



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:21 PM UTC

PDB ID : 8KBB / pdb_00008kbb
Title : Structure of apo-CmTad1
Authors : Xiao, Y.; Feng, Y.
Deposited on : 2023-08-04
Resolution : 2.56 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:

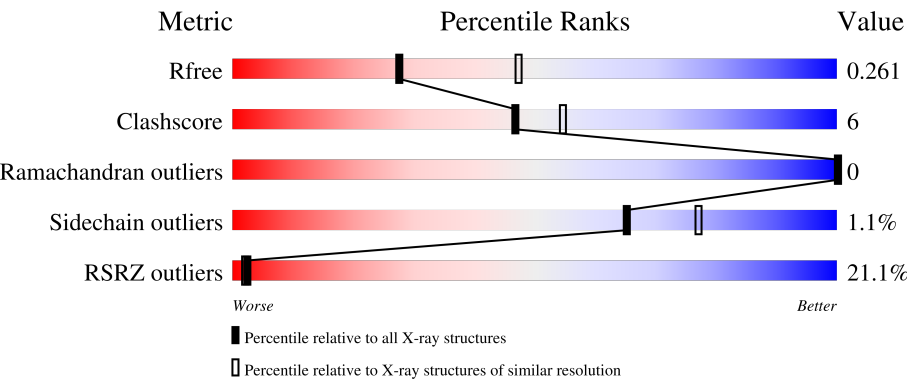
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.56 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	1853 (2.58-2.54)
Clashscore	190562	1897 (2.58-2.54)
Ramachandran outliers	187476	1875 (2.58-2.54)
Sidechain outliers	187428	1875 (2.58-2.54)
RSRZ outliers	180081	1853 (2.58-2.54)

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	132	14% 77% 18% 5%
1	B	132	18% 84% 16%
1	C	132	11% 88% 12%
1	D	132	15% 86% 11% ..
1	E	132	20% 85% 13% ..

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Mol	Chain	Length	Quality of chain
1	F	132	18% 86% 14%
1	G	132	14% 83% 14% ••
1	H	132	21% 82% 16% •
1	I	132	28% 80% 18% •
1	J	132	35% 78% 20% •
1	K	132	25% 78% 20% •
1	L	132	29% 82% 14% 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 13140 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 Thoeris anti-defense 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	126	Total	C	N	O	S	0	0	0
			1021	649	173	191	8			
1	B	132	Total	C	N	O	S	0	0	0
			1064	673	179	204	8			
1	G	131	Total	C	N	O	S	0	0	0
			1058	670	178	202	8			
1	H	129	Total	C	N	O	S	0	0	0
			1040	659	176	197	8			
1	I	129	Total	C	N	O	S	0	0	0
			1043	661	176	198	8			
1	J	129	Total	C	N	O	S	0	0	0
			1043	661	176	198	8			
1	K	130	Total	C	N	O	S	0	0	0
			1052	667	177	200	8			
1	L	126	Total	C	N	O	S	0	0	0
			1024	651	173	192	8			
1	C	132	Total	C	N	O	S	0	0	0
			1064	673	179	204	8			
1	D	131	Total	C	N	O	S	0	0	0
			1058	670	178	202	8			
1	E	130	Total	C	N	O	S	0	0	0
			1052	667	177	200	8			
1	F	132	Total	C	N	O	S	0	0	0
			1064	673	179	204	8			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-9	SER	-	expression tag	UNP P0DW61
A	-8	SER	-	expression tag	UNP P0DW61
A	-7	GLY	-	expression tag	UNP P0DW61
A	-6	LEU	-	expression tag	UNP P0DW61
A	-5	VAL	-	expression tag	UNP P0DW61

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	PRO	-	expression tag	UNP P0DW61
A	-3	ARG	-	expression tag	UNP P0DW61
A	-2	GLY	-	expression tag	UNP P0DW61
A	-1	SER	-	expression tag	UNP P0DW61
A	0	HIS	-	expression tag	UNP P0DW61
B	-9	SER	-	expression tag	UNP P0DW61
B	-8	SER	-	expression tag	UNP P0DW61
B	-7	GLY	-	expression tag	UNP P0DW61
B	-6	LEU	-	expression tag	UNP P0DW61
B	-5	VAL	-	expression tag	UNP P0DW61
B	-4	PRO	-	expression tag	UNP P0DW61
B	-3	ARG	-	expression tag	UNP P0DW61
B	-2	GLY	-	expression tag	UNP P0DW61
B	-1	SER	-	expression tag	UNP P0DW61
B	0	HIS	-	expression tag	UNP P0DW61
G	-9	SER	-	expression tag	UNP P0DW61
G	-8	SER	-	expression tag	UNP P0DW61
G	-7	GLY	-	expression tag	UNP P0DW61
G	-6	LEU	-	expression tag	UNP P0DW61
G	-5	VAL	-	expression tag	UNP P0DW61
G	-4	PRO	-	expression tag	UNP P0DW61
G	-3	ARG	-	expression tag	UNP P0DW61
G	-2	GLY	-	expression tag	UNP P0DW61
G	-1	SER	-	expression tag	UNP P0DW61
G	0	HIS	-	expression tag	UNP P0DW61
H	-9	SER	-	expression tag	UNP P0DW61
H	-8	SER	-	expression tag	UNP P0DW61
H	-7	GLY	-	expression tag	UNP P0DW61
H	-6	LEU	-	expression tag	UNP P0DW61
H	-5	VAL	-	expression tag	UNP P0DW61
H	-4	PRO	-	expression tag	UNP P0DW61
H	-3	ARG	-	expression tag	UNP P0DW61
H	-2	GLY	-	expression tag	UNP P0DW61
H	-1	SER	-	expression tag	UNP P0DW61
H	0	HIS	-	expression tag	UNP P0DW61
I	-9	SER	-	expression tag	UNP P0DW61
I	-8	SER	-	expression tag	UNP P0DW61
I	-7	GLY	-	expression tag	UNP P0DW61
I	-6	LEU	-	expression tag	UNP P0DW61
I	-5	VAL	-	expression tag	UNP P0DW61
I	-4	PRO	-	expression tag	UNP P0DW61
I	-3	ARG	-	expression tag	UNP P0DW61

