



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 03:33 pm BST

PDB ID : 4EMP
Title : Crystal structure of the mutant of ClpP E137A from Staphylococcus aureus
Authors : Ye, F.; Zhang, J.; Liu, H.; Luo, C.; Yang, C.-G.
Deposited on : 2012-04-12
Resolution : 2.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.11
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

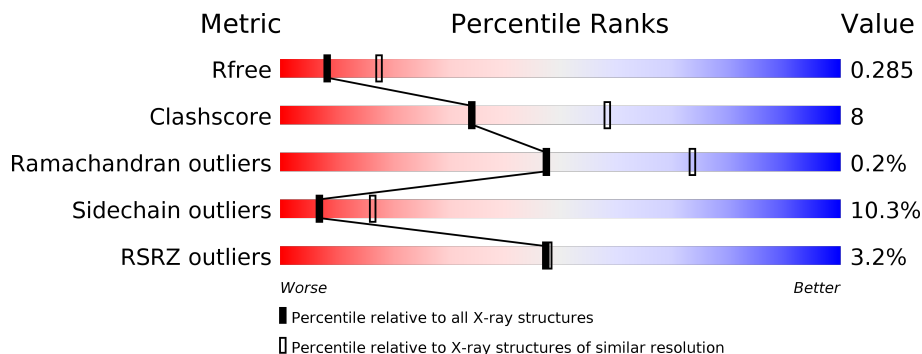
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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 on 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	200	<p>2% 70% 21% • 6%</p>
1	B	200	<p>0% 72% 17% • 9%</p>
1	C	200	<p>2% 70% 18% 5% 8%</p>
1	E	200	<p>5% 76% 13% • 8%</p>
1	F	200	<p>4% 72% 19% • 8%</p>
1	G	200	<p>2% 72% 16% • 10%</p>

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Mol	Chain	Length	Quality of chain
1	I	200	<p>3% 72% 17% 6% 6%</p>
1	K	200	<p>4% 72% 18% • 7%</p>
1	L	200	<p>6% 74% 15% • 8%</p>
1	M	200	<p>5% 69% 21% • 8%</p>
1	N	200	<p>3% 72% 18% • 8%</p>
1	S	200	<p>3% 73% 15% 5% 8%</p>
1	T	200	<p>4% 72% 18% • 7%</p>
1	V	200	<p>3% 78% 15% • 6%</p>

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 20028 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 ATP-dependent Clp protease proteolytic subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	V	189	1462	919	251	284	8	0	2	0
1	A	189	1462	919	251	284	8	0	2	0
1	B	182	1400	884	237	272	7	0	1	0
1	C	185	1430	902	243	277	8	0	2	0
1	E	185	1432	903	243	278	8	0	2	0
1	F	185	1427	899	243	278	7	0	1	0
1	G	180	1386	876	235	268	7	0	1	0
1	I	189	1452	911	251	284	6	0	0	0
1	K	186	1431	901	244	279	7	0	1	0
1	L	184	1413	890	242	275	6	0	0	0
1	M	185	1427	899	243	278	7	0	1	0
1	N	185	1427	899	243	278	7	0	1	0
1	S	184	1425	899	242	276	8	0	2	0
1	T	186	1431	901	244	279	7	0	1	0

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
V	-4	GLY	-	EXPRESSION TAG	UNP P63786

