	SER
1	F	24	ILE
1	F	59	ARG
1	F	66	ARG
1	F	67	LEU
1	F	74	LEU
1	F	75	SER
1	F	78	SER
1	F	79	VAL
1	F	86	LYS
1	F	91	ARG
1	F	106	GLN
1	F	111	ARG
1	F	112	ARG
1	F	147	LEU
1	F	161	LEU
1	F	176	PRO
1	F	181	ILE
1	F	194	ASP
1	F	199	ARG
1	F	208	LEU
1	F	213	SER
1	F	220	ARG
1	F	227	ARG
1	G	11	ARG
1	G	16	GLU
1	G	32	SER
1	G	72	ILE

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Mol	Chain	Res	Type
1	G	79	VAL
1	G	91	ARG
1	G	112	ARG
1	G	116	LEU
1	G	124	THR
1	G	147	LEU
1	G	161	LEU
1	G	163	LEU
1	G	167	LEU
1	G	193	GLU
1	G	199	ARG
1	G	206	ARG
1	G	208	LEU
1	G	220	ARG
1	G	221	ASP
1	G	225	GLU
1	G	227	ARG
1	G	239	LYS
1	H	2	PRO
1	H	4	SER
1	H	28	ASP
1	H	32	SER
1	H	79	VAL
1	H	93	ILE
1	H	96	ARG
1	H	112	ARG
1	H	147	LEU
1	H	161	LEU
1	H	165	ASP
1	H	176	PRO
1	H	186	ILE
1	H	193	GLU
1	H	199	ARG
1	H	208	LEU
1	H	213	SER
1	H	220	ARG
1	H	227	ARG
1	H	242	TYR
1	I	5	ILE
1	I	18	THR
1	I	19	THR
1	I	24	ILE

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Mol	Chain	Res	Type
1	I	59	ARG
1	I	62	GLU
1	I	79	VAL
1	I	86	LYS
1	I	87	GLU
1	I	93	ILE
1	I	96	ARG
1	I	116	LEU
1	I	124	THR
1	I	141	LEU
1	I	145	MET
1	I	147	LEU
1	I	159	LYS
1	I	163	LEU
1	I	167	LEU
1	I	179	GLU
1	I	180	ILE
1	I	185	LEU
1	I	192	THR
1	I	199	ARG
1	I	208	LEU
1	I	242	TYR
1	J	2	PRO
1	J	14	GLU
1	J	33	GLN
1	J	74	LEU
1	J	75	SER
1	J	76	VAL
1	J	78	SER
1	J	79	VAL
1	J	91	ARG
1	J	93	ILE
1	J	147	LEU
1	J	161	LEU
1	J	199	ARG
1	J	213	SER
1	J	220	ARG
1	J	227	ARG
1	J	235	GLU
1	K	11	ARG
1	K	32	SER
1	K	62	GLU

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Mol	Chain	Res	Type
1	K	66	ARG
1	K	67	LEU
1	K	79	VAL
1	K	91	ARG
1	K	96	ARG
1	K	112	ARG
1	K	116	LEU
1	K	124	THR
1	K	147	LEU
1	K	161	LEU
1	K	168	LYS
1	K	199	ARG
1	K	211	TRP
1	K	213	SER
1	K	219	SER
1	K	228	ARG
1	K	230	LEU
1	K	239	LYS
1	K	244	GLU
1	L	28	ASP
1	L	59	ARG
1	L	75	SER
1	L	124	THR
1	L	147	LEU
1	L	161	LEU
1	L	179	GLU
1	L	193	GLU
1	L	197	ARG
1	L	199	ARG
1	L	206	ARG
1	L	208	LEU
1	L	220	ARG
1	L	225	GLU
1	L	227	ARG
1	L	242	TYR
1	L	243	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	92	HIS

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Mol	Chain	Res	Type
1	B	33	GLN
1	C	29	HIS
1	C	33	GLN
1	C	195	GLN
1	D	65	GLN
1	E	21	HIS
1	E	33	GLN
1	E	42	HIS
1	E	92	HIS
1	E	204	GLN
1	F	33	GLN
1	F	42	HIS
1	F	65	GLN
1	F	123	HIS
1	F	204	GLN
1	G	195	GLN
1	H	42	HIS
1	H	65	GLN
1	H	195	GLN
1	I	42	HIS
1	I	65	GLN
1	I	204	GLN
1	J	29	HIS
1	J	33	GLN
1	J	195	GLN
1	J	204	GLN
1	K	42	HIS
1	K	204	GLN
1	L	21	HIS
1	L	65	GLN
1	L	204	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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	244/250 (97%)	-0.39	2 (0%) 82 77	41, 59, 88, 111	0
1	B	244/250 (97%)	-0.50	1 (0%) 88 85	39, 56, 82, 119	0
1	C	244/250 (97%)	-0.60	1 (0%) 88 85	35, 50, 75, 108	0
1	D	244/250 (97%)	-0.66	1 (0%) 88 85	33, 48, 76, 120	0
1	E	244/250 (97%)	-0.63	0 100 100	33, 47, 76, 105	0
1	F	244/250 (97%)	-0.62	1 (0%) 88 85	30, 46, 81, 109	0
1	G	244/250 (97%)	-0.67	0 100 100	32, 46, 71, 96	0
1	H	244/250 (97%)	-0.59	1 (0%) 88 85	30, 49, 77, 108	0
1	I	244/250 (97%)	-0.51	1 (0%) 88 85	40, 56, 81, 107	0
1	J	244/250 (97%)	-0.49	1 (0%) 88 85	39, 57, 87, 112	0
1	K	244/250 (97%)	-0.47	1 (0%) 88 85	36, 58, 88, 116	0
1	L	244/250 (97%)	-0.50	1 (0%) 88 85	41, 57, 82, 119	0
All	All	2928/3000 (97%)	-0.55	11 (0%) 88 85	30, 53, 82, 120	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	245	ALA	4.3
1	K	245	ALA	3.3
1	H	245	ALA	3.2
1	B	245	ALA	3.1
1	J	245	ALA	3.0
1	A	245	ALA	2.7
1	F	245	ALA	2.7
1	C	245	ALA	2.6
1	D	245	ALA	2.5
1	L	245	ALA	2.2
1	A	25	LYS	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.