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Chain	Residue	Modelled	Actual	Comment	Reference
I	-2	GLY	-	expression tag	UNP P0DW61
I	-1	SER	-	expression tag	UNP P0DW61
I	0	HIS	-	expression tag	UNP P0DW61
J	-9	SER	-	expression tag	UNP P0DW61
J	-8	SER	-	expression tag	UNP P0DW61
J	-7	GLY	-	expression tag	UNP P0DW61
J	-6	LEU	-	expression tag	UNP P0DW61
J	-5	VAL	-	expression tag	UNP P0DW61
J	-4	PRO	-	expression tag	UNP P0DW61
J	-3	ARG	-	expression tag	UNP P0DW61
J	-2	GLY	-	expression tag	UNP P0DW61
J	-1	SER	-	expression tag	UNP P0DW61
J	0	HIS	-	expression tag	UNP P0DW61
K	-9	SER	-	expression tag	UNP P0DW61
K	-8	SER	-	expression tag	UNP P0DW61
K	-7	GLY	-	expression tag	UNP P0DW61
K	-6	LEU	-	expression tag	UNP P0DW61
K	-5	VAL	-	expression tag	UNP P0DW61
K	-4	PRO	-	expression tag	UNP P0DW61
K	-3	ARG	-	expression tag	UNP P0DW61
K	-2	GLY	-	expression tag	UNP P0DW61
K	-1	SER	-	expression tag	UNP P0DW61
K	0	HIS	-	expression tag	UNP P0DW61
L	-9	SER	-	expression tag	UNP P0DW61
L	-8	SER	-	expression tag	UNP P0DW61
L	-7	GLY	-	expression tag	UNP P0DW61
L	-6	LEU	-	expression tag	UNP P0DW61
L	-5	VAL	-	expression tag	UNP P0DW61
L	-4	PRO	-	expression tag	UNP P0DW61
L	-3	ARG	-	expression tag	UNP P0DW61
L	-2	GLY	-	expression tag	UNP P0DW61
L	-1	SER	-	expression tag	UNP P0DW61
L	0	HIS	-	expression tag	UNP P0DW61
C	-9	SER	-	expression tag	UNP P0DW61
C	-8	SER	-	expression tag	UNP P0DW61
C	-7	GLY	-	expression tag	UNP P0DW61
C	-6	LEU	-	expression tag	UNP P0DW61
C	-5	VAL	-	expression tag	UNP P0DW61
C	-4	PRO	-	expression tag	UNP P0DW61
C	-3	ARG	-	expression tag	UNP P0DW61
C	-2	GLY	-	expression tag	UNP P0DW61
C	-1	SER	-	expression tag	UNP P0DW61

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Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	expression tag	UNP P0DW61
D	-9	SER	-	expression tag	UNP P0DW61
D	-8	SER	-	expression tag	UNP P0DW61
D	-7	GLY	-	expression tag	UNP P0DW61
D	-6	LEU	-	expression tag	UNP P0DW61
D	-5	VAL	-	expression tag	UNP P0DW61
D	-4	PRO	-	expression tag	UNP P0DW61
D	-3	ARG	-	expression tag	UNP P0DW61
D	-2	GLY	-	expression tag	UNP P0DW61
D	-1	SER	-	expression tag	UNP P0DW61
D	0	HIS	-	expression tag	UNP P0DW61
E	-9	SER	-	expression tag	UNP P0DW61
E	-8	SER	-	expression tag	UNP P0DW61
E	-7	GLY	-	expression tag	UNP P0DW61
E	-6	LEU	-	expression tag	UNP P0DW61
E	-5	VAL	-	expression tag	UNP P0DW61
E	-4	PRO	-	expression tag	UNP P0DW61
E	-3	ARG	-	expression tag	UNP P0DW61
E	-2	GLY	-	expression tag	UNP P0DW61
E	-1	SER	-	expression tag	UNP P0DW61
E	0	HIS	-	expression tag	UNP P0DW61
F	-9	SER	-	expression tag	UNP P0DW61
F	-8	SER	-	expression tag	UNP P0DW61
F	-7	GLY	-	expression tag	UNP P0DW61
F	-6	LEU	-	expression tag	UNP P0DW61
F	-5	VAL	-	expression tag	UNP P0DW61
F	-4	PRO	-	expression tag	UNP P0DW61
F	-3	ARG	-	expression tag	UNP P0DW61
F	-2	GLY	-	expression tag	UNP P0DW61
F	-1	SER	-	expression tag	UNP P0DW61
F	0	HIS	-	expression tag	UNP P0DW61

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0

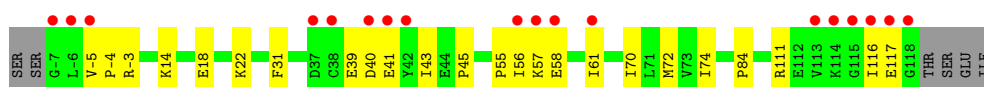
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	56	Total 56	O 56	0	0
3	B	51	Total 51	O 51	0	0
3	G	49	Total 49	O 49	0	0
3	H	57	Total 57	O 57	0	0
3	I	33	Total 33	O 33	0	0
3	J	28	Total 28	O 28	0	0
3	K	30	Total 30	O 30	0	0
3	L	30	Total 30	O 30	0	0
3	C	58	Total 58	O 58	0	0
3	D	59	Total 59	O 59	0	0
3	E	46	Total 46	O 46	0	0
3	F	58	Total 58	O 58	0	0

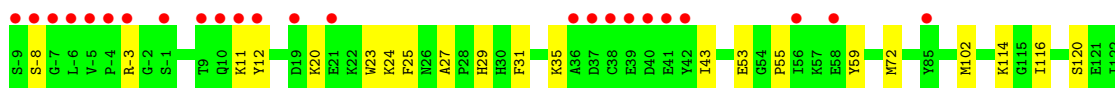
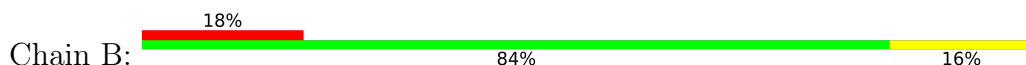
3 Residue-property plots [i](#)

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.