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Chain	Residue	Modelled	Actual	Comment	Reference
V	-3	PRO	-	EXPRESSION TAG	UNP P63786
V	-2	LEU	-	EXPRESSION TAG	UNP P63786
V	-1	GLY	-	EXPRESSION TAG	UNP P63786
V	0	SER	-	EXPRESSION TAG	UNP P63786
V	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
A	-4	GLY	-	EXPRESSION TAG	UNP P63786
A	-3	PRO	-	EXPRESSION TAG	UNP P63786
A	-2	LEU	-	EXPRESSION TAG	UNP P63786
A	-1	GLY	-	EXPRESSION TAG	UNP P63786
A	0	SER	-	EXPRESSION TAG	UNP P63786
A	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
B	-4	GLY	-	EXPRESSION TAG	UNP P63786
B	-3	PRO	-	EXPRESSION TAG	UNP P63786
B	-2	LEU	-	EXPRESSION TAG	UNP P63786
B	-1	GLY	-	EXPRESSION TAG	UNP P63786
B	0	SER	-	EXPRESSION TAG	UNP P63786
B	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
C	-4	GLY	-	EXPRESSION TAG	UNP P63786
C	-3	PRO	-	EXPRESSION TAG	UNP P63786
C	-2	LEU	-	EXPRESSION TAG	UNP P63786
C	-1	GLY	-	EXPRESSION TAG	UNP P63786
C	0	SER	-	EXPRESSION TAG	UNP P63786
C	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
E	-4	GLY	-	EXPRESSION TAG	UNP P63786
E	-3	PRO	-	EXPRESSION TAG	UNP P63786
E	-2	LEU	-	EXPRESSION TAG	UNP P63786
E	-1	GLY	-	EXPRESSION TAG	UNP P63786
E	0	SER	-	EXPRESSION TAG	UNP P63786
E	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
F	-4	GLY	-	EXPRESSION TAG	UNP P63786
F	-3	PRO	-	EXPRESSION TAG	UNP P63786
F	-2	LEU	-	EXPRESSION TAG	UNP P63786
F	-1	GLY	-	EXPRESSION TAG	UNP P63786
F	0	SER	-	EXPRESSION TAG	UNP P63786
F	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
G	-4	GLY	-	EXPRESSION TAG	UNP P63786
G	-3	PRO	-	EXPRESSION TAG	UNP P63786
G	-2	LEU	-	EXPRESSION TAG	UNP P63786
G	-1	GLY	-	EXPRESSION TAG	UNP P63786
G	0	SER	-	EXPRESSION TAG	UNP P63786
G	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
I	-4	GLY	-	EXPRESSION TAG	UNP P63786

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Chain	Residue	Modelled	Actual	Comment	Reference
I	-3	PRO	-	EXPRESSION TAG	UNP P63786
I	-2	LEU	-	EXPRESSION TAG	UNP P63786
I	-1	GLY	-	EXPRESSION TAG	UNP P63786
I	0	SER	-	EXPRESSION TAG	UNP P63786
I	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
K	-4	GLY	-	EXPRESSION TAG	UNP P63786
K	-3	PRO	-	EXPRESSION TAG	UNP P63786
K	-2	LEU	-	EXPRESSION TAG	UNP P63786
K	-1	GLY	-	EXPRESSION TAG	UNP P63786
K	0	SER	-	EXPRESSION TAG	UNP P63786
K	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
L	-4	GLY	-	EXPRESSION TAG	UNP P63786
L	-3	PRO	-	EXPRESSION TAG	UNP P63786
L	-2	LEU	-	EXPRESSION TAG	UNP P63786
L	-1	GLY	-	EXPRESSION TAG	UNP P63786
L	0	SER	-	EXPRESSION TAG	UNP P63786
L	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
M	-4	GLY	-	EXPRESSION TAG	UNP P63786
M	-3	PRO	-	EXPRESSION TAG	UNP P63786
M	-2	LEU	-	EXPRESSION TAG	UNP P63786
M	-1	GLY	-	EXPRESSION TAG	UNP P63786
M	0	SER	-	EXPRESSION TAG	UNP P63786
M	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
N	-4	GLY	-	EXPRESSION TAG	UNP P63786
N	-3	PRO	-	EXPRESSION TAG	UNP P63786
N	-2	LEU	-	EXPRESSION TAG	UNP P63786
N	-1	GLY	-	EXPRESSION TAG	UNP P63786
N	0	SER	-	EXPRESSION TAG	UNP P63786
N	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
S	-4	GLY	-	EXPRESSION TAG	UNP P63786
S	-3	PRO	-	EXPRESSION TAG	UNP P63786
S	-2	LEU	-	EXPRESSION TAG	UNP P63786
S	-1	GLY	-	EXPRESSION TAG	UNP P63786
S	0	SER	-	EXPRESSION TAG	UNP P63786
S	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786
T	-4	GLY	-	EXPRESSION TAG	UNP P63786
T	-3	PRO	-	EXPRESSION TAG	UNP P63786
T	-2	LEU	-	EXPRESSION TAG	UNP P63786
T	-1	GLY	-	EXPRESSION TAG	UNP P63786
T	0	SER	-	EXPRESSION TAG	UNP P63786
T	137	ALA	GLU	ENGINEERED MUTATION	UNP P63786


- Molecule 2 is water.