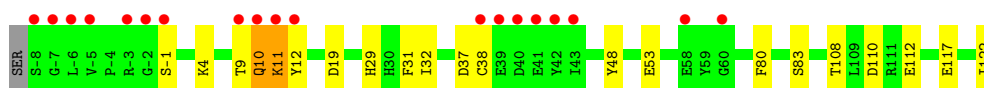
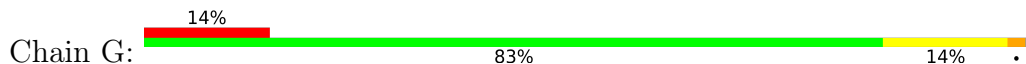
- Molecule 1: Thoeris anti-defense 1



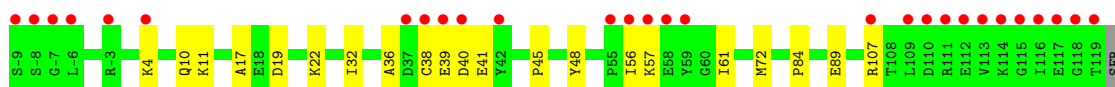
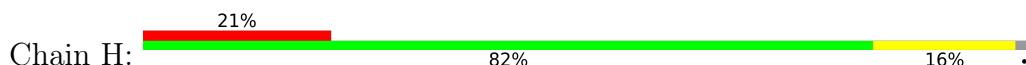
- Molecule 1: Thoeris anti-defense 1



- Molecule 1: Thoeris anti-defense 1

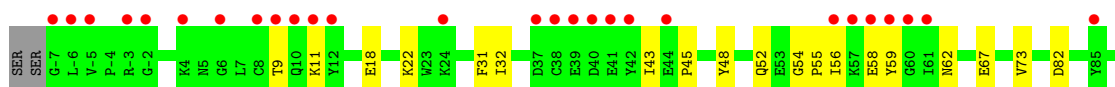


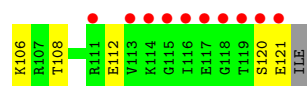
- Molecule 1: Thoeris anti-defense 1



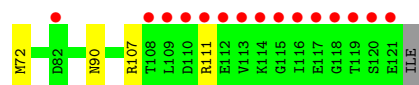
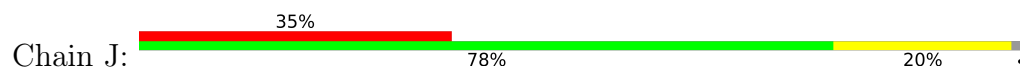
GLU
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- Molecule 1: Thoeris anti-defense 1

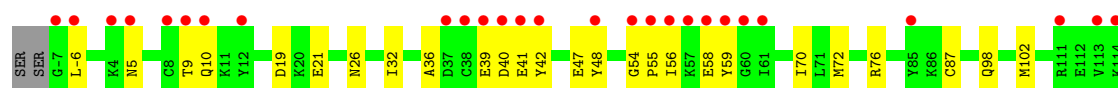
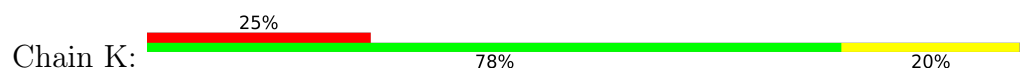




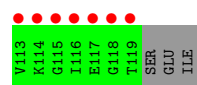
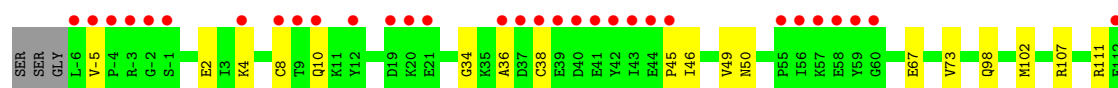
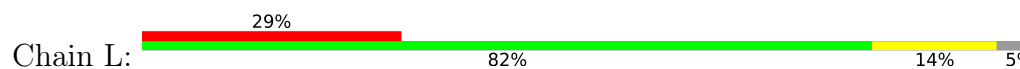
● Molecule 1: Thoeris anti-defense 1



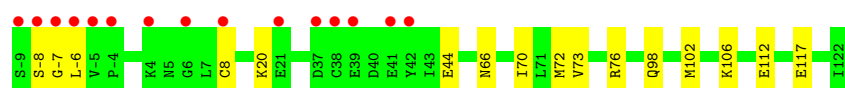
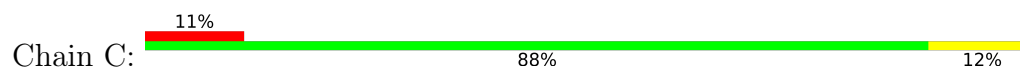
● Molecule 1: Thoeris anti-defense 1



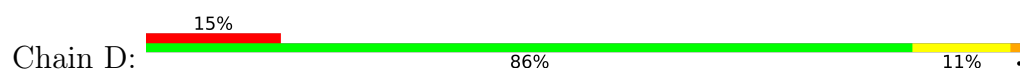
● Molecule 1: Thoeris anti-defense 1



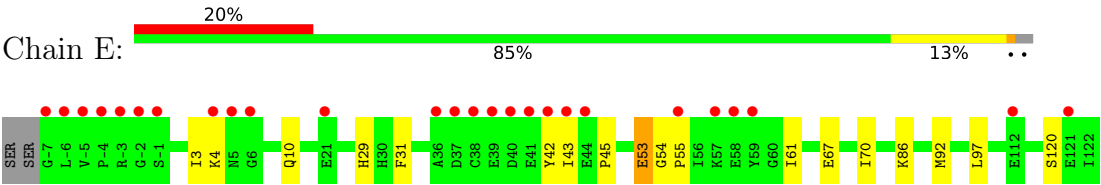
● Molecule 1: Thoeris anti-defense 1



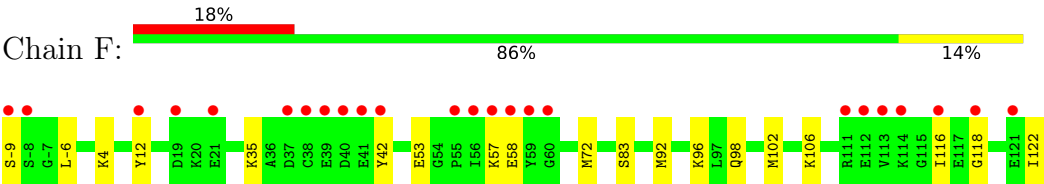
● Molecule 1: Thoeris anti-defense 1



● Molecule 1: Thoeris anti-defense 1



● Molecule 1: Thoeris anti-defense 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	138.85Å 144.52Å 150.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.13 – 2.56 41.13 – 2.56	Depositor EDS
% Data completeness (in resolution range)	99.2 (41.13-2.56) 99.3 (41.13-2.56)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.54Å)	Xtriage
Refinement program	PHENIX (1.18.2_3874: ???)	Depositor
R, R_{free}	0.227 , 0.263 0.229 , 0.261	Depositor DCC
R_{free} test set	2014 reflections (2.06%)	wwPDB-VP
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.224	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	13140	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.74% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.61	1/1042 (0.1%)	0.64	0/1402
1	B	0.61	2/1085 (0.2%)	0.69	1/1459 (0.1%)
1	C	0.74	0/1085	0.78	4/1459 (0.3%)
1	D	0.79	0/1079	0.80	5/1451 (0.3%)
1	E	0.82	2/1073 (0.2%)	0.68	1/1443 (0.1%)
1	F	0.82	1/1085 (0.1%)	0.73	1/1459 (0.1%)
1	G	0.66	1/1079 (0.1%)	0.81	4/1451 (0.3%)
1	H	0.97	3/1061 (0.3%)	0.82	4/1428 (0.3%)
1	I	0.68	1/1064 (0.1%)	0.72	0/1432
1	J	0.46	0/1064	0.67	2/1432 (0.1%)
1	K	0.91	0/1073	0.95	6/1443 (0.4%)
1	L	0.77	0/1045	0.78	5/1407 (0.4%)
All	All	0.75	11/12835 (0.1%)	0.76	33/17266 (0.2%)

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	27	ALA	CA-C	-8.12	1.44	1.52
1	H	22	LYS	C-O	-6.04	1.16	1.24
1	F	53	GLU	CA-C	-5.79	1.45	1.52
1	G	37	ASP	CA-C	-5.75	1.47	1.53
1	E	53	GLU	CA-C	-5.66	1.45	1.52
1	B	27	ALA	C-O	5.31	1.28	1.24
1	A	22	LYS	C-O	-5.27	1.16	1.24
1	H	19	ASP	CA-C	-5.20	1.46	1.52
1	I	22	LYS	C-O	-5.15	1.17	1.24
1	H	38	CYS	CA-C	-5.03	1.46	1.52
1	E	3	ILE	N-CA	-5.01	1.40	1.46

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	9	THR	N-CA-C	12.92	125.36	111.28
1	G	9	THR	N-CA-C	12.43	124.83	111.28
1	H	38	CYS	N-CA-C	-10.48	93.14	109.07
1	D	40	ASP	N-CA-C	10.33	122.54	111.28
1	B	27	ALA	O-C-N	-9.68	115.51	121.71
1	C	-8	SER	N-CA-C	-9.64	99.88	112.41
1	D	27	ALA	N-CA-C	9.38	119.13	108.25
1	L	38	CYS	N-CA-C	-8.92	95.67	108.96
1	H	45	PRO	N-CA-C	8.37	123.64	110.50
1	F	116	ILE	N-CA-C	8.23	120.03	112.29
1	K	116	ILE	N-CA-C	8.10	119.90	112.29
1	J	38	CYS	N-CA-C	-7.56	97.56	108.60
1	K	56	ILE	N-CA-C	7.25	117.97	110.36
1	L	36	ALA	N-CA-C	7.22	119.23	111.36
1	G	38	CYS	N-CA-C	-6.61	98.98	109.23
1	K	10	GLN	N-CA-C	6.34	120.36	112.24
1	H	36	ALA	N-CA-C	6.13	117.97	111.28
1	G	10	GLN	N-CA-C	6.04	120.60	113.23
1	J	36	ALA	N-CA-C	6.03	117.94	111.36
1	L	-5	VAL	CA-C-N	-5.90	113.89	119.85
1	L	-5	VAL	C-N-CA	-5.90	113.89	119.85
1	D	42	TYR	N-CA-C	5.84	118.57	109.52
1	K	40	ASP	N-CA-C	5.82	118.40	111.71
1	D	17	ALA	N-CA-C	-5.55	101.04	108.86
1	G	11	LYS	N-CA-C	5.54	117.40	111.36
1	C	20	LYS	CA-C-N	5.50	127.91	120.38
1	C	20	LYS	C-N-CA	5.50	127.91	120.38
1	H	17	ALA	N-CA-C	-5.41	101.23	108.86
1	E	42	TYR	N-CA-C	5.41	118.04	109.50
1	D	36	ALA	N-CA-C	5.38	117.22	111.36
1	L	45	PRO	N-CA-C	5.38	119.94	111.38
1	C	-7	GLY	N-CA-C	-5.30	107.90	115.64
1	K	36	ALA	N-CA-C	5.08	116.82	111.28

There are no chirality outliers.

There are no planarity outliers.

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	1021	0	1012	16	0
1	B	1064	0	1051	12	0
1	C	1064	0	1051	12	0
1	D	1058	0	1046	12	0
1	E	1052	0	1041	13	0
1	F	1064	0	1051	14	0
1	G	1058	0	1046	12	0
1	H	1040	0	1029	10	0
1	I	1043	0	1030	19	0
1	J	1043	0	1030	20	0
1	K	1052	0	1041	16	0
1	L	1024	0	1016	11	0
2	A	1	0	0	0	0
2	D	1	0	0	0	0
3	A	56	0	0	2	0
3	B	51	0	0	1	0
3	C	58	0	0	1	0
3	D	59	0	0	2	0
3	E	46	0	0	3	0
3	F	58	0	0	2	0
3	G	49	0	0	1	0
3	H	57	0	0	1	0
3	I	33	0	0	4	0
3	J	28	0	0	5	0
3	K	30	0	0	2	0
3	L	30	0	0	6	0
All	All	13140	0	12444	143	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 6.