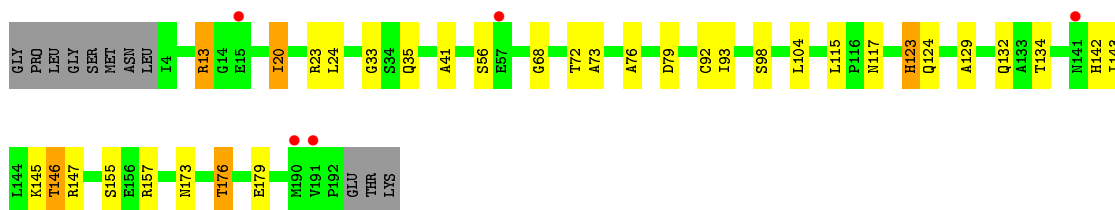
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	V	1	Total O 1 1	0	0
2	A	3	Total O 3 3	0	0
2	B	4	Total O 4 4	0	0
2	C	1	Total O 1 1	0	0
2	E	2	Total O 2 2	0	0
2	F	1	Total O 1 1	0	0
2	G	2	Total O 2 2	0	0
2	I	2	Total O 2 2	0	0
2	K	1	Total O 1 1	0	0
2	N	1	Total O 1 1	0	0
2	S	4	Total O 4 4	0	0
2	T	1	Total O 1 1	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

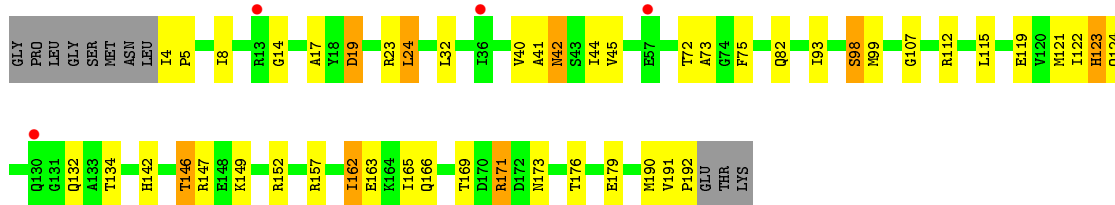
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain V: 



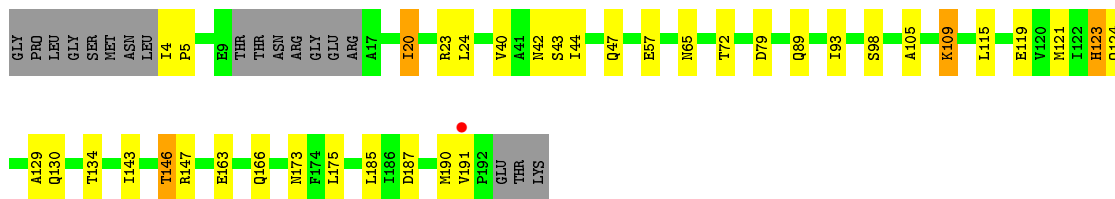
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain A: 



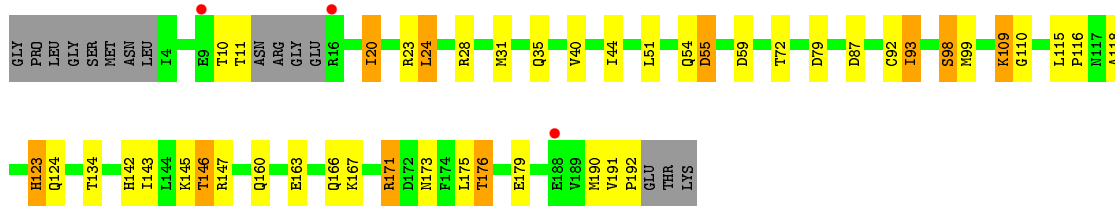
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain B: 

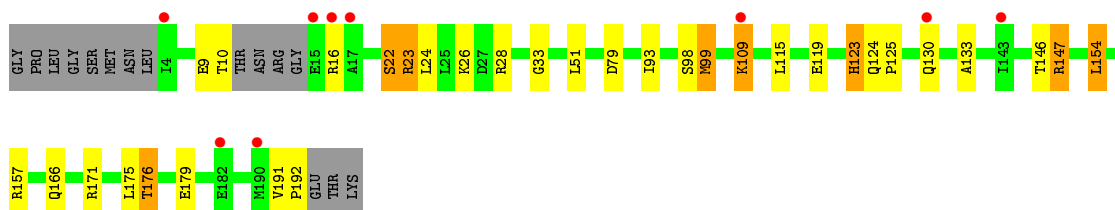
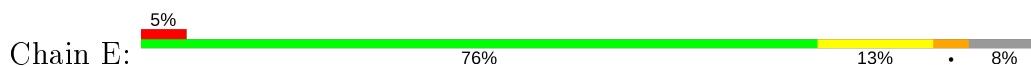