All (143) 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:F:4:LYS:NZ	3:F:201:HOH:O	2.02	0.93
1:K:59:TYR:O	3:K:201:HOH:O	1.91	0.89
1:J:10:GLN:NE2	3:J:201:HOH:O	2.08	0.85
1:L:67:GLU:OE2	3:L:202:HOH:O	2.03	0.76
1:A:111:ARG:HE	1:A:117:GLU:HG2	1.52	0.75
1:L:50:ASN:N	3:L:201:HOH:O	1.94	0.70
1:E:61:ILE:HD11	1:F:12:TYR:HB3	1.75	0.69
1:F:-9:SER:N	3:F:202:HOH:O	2.28	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:110:ASP:O	3:G:201:HOH:O	2.12	0.67
1:B:59:TYR:O	3:B:201:HOH:O	2.11	0.67
1:H:4:LYS:NZ	1:H:10:GLN:OE1	2.21	0.66
1:C:72:MET:HG2	1:D:31:PHE:CD2	2.31	0.65
1:A:84:PRO:O	3:A:301:HOH:O	2.14	0.65
1:D:5:ASN:ND2	3:D:302:HOH:O	2.29	0.65
1:J:67:GLU:OE2	3:J:202:HOH:O	2.14	0.64
1:I:59:TYR:OH	3:I:201:HOH:O	2.01	0.63
1:D:2:GLU:OE1	1:D:4:LYS:HD3	1.99	0.63
1:F:-6:LEU:HD12	1:F:42:TYR:HB3	1.82	0.62
1:J:19:ASP:OD1	1:J:20:LYS:N	2.33	0.61
1:J:90:ASN:OD1	3:J:203:HOH:O	2.15	0.61
1:K:-6:LEU:HD22	1:K:42:TYR:HB3	1.82	0.61
1:A:61:ILE:HB	1:B:35:LYS:HE3	1.83	0.61
1:A:14:LYS:NZ	3:A:303:HOH:O	2.28	0.61
1:B:114:LYS:HB2	1:B:116:ILE:HG12	1.82	0.61
1:E:31:PHE:CD2	1:F:72:MET:HG2	2.35	0.60
1:I:56:ILE:H	1:I:56:ILE:HD12	1.67	0.60
1:A:-5:VAL:HG22	1:A:-4:PRO:HD2	1.83	0.59
1:G:122:ILE:HD11	1:H:84:PRO:HB2	1.84	0.59
1:C:98:GLN:O	1:C:102:MET:HG3	2.03	0.59
1:A:43:ILE:O	1:A:45:PRO:HD3	2.02	0.59
1:J:107:ARG:HH22	1:J:111:ARG:NH1	2.00	0.59
1:A:31:PHE:CD2	1:B:72:MET:HG2	2.38	0.58
1:D:4:LYS:HD2	1:D:10:GLN:HE22	1.69	0.58
3:C:242:HOH:O	1:E:92:MET:SD	2.57	0.58
1:A:111:ARG:HB3	1:A:116:ILE:O	2.04	0.58
1:L:4:LYS:HE2	1:L:10:GLN:OE1	2.04	0.58
1:B:102:MET:HE2	1:I:106:LYS:HB2	1.87	0.56
1:F:98:GLN:O	1:F:102:MET:HG3	2.06	0.56
1:L:107:ARG:O	1:L:111:ARG:HG3	2.07	0.55
1:D:5:ASN:ND2	3:D:301:HOH:O	2.12	0.55
1:K:58:GLU:HG3	1:K:59:TYR:HD1	1.71	0.55
1:A:-3:ARG:HG3	1:A:43:ILE:HG13	1.89	0.54
1:G:-1:SER:HB3	1:G:19:ASP:HB2	1.88	0.54
1:A:72:MET:HG2	1:B:31:PHE:CD2	2.44	0.53
1:G:122:ILE:HD12	1:F:83:SER:C	2.35	0.52
1:I:120:SER:O	1:I:121:GLU:HB2	2.09	0.52
1:K:98:GLN:O	1:K:102:MET:HG3	2.09	0.52
1:C:73:VAL:HG11	1:D:70:ILE:HD11	1.92	0.52
1:A:56:ILE:H	1:A:56:ILE:HD12	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:4:LYS:HG2	1:G:10:GLN:HE22	1.75	0.51
1:J:39:GLU:HB2	1:J:42:TYR:HB2	1.90	0.51
1:L:34:GLY:HA2	1:L:46:ILE:HG13	1.91	0.51
1:I:9:THR:HG23	1:I:11:LYS:H	1.73	0.51
1:K:26:ASN:ND2	3:K:203:HOH:O	2.27	0.51
1:G:108:THR:O	1:G:112:GLU:HG2	2.11	0.50
1:H:89:GLU:OE2	1:H:89:GLU:N	2.38	0.50
1:A:-3:ARG:HD3	1:A:40:ASP:O	2.11	0.50
1:J:18:GLU:OE1	3:J:204:HOH:O	2.19	0.50
1:J:29:HIS:CG	1:J:53:GLU:HG2	2.47	0.50
1:K:32:ILE:HG12	1:K:48:TYR:CE2	2.47	0.50
1:F:122:ILE:H	1:F:122:ILE:HD12	1.77	0.49
1:J:62:ASN:O	3:J:205:HOH:O	2.19	0.49
1:E:86:LYS:NZ	3:E:207:HOH:O	2.44	0.49
1:K:39:GLU:HB2	1:K:42:TYR:HD2	1.78	0.49
1:I:55:PRO:HD2	1:I:59:TYR:CE2	2.48	0.49
1:G:31:PHE:CD2	1:H:72:MET:HG2	2.47	0.49
1:E:29:HIS:CG	1:E:53:GLU:HG2	2.48	0.49
1:L:8:CYS:HA	3:L:211:HOH:O	2.12	0.49
1:I:67:GLU:OE2	3:I:203:HOH:O	2.20	0.48
1:I:73:VAL:HG11	1:J:70:ILE:HD11	1.94	0.48
1:K:87:CYS:SG	1:L:107:ARG:HD2	2.54	0.48
1:H:32:ILE:HG12	1:H:48:TYR:CD1	2.48	0.48
1:I:54:GLY:HA3	1:I:59:TYR:CD2	2.48	0.48
1:C:72:MET:O	1:C:76:ARG:HG3	2.13	0.47
1:I:82:ASP:OD2	3:I:202:HOH:O	2.20	0.47
1:J:57:LYS:HD2	1:J:57:LYS:N	2.29	0.47
1:C:102:MET:HE2	1:F:106:LYS:HB2	1.96	0.47
1:E:67:GLU:OE1	3:E:202:HOH:O	2.19	0.47
1:H:56:ILE:H	1:H:56:ILE:HD12	1.79	0.47
1:I:31:PHE:CD2	1:J:72:MET:HG2	2.50	0.47
1:B:-3:ARG:HA	1:B:43:ILE:HD12	1.97	0.46
1:E:10:GLN:N	3:E:201:HOH:O	2.11	0.46
1:K:47:GLU:CD	3:L:201:HOH:O	2.58	0.46
1:J:1:MET:HE1	1:J:23:TRP:CZ2	2.51	0.46
1:L:98:GLN:O	1:L:102:MET:HG3	2.15	0.46
1:D:19:ASP:HB3	1:D:22:LYS:HD3	1.96	0.46
1:I:32:ILE:HG12	1:I:48:TYR:CE1	2.51	0.46