- Molecule 1: ATP-dependent Clp protease proteolytic subunit

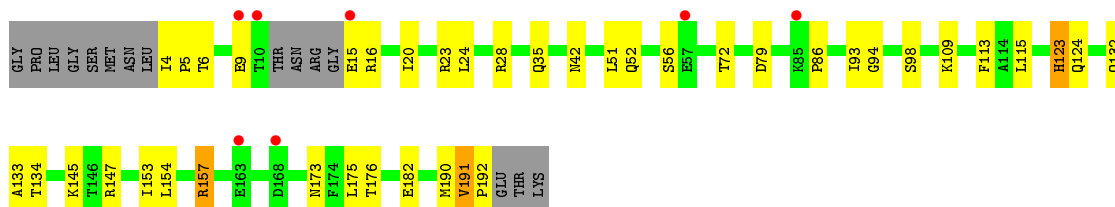
Chain C: 



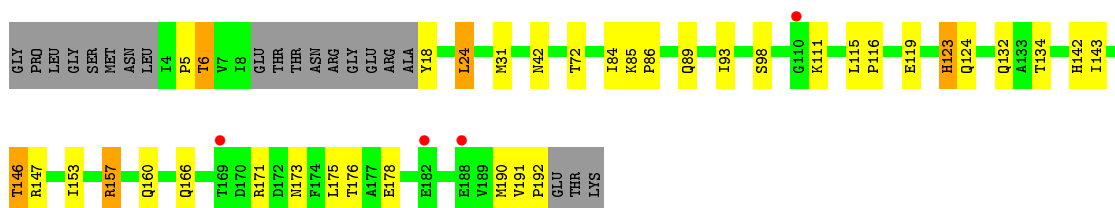
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



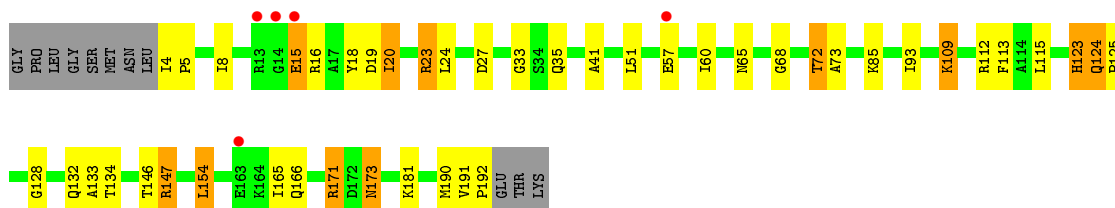
- Molecule 1: ATP-dependent Clp protease proteolytic subunit



- Molecule 1: ATP-dependent Clp protease proteolytic subunit

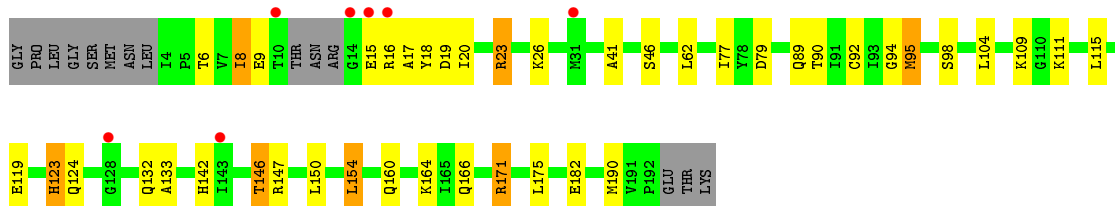


- Molecule 1: ATP-dependent Clp protease proteolytic subunit



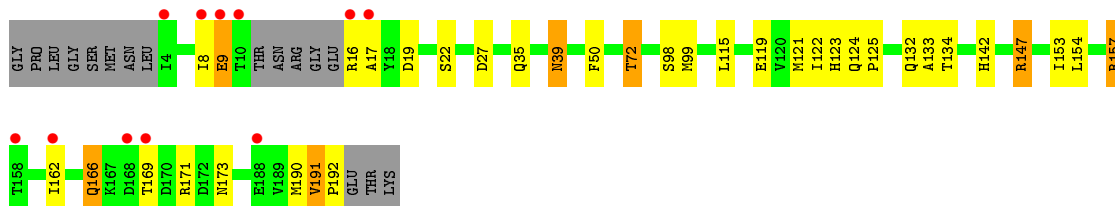
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain K: 



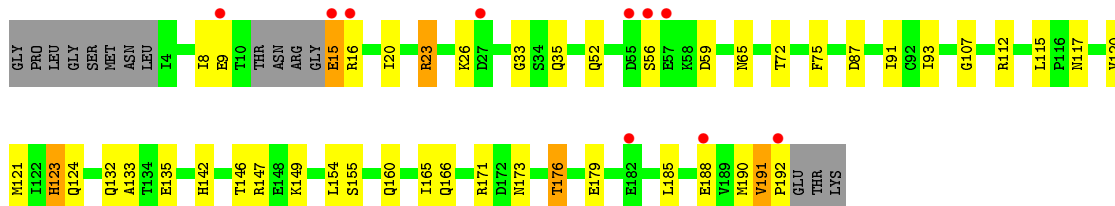
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain L: 



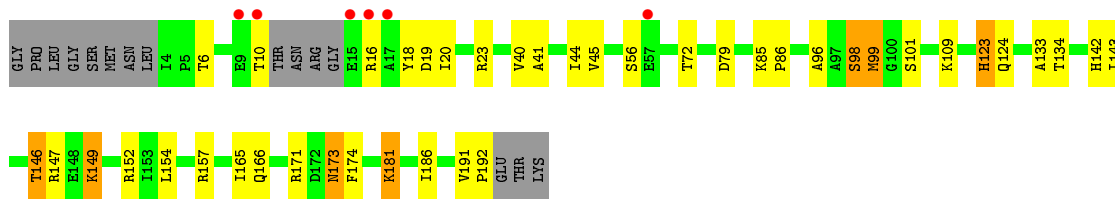
- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain M: 



- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain N: 



- Molecule 1: ATP-dependent Clp protease proteolytic subunit

Chain S: 