1:G:117:GLU:OE1	1:H:11:LYS:NZ	2.45	0.46
1:E:43:ILE:O	1:E:45:PRO:HD3	2.15	0.46
1:K:70:ILE:HD11	1:L:73:VAL:HG11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:108:THR:O	1:I:112:GLU:HG3	2.17	0.45
1:B:24:LYS:HB3	1:B:25:PHE:CD2	2.52	0.45
1:B:29:HIS:CD2	1:B:53:GLU:HA	2.52	0.45
1:E:61:ILE:HG21	1:F:35:LYS:HG2	1.98	0.45
1:A:55:PRO:HG2	1:A:58:GLU:HB2	1.98	0.45
1:E:70:ILE:HG23	1:E:97:LEU:HD22	1.99	0.45
1:I:58:GLU:C	3:I:204:HOH:O	2.60	0.44
1:A:39:GLU:C	1:A:41:GLU:H	2.24	0.44
1:I:31:PHE:CG	1:J:72:MET:HG2	2.52	0.44
1:K:72:MET:O	1:K:76:ARG:HG3	2.18	0.44
1:C:8:CYS:SG	1:D:71:LEU:HD22	2.57	0.44
1:B:11:LYS:HE2	1:B:12:TYR:CZ	2.53	0.44
1:C:66:ASN:O	1:C:70:ILE:HG12	2.18	0.44
1:C:102:MET:HB3	1:F:102:MET:HB3	2.00	0.44
1:D:11:LYS:HD3	1:D:12:TYR:CE2	2.52	0.44
1:J:2:GLU:HG2	1:J:4:LYS:HG3	2.00	0.44
1:L:2:GLU:OE2	1:L:4:LYS:HE3	2.18	0.44
1:I:32:ILE:HG12	1:I:48:TYR:CD1	2.53	0.43
1:C:112:GLU:HG3	1:C:117:GLU:HG3	2.00	0.43
1:B:55:PRO:HG2	1:B:120:SER:HB3	2.00	0.43
1:G:29:HIS:CG	1:G:53:GLU:HG2	2.52	0.43
1:G:80:PHE:O	1:G:83:SER:OG	2.24	0.43
1:C:106:LYS:HB2	1:F:102:MET:HE2	2.00	0.43
1:H:107:ARG:NH2	3:H:205:HOH:O	2.50	0.43
1:I:55:PRO:HD2	1:I:59:TYR:CD2	2.54	0.43
1:J:57:LYS:HD2	1:J:57:LYS:H	1.84	0.43
1:G:12:TYR:HB3	1:H:61:ILE:HD11	2.00	0.43
1:C:72:MET:HG2	1:D:31:PHE:CG	2.54	0.43
1:K:47:GLU:HG2	3:L:201:HOH:O	2.19	0.42
1:J:107:ARG:NH2	1:J:111:ARG:NH1	2.65	0.42
1:A:70:ILE:O	1:A:74:ILE:HG13	2.19	0.42
1:D:9:THR:HG23	1:D:11:LYS:H	1.85	0.42
1:E:55:PRO:HG2	1:E:120:SER:HB3	2.00	0.42
1:F:92:MET:O	1:F:96:LYS:HG2	2.20	0.42
1:I:43:ILE:HG22	1:I:45:PRO:HD3	2.01	0.42
1:J:19:ASP:HB3	1:J:22:LYS:HG3	2.02	0.42
1:K:54:GLY:HA2	1:K:55:PRO:HD3	1.94	0.42
1:K:32:ILE:HG12	1:K:48:TYR:CD2	2.55	0.42
1:H:57:LYS:HE2	1:H:57:LYS:HA	2.02	0.41
1:I:52:GLN:HA	1:I:62:ASN:OD1	2.20	0.41
1:G:32:ILE:HG12	1:G:48:TYR:CE2	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:54:GLY:HA2	1:E:55:PRO:HD3	1.91	0.41
1:C:-6:LEU:HD12	1:C:44:GLU:HG3	2.03	0.41
1:D:18:GLU:H	1:D:18:GLU:HG3	1.25	0.41
1:B:20:LYS:HA	1:B:23:TRP:CE2	2.56	0.41
1:K:58:GLU:HG3	1:K:59:TYR:CD1	2.51	0.41
1:E:67:GLU:OE1	1:E:67:GLU:N	2.50	0.41
1:J:23:TRP:NE1	1:J:28:PRO:HB3	2.36	0.40
1:K:19:ASP:OD1	1:K:21:GLU:HG2	2.21	0.40
1:L:49:VAL:HA	3:L:201:HOH:O	2.21	0.40
1:A:57:LYS:HA	1:A:57:LYS:HE2	2.03	0.40
1:J:-1:SER:HB2	1:J:19:ASP:HB2	2.04	0.40
1:F:57:LYS:HG3	1:F:118:GLY:HA3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	124/132 (94%)	118 (95%)	6 (5%)	0	100	100
1	B	130/132 (98%)	127 (98%)	3 (2%)	0	100	100
1	C	130/132 (98%)	127 (98%)	3 (2%)	0	100	100
1	D	129/132 (98%)	127 (98%)	2 (2%)	0	100	100
1	E	128/132 (97%)	122 (95%)	6 (5%)	0	100	100
1	F	130/132 (98%)	127 (98%)	3 (2%)	0	100	100
1	G	129/132 (98%)	124 (96%)	5 (4%)	0	100	100
1	H	127/132 (96%)	124 (98%)	3 (2%)	0	100	100
1	I	127/132 (96%)	125 (98%)	2 (2%)	0	100	100
1	J	127/132 (96%)	122 (96%)	5 (4%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	K	128/132 (97%)	126 (98%)	2 (2%)	0	100	100
1	L	124/132 (94%)	119 (96%)	5 (4%)	0	100	100
All	All	1533/1584 (97%)	1488 (97%)	45 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	110/116 (95%)	109 (99%)	1 (1%)	70	80
1	B	116/116 (100%)	115 (99%)	1 (1%)	70	80
1	C	116/116 (100%)	116 (100%)	0	100	100
1	D	115/116 (99%)	112 (97%)	3 (3%)	40	57
1	E	114/116 (98%)	113 (99%)	1 (1%)	70	80
1	F	116/116 (100%)	115 (99%)	1 (1%)	70	80
1	G	115/116 (99%)	114 (99%)	1 (1%)	70	80
1	H	113/116 (97%)	110 (97%)	3 (3%)	39	55
1	I	113/116 (97%)	112 (99%)	1 (1%)	70	80
1	J	113/116 (97%)	112 (99%)	1 (1%)	70	80
1	K	114/116 (98%)	112 (98%)	2 (2%)	51	67
1	L	111/116 (96%)	111 (100%)	0	100	100
All	All	1366/1392 (98%)	1351 (99%)	15 (1%)	65	77

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	GLU
1	B	-8	SER
1	G	11	LYS
1	H	39	GLU

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Mol	Chain	Res	Type
1	H	40	ASP
1	H	41	GLU
1	I	18	GLU
1	J	41	GLU
1	K	5	ASN
1	K	41	GLU
1	D	18	GLU
1	D	22	LYS
1	D	40	ASP
1	E	4	LYS
1	F	58	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	29	HIS
1	G	65	ASN
1	J	78	GLN
1	K	10	GLN
1	K	50	ASN
1	D	65	ASN
1	D	78	GLN
1	E	26	ASN
1	F	29	HIS

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	126/132 (95%)	0.67	18 (14%) 6 6	19, 37, 82, 104	0
1	B	132/132 (100%)	0.86	24 (18%) 3 3	21, 46, 83, 105	0
1	C	132/132 (100%)	0.59	15 (11%) 10 9	22, 40, 78, 99	0
1	D	131/132 (99%)	0.65	20 (15%) 5 5	20, 39, 73, 85	0
1	E	130/132 (98%)	0.85	26 (20%) 3 2	21, 46, 84, 104	0
1	F	132/132 (100%)	0.74	24 (18%) 3 3	20, 41, 75, 88	0
1	G	131/132 (99%)	0.69	19 (14%) 6 5	23, 42, 81, 98	0
1	H	129/132 (97%)	0.87	28 (21%) 2 2	23, 39, 92, 118	0
1	I	129/132 (97%)	1.45	37 (28%) 1 1	25, 48, 94, 108	0
1	J	129/132 (97%)	1.57	46 (35%) 1 0	26, 57, 99, 123	0
1	K	130/132 (98%)	1.20	33 (25%) 1 1	22, 46, 84, 99	0
1	L	126/132 (95%)	1.29	38 (30%) 1 1	22, 53, 98, 113	0
All	All	1557/1584 (98%)	0.95	328 (21%) 2 2	19, 44, 90, 123	0

All (328) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	-6	LEU	7.8
1	A	116	ILE	7.6
1	A	118	GLY	7.5
1	L	116	ILE	7.4
1	H	116	ILE	7.3
1	B	-8	SER	6.4
1	I	56	ILE	6.1
1	I	116	ILE	6.1
1	K	59	TYR	6.0
1	I	-7	GLY	6.0
1	I	115	GLY	6.0

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Mol	Chain	Res	Type	RSRZ
1	H	119	THR	5.8
1	C	-8	SER	5.8
1	H	-7	GLY	5.8
1	B	-9	SER	5.7
1	E	-7	GLY	5.6
1	J	113	VAL	5.6
1	J	9	THR	5.6
1	H	118	GLY	5.5
1	I	117	GLU	5.5
1	K	56	ILE	5.4
1	D	58	GLU	5.3
1	I	118	GLY	5.3
1	L	42	TYR	5.3
1	D	-8	SER	5.3
1	K	55	PRO	5.1
1	I	9	THR	5.1
1	A	-7	GLY	5.1
1	J	118	GLY	5.1
1	K	5	ASN	5.1
1	B	-7	GLY	5.0
1	I	39	GLU	4.9
1	B	12	TYR	4.8
1	J	114	LYS	4.8
1	I	121	GLU	4.8
1	H	115	GLY	4.8
1	D	41	GLU	4.8
1	H	113	VAL	4.8
1	A	56	ILE	4.8
1	J	5	ASN	4.7
1	K	58	GLU	4.7
1	I	12	TYR	4.6
1	H	-6	LEU	4.6
1	G	-7	GLY	4.5
1	A	114	LYS	4.5
1	H	117	GLU	4.5
1	L	119	THR	4.4
1	H	39	GLU	4.4
1	I	119	THR	4.4
1	L	8	CYS	4.4
1	C	-7	GLY	4.3
1	I	38	CYS	4.3
1	B	9	THR	4.3

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Mol	Chain	Res	Type	RSRZ
1	A	117	GLU	4.3
1	K	9	THR	4.3
1	G	11	LYS	4.3
1	C	-9	SER	4.2
1	J	57	LYS	4.2
1	L	118	GLY	4.2
1	E	41	GLU	4.2
1	B	39	GLU	4.1
1	K	61	ILE	4.1
1	F	55	PRO	4.1
1	J	42	TYR	4.1
1	H	58	GLU	4.1
1	J	121	GLU	4.1
1	K	60	GLY	4.1
1	J	39	GLU	4.1
1	B	37	ASP	4.1
1	J	-1	SER	4.1
1	E	39	GLU	4.0
1	B	42	TYR	4.0
1	K	8	CYS	4.0
1	L	117	GLU	4.0
1	H	-8	SER	4.0
1	L	-3	ARG	4.0
1	B	-6	LEU	4.0
1	J	10	GLN	4.0
1	I	41	GLU	3.9
1	J	120	SER	3.9
1	F	59	TYR	3.9
1	I	10	GLN	3.9
1	J	41	GLU	3.9
1	K	10	GLN	3.9
1	K	-7	GLY	3.9
1	E	38	CYS	3.9
1	E	42	TYR	3.9
1	I	60	GLY	3.8
1	H	114	LYS	3.8
1	K	12	TYR	3.8
1	G	-8	SER	3.8
1	E	-1	SER	3.8
1	L	115	GLY	3.8
1	D	-7	GLY	3.8
1	A	37	ASP	3.8

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Mol	Chain	Res	Type	RSRZ
1	E	59	TYR	3.8
1	I	40	ASP	3.8
1	J	40	ASP	3.7
1	L	114	LYS	3.7
1	H	-9	SER	3.7
1	E	-6	LEU	3.7
1	E	5	ASN	3.7
1	I	-6	LEU	3.7
1	L	12	TYR	3.7
1	I	120	SER	3.7
1	F	41	GLU	3.7
1	J	38	CYS	3.7
1	H	112	GLU	3.6
1	K	122	ILE	3.6
1	H	57	LYS	3.6
1	J	12	TYR	3.6
1	G	-2	GLY	3.6
1	J	56	ILE	3.6
1	L	-4	PRO	3.6
1	A	58	GLU	3.5
1	J	37	ASP	3.5
1	L	41	GLU	3.5
1	C	38	CYS	3.5
1	A	42	TYR	3.5
1	F	56	ILE	3.5
1	J	-4	PRO	3.4
1	G	-1	SER	3.4
1	I	114	LYS	3.4
1	J	59	TYR	3.4
1	C	42	TYR	3.4
1	J	116	ILE	3.4
1	A	-6	LEU	3.4
1	L	58	GLU	3.4
1	B	-5	VAL	3.3
1	F	60	GLY	3.3
1	B	85	TYR	3.3
1	D	38	CYS	3.3
1	E	-2	GLY	3.3
1	A	41	GLU	3.3
1	K	41	GLU	3.3
1	J	-6	LEU	3.2
1	J	4	LYS	3.2