- Molecule 1: ATP-dependent Clp protease proteolytic subunit



4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	168.73Å 96.06Å 192.43Å 90.00° 91.17° 90.00°	Depositor
Resolution (Å)	30.00 – 2.70 46.60 – 2.61	Depositor EDS
% Data completeness (in resolution range)	95.2 (30.00-2.70) 94.6 (46.60-2.61)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.61Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.235 , 0.290 0.231 , 0.285	Depositor DCC
R_{free} test set	4444 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	52.9	Xtriage
Anisotropy	0.245	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 29.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.012 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.010 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.014 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.015 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.016 for -h,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	20028	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.49	0/1487	0.62	0/2007
1	B	0.53	0/1421	0.66	1/1918 (0.1%)
1	C	0.49	0/1454	0.61	0/1962
1	E	0.48	0/1456	0.59	0/1964
1	F	0.47	0/1448	0.61	0/1954
1	G	0.48	0/1407	0.61	0/1899
1	I	0.49	0/1471	0.60	0/1987
1	K	0.45	0/1452	0.60	0/1959
1	L	0.44	0/1431	0.59	0/1932
1	M	0.46	0/1448	0.59	0/1954
1	N	0.47	0/1448	0.59	0/1954
1	S	0.50	0/1449	0.65	0/1954
1	T	0.48	0/1452	0.61	0/1959
1	V	0.51	0/1487	0.63	0/2007
All	All	0.48	0/20311	0.61	1/27410 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	175	LEU	CA-CB-CG	5.12	127.07	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	4	ILE	Peptide
1	B	5	PRO	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	1462	0	1482	46	0
1	B	1400	0	1417	25	0
1	C	1430	0	1453	32	0
1	E	1432	0	1452	28	0
1	F	1427	0	1443	24	0
1	G	1386	0	1406	31	0
1	I	1452	0	1464	34	0
1	K	1431	0	1446	37	0
1	L	1413	0	1428	32	0
1	M	1427	0	1443	30	0
1	N	1427	0	1443	27	0
1	S	1425	0	1445	32	0
1	T	1431	0	1446	30	0
1	V	1462	0	1482	25	0
2	A	3	0	0	0	0
2	B	4	0	0	0	0
2	C	1	0	0	0	0
2	E	2	0	0	0	0
2	F	1	0	0	0	0
2	G	2	0	0	0	0
2	I	2	0	0	0	0
2	K	1	0	0	0	0
2	N	1	0	0	0	0
2	S	4	0	0	0	0
2	T	1	0	0	0	0
2	V	1	0	0	0	0
All	All	20028	0	20250	336	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 (336) 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:L:191:VAL:HG22	1:L:192:PRO:HA	1.27	1.14
1:E:93:ILE:HG22	1:E:115:LEU:CD1	1.83	1.09
1:M:191:VAL:HG22	1:M:192:PRO:HA	1.34	1.08
1:E:93:ILE:HG22	1:E:115:LEU:HD12	1.37	1.04
1:F:191:VAL:HG22	1:F:192:PRO:HA	1.37	1.02
1:G:191:VAL:HG22	1:G:192:PRO:HA	1.45	0.98
1:N:98:SER:HB3	1:N:123:HIS:CE1	2.03	0.93
1:S:171:ARG:HH22	1:T:132:GLN:NE2	1.68	0.91
1:A:132:GLN:HE22	1:G:171:ARG:NH2	1.68	0.91
1:V:33:GLY:HA3	1:B:42:ASN:HD21	1.39	0.87
1:B:98:SER:HB3	1:B:123:HIS:CE1	2.09	0.87
1:C:171:ARG:HH22	1:F:132:GLN:NE2	1.73	0.86
1:F:98:SER:HB3	1:F:123:HIS:CE1	2.11	0.85
1:F:93:ILE:HG22	1:F:115:LEU:HD12	1.58	0.85
1:L:171:ARG:HH22	1:M:132:GLN:NE2	1.75	0.85
1:K:6:THR:O	1:L:22:SER:HB3	1.77	0.84
1:N:191:VAL:HG22	1:N:192:PRO:HA	1.59	0.83
1:B:119:GLU:OE2	1:C:142:HIS:HE1	1.61	0.83
1:A:152:ARG:HA	1:A:162:ILE:HD11	1.59	0.82
1:A:132:GLN:NE2	1:G:171:ARG:NH2	2.29	0.80
1:I:132:GLN:HE22	1:T:171:ARG:HH22	1.28	0.79
1:I:191:VAL:HB	1:I:192:PRO:HA	1.65	0.78
1:A:132:GLN:HE22	1:G:171:ARG:HH22	1.30	0.77
1:K:98:SER:HB2	1:K:123:HIS:CE1	2.19	0.77
1:S:191:VAL:HG22	1:S:192:PRO:HA	1.65	0.77
1:F:9:GLU:HB2	1:F:16:ARG:HB2	1.67	0.76
1:K:98:SER:CB	1:K:123:HIS:CE1	2.68	0.76
1:K:171:ARG:HH22	1:L:132:GLN:NE2	1.84	0.76
1:A:98:SER:HB2	1:A:123:HIS:CE1	2.23	0.74
1:F:93:ILE:HG22	1:F:115:LEU:CD1	2.18	0.73
1:C:176:THR:HG22	1:C:179:GLU:H	1.51	0.73
1:F:134:THR:H	1:N:124:GLN:HE22	1.35	0.73
1:L:171:ARG:NH2	1:M:132:GLN:NE2	2.36	0.73
1:C:40:VAL:O	1:C:44:ILE:HG12	1.89	0.72
1:I:93:ILE:HG22	1:I:115:LEU:HD12	1.69	0.72
1:S:98:SER:HB2	1:S:123:HIS:CE1	2.25	0.72
1:T:16:ARG:HB3	1:T:18:TYR:HE1	1.53	0.72
1:E:99:MET:HE2	1:E:99:MET:HA	1.71	0.71
1:M:155:SER:OG	1:M:160:GLN:O	2.08	0.71
1:S:171:ARG:NH2	1:T:132:GLN:NE2	2.38	0.71
1:B:134:THR:H	1:T:124:GLN:HE22	1.37	0.71
1:C:171:ARG:HH22	1:F:132:GLN:HE22	1.39	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:142:HIS:O	1:K:146:THR:HG23	1.92	0.70
1:A:134:THR:H	1:K:124:GLN:HE22	1.39	0.69
1:K:160:GLN:HE21	1:K:164:LYS:HD3	1.57	0.69
1:T:16:ARG:HB3	1:T:18:TYR:CE1	2.27	0.69
1:I:132:GLN:NE2	1:T:171:ARG:HH22	1.89	0.69
1:V:132:GLN:NE2	1:A:171:ARG:HH22	1.90	0.69
1:C:124:GLN:HE21	1:S:133:ALA:HB3	1.58	0.69
1:S:6:THR:O	1:T:22:SER:HB3	1.93	0.68
1:V:33:GLY:HA3	1:B:42:ASN:ND2	2.06	0.68
1:A:41:ALA:HB2	1:A:73:ALA:HB1	1.74	0.68
1:L:115:LEU:HD21	1:L:190:MET:HE3	1.76	0.68
1:K:9:GLU:HB2	1:K:16:ARG:HB2	1.74	0.68
1:K:160:GLN:NE2	1:K:164:LYS:HD3	2.09	0.68
1:M:9:GLU:HB2	1:M:16:ARG:HG3	1.75	0.68
1:T:98:SER:HB2	1:T:123:HIS:CE1	2.29	0.68
1:E:109:LYS:HA	1:E:109:LYS:HE2	1.74	0.67
1:B:124:GLN:HE22	1:T:134:THR:H	1.42	0.67
1:B:40:VAL:O	1:B:44:ILE:HG12	1.95	0.67
1:C:93:ILE:HG22	1:C:115:LEU:HD12	1.77	0.67
1:K:8:ILE:HG21	1:L:16:ARG:HG3	1.75	0.67
1:T:164:LYS:NZ	1:T:168:ASP:OD2	2.27	0.67
1:E:176:THR:HG22	1:E:179:GLU:H	1.60	0.66
1:G:98:SER:HB3	1:G:123:HIS:CE1	2.31	0.66
1:M:191:VAL:CG2	1:M:192:PRO:HA	2.19	0.66
1:K:98:SER:HB3	1:K:123:HIS:CE1	2.31	0.65
1:N:191:VAL:CG2	1:N:192:PRO:HA	2.26	0.65
1:V:98:SER:HB2	1:V:123:HIS:CE1	2.32	0.65
1:L:191:VAL:CG2	1:L:192:PRO:HA	2.17	0.64
1:A:123:HIS:HD1	1:A:123:HIS:C	2.00	0.64
1:E:93:ILE:CG2	1:E:115:LEU:HD12	2.22	0.64
1:I:171:ARG:HH22	1:K:132:GLN:NE2	1.95	0.64
1:T:143:ILE:HA	1:T:146:THR:HG23	1.80	0.64
1:N:142:HIS:O	1:N:146:THR:HG22	1.99	0.63
1:C:54:GLN:O	1:C:55:ASP:HB2	1.98	0.63
1:A:124:GLN:HE21	1:K:133:ALA:HB3	1.63	0.63
1:C:98:SER:OG	1:C:99:MET:N	2.32	0.63
1:A:142:HIS:HE1	1:G:119:GLU:OE2	1.80	0.63