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Mol	Chain	Res	Type	RSRZ
1	C	4	LYS	3.2
1	G	40	ASP	3.2
1	J	119	THR	3.2
1	H	56	ILE	3.2
1	K	116	ILE	3.2
1	D	22	LYS	3.2
1	J	-2	GLY	3.2
1	C	37	ASP	3.2
1	B	40	ASP	3.1
1	L	39	GLU	3.1
1	D	39	GLU	3.1
1	F	40	ASP	3.1
1	G	-5	VAL	3.1
1	K	113	VAL	3.1
1	K	115	GLY	3.1
1	J	112	GLU	3.1
1	F	39	GLU	3.1
1	F	-9	SER	3.1
1	J	60	GLY	3.1
1	F	116	ILE	3.1
1	A	57	LYS	3.0
1	G	9	THR	3.0
1	K	57	LYS	3.0
1	L	57	LYS	3.0
1	E	4	LYS	3.0
1	I	85	TYR	3.0
1	D	57	LYS	3.0
1	L	44	GLU	3.0
1	C	8	CYS	3.0
1	J	43	ILE	3.0
1	K	119	THR	3.0
1	C	41	GLU	3.0
1	I	59	TYR	3.0
1	K	85	TYR	3.0
1	L	40	ASP	3.0
1	K	54	GLY	2.9
1	B	41	GLU	2.9
1	K	4	LYS	2.9
1	L	37	ASP	2.9
1	E	40	ASP	2.9
1	J	-7	GLY	2.9
1	C	21	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	J	-3	ARG	2.9
1	H	40	ASP	2.9
1	F	37	ASP	2.9
1	B	-1	SER	2.9
1	I	44	GLU	2.9
1	I	42	TYR	2.9
1	D	59	TYR	2.9
1	L	43	ILE	2.8
1	B	38	CYS	2.8
1	L	38	CYS	2.8
1	F	38	CYS	2.8
1	E	36	ALA	2.8
1	C	-6	LEU	2.8
1	B	10	GLN	2.8
1	K	37	ASP	2.8
1	J	61	ILE	2.8
1	L	56	ILE	2.8
1	L	-5	VAL	2.8
1	G	-6	LEU	2.8
1	J	111	ARG	2.8
1	A	115	GLY	2.8
1	G	12	TYR	2.8
1	D	122	ILE	2.8
1	D	21	GLU	2.8
1	F	111	ARG	2.8
1	J	109	LEU	2.8
1	I	4	LYS	2.8
1	L	4	LYS	2.8
1	A	40	ASP	2.8
1	L	59	TYR	2.7
1	D	12	TYR	2.7
1	A	-5	VAL	2.7
1	A	113	VAL	2.7
1	G	58	GLU	2.7
1	J	8	CYS	2.7
1	J	55	PRO	2.7
1	H	111	ARG	2.7
1	L	112	GLU	2.7
1	E	-4	PRO	2.7
1	J	-5	VAL	2.7
1	K	118	GLY	2.7
1	H	4	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	F	12	TYR	2.7
1	J	82	ASP	2.6
1	J	110	ASP	2.6
1	E	37	ASP	2.6
1	K	111	ARG	2.6
1	G	38	CYS	2.6
1	J	58	GLU	2.6
1	D	55	PRO	2.6
1	J	117	GLU	2.6
1	F	121	GLU	2.6
1	E	43	ILE	2.6
1	L	9	THR	2.6
1	I	-2	GLY	2.5
1	G	41	GLU	2.5
1	E	21	GLU	2.5
1	H	107	ARG	2.5
1	G	42	TYR	2.5
1	G	39	GLU	2.5
1	K	-6	LEU	2.5
1	E	-5	VAL	2.5
1	L	36	ALA	2.5
1	I	8	CYS	2.5
1	I	111	ARG	2.5
1	F	118	GLY	2.5
1	D	37	ASP	2.5
1	G	10	GLN	2.5
1	L	-1	SER	2.5
1	B	-4	PRO	2.5
1	I	-3	ARG	2.5
1	B	21	GLU	2.4
1	B	58	GLU	2.4
1	E	58	GLU	2.4
1	F	114	LYS	2.4
1	K	40	ASP	2.4
1	L	19	ASP	2.4
1	B	11	LYS	2.4
1	K	117	GLU	2.4
1	L	-2	GLY	2.4
1	A	38	CYS	2.4
1	F	19	ASP	2.4
1	H	42	TYR	2.4
1	D	42	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	36	ALA	2.4
1	L	55	PRO	2.4
1	E	44	GLU	2.4
1	F	57	LYS	2.4
1	F	-8	SER	2.4
1	H	110	ASP	2.4
1	H	38	CYS	2.4
1	K	38	CYS	2.4
1	C	39	GLU	2.4
1	J	11	LYS	2.3
1	L	113	VAL	2.3
1	E	121	GLU	2.3
1	E	-3	ARG	2.3
1	E	57	LYS	2.3
1	L	10	GLN	2.3
1	D	121	GLU	2.3
1	F	42	TYR	2.3
1	C	-4	PRO	2.3
1	J	6	GLY	2.3
1	I	113	VAL	2.3
1	J	115	GLY	2.2
1	B	56	ILE	2.2
1	I	57	LYS	2.2
1	H	37	ASP	2.2
1	I	58	GLU	2.2
1	L	45	PRO	2.2
1	I	24	LYS	2.2
1	E	6	GLY	2.2
1	D	36	ALA	2.2
1	D	-1	SER	2.2
1	I	37	ASP	2.2
1	E	112	GLU	2.2
1	F	112	GLU	2.2
1	J	45	PRO	2.2
1	E	55	PRO	2.2
1	I	-5	VAL	2.2
1	G	43	ILE	2.2
1	I	61	ILE	2.2
1	B	19	ASP	2.2
1	H	-3	ARG	2.2
1	H	59	TYR	2.1
1	L	20	LYS	2.1

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Mol	Chain	Res	Type	RSRZ
1	F	113	VAL	2.1
1	I	6	GLY	2.1
1	C	6	GLY	2.1
1	B	-3	ARG	2.1
1	L	21	GLU	2.1
1	F	21	GLU	2.1
1	G	-3	ARG	2.1
1	H	109	LEU	2.1
1	L	60	GLY	2.1
1	F	58	GLU	2.1
1	D	56	ILE	2.1
1	K	114	LYS	2.1
1	K	42	TYR	2.1
1	H	55	PRO	2.1
1	C	-5	VAL	2.0
1	G	60	GLY	2.0
1	K	39	GLU	2.0
1	A	61	ILE	2.0
1	J	13	THR	2.0
1	J	108	THR	2.0
1	K	48	TYR	2.0
1	I	11	LYS	2.0
1	D	40	ASP	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](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q < 0.9' lists the number of atoms with occupancy less than 0.9.

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	A	201	1/1	0.99	0.12	54,54,54,54	0
2	ZN	D	201	1/1	0.99	0.09	55,55,55,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.