1:M:93:ILE:HG13	1:M:93:ILE:O	1.99	0.63
1:T:176:THR:HG22	1:T:179:GLU:H	1.62	0.62
1:B:115:LEU:HD13	1:C:79:ASP:HB3	1.81	0.62
1:F:153:ILE:O	1:F:157:ARG:HG3	1.98	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:119:GLU:OE2	1:C:142:HIS:CE1	2.48	0.61
1:B:163:GLU:CD	1:B:163:GLU:H	2.03	0.61
1:F:124:GLN:HE22	1:N:134:THR:H	1.49	0.61
1:K:98:SER:HB2	1:K:123:HIS:HE1	1.63	0.61
1:E:79:ASP:HB3	1:F:115:LEU:HD13	1.82	0.61
1:K:89:GLN:HG2	1:K:111:LYS:HB3	1.83	0.61
1:V:134:THR:HG23	1:I:124:GLN:HE22	1.65	0.60
1:I:93:ILE:HG22	1:I:115:LEU:CD1	2.31	0.60
1:A:121:MET:HE3	1:A:123:HIS:HB3	1.82	0.60
1:A:93:ILE:HG22	1:A:115:LEU:HD12	1.83	0.60
1:E:22:SER:HB3	1:F:6:THR:O	2.00	0.60
1:I:15:GLU:HG2	1:K:16:ARG:HG2	1.82	0.60
1:C:171:ARG:NH2	1:F:132:GLN:NE2	2.47	0.60
1:C:124:GLN:NE2	1:S:134:THR:HG23	2.16	0.60
1:N:149:LYS:HG2	1:N:152:ARG:NH2	2.17	0.60
1:A:142:HIS:O	1:A:146:THR:HG23	2.02	0.60
1:V:124:GLN:NE2	1:I:134:THR:H	2.00	0.60
1:I:171:ARG:HH22	1:K:132:GLN:HE21	1.50	0.59
1:A:98:SER:OG	1:A:99:MET:N	2.34	0.59
1:A:115:LEU:HD21	1:A:190:MET:HE3	1.84	0.59
1:E:98:SER:OG	1:E:99:MET:N	2.36	0.59
1:L:9:GLU:HB2	1:L:16:ARG:HD3	1.85	0.59
1:G:176:THR:HG22	1:G:178:GLU:H	1.67	0.59
1:S:93:ILE:HG22	1:S:115:LEU:HD12	1.85	0.59
1:L:98:SER:H	1:L:121:MET:HE2	1.67	0.59
1:E:99:MET:CE	1:E:99:MET:HA	2.32	0.59
1:S:171:ARG:HH22	1:T:132:GLN:HE22	1.50	0.59
1:C:191:VAL:HA	1:C:192:PRO:C	2.23	0.58
1:L:153:ILE:O	1:L:157:ARG:HG3	2.03	0.58
1:C:93:ILE:HG22	1:C:115:LEU:CD1	2.33	0.58
1:G:93:ILE:HG22	1:G:115:LEU:HD12	1.84	0.58
1:G:142:HIS:O	1:G:146:THR:HG23	2.04	0.57
1:G:124:GLN:HE21	1:L:133:ALA:HB3	1.69	0.57
1:S:5:PRO:HD2	1:S:20:ILE:HG12	1.87	0.57
1:G:93:ILE:HG22	1:G:115:LEU:CD1	2.34	0.57
1:C:115:LEU:HD13	1:F:79:ASP:HB3	1.87	0.57
1:S:143:ILE:HA	1:S:146:THR:HG23	1.87	0.57
1:G:134:THR:HG23	1:L:124:GLN:HE22	1.70	0.57
1:E:33:GLY:HA3	1:G:42:ASN:HD21	1.69	0.56
1:G:116:PRO:HD3	1:G:190:MET:O	2.05	0.56
1:V:93:ILE:HG22	1:V:115:LEU:HD12	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:GLN:NE2	1:G:171:ARG:HH21	2.01	0.56
1:B:143:ILE:HA	1:B:146:THR:HG23	1.87	0.56
1:G:124:GLN:HE22	1:L:134:THR:H	1.53	0.56
1:K:92:CYS:HB2	1:K:104:LEU:HD13	1.87	0.56
1:S:23:ARG:C	1:S:23:ARG:HD3	2.26	0.56
1:C:124:GLN:NE2	1:S:134:THR:H	2.04	0.56
1:V:98:SER:CB	1:V:123:HIS:CE1	2.90	0.55
1:N:40:VAL:O	1:N:44:ILE:HG12	2.07	0.55
1:C:134:THR:H	1:S:124:GLN:HE22	1.55	0.55
1:T:98:SER:OG	1:T:99:MET:N	2.41	0.54
1:K:171:ARG:HH22	1:L:132:GLN:HE21	1.55	0.54
1:T:23:ARG:HD3	1:T:23:ARG:O	2.08	0.54
1:A:123:HIS:ND1	1:A:123:HIS:C	2.60	0.54
1:V:142:HIS:O	1:V:146:THR:HG22	2.08	0.54
1:E:124:GLN:HE21	1:M:133:ALA:HB3	1.72	0.53
1:M:154:LEU:HB3	1:M:165:ILE:HD13	1.89	0.53
1:S:40:VAL:O	1:S:44:ILE:HG12	2.08	0.53
1:G:98:SER:CB	1:G:123:HIS:CE1	2.91	0.53
1:V:124:GLN:HE21	1:I:133:ALA:HB3	1.72	0.53
1:E:123:HIS:HD1	1:E:123:HIS:C	2.12	0.53
1:L:166:GLN:HE21	1:L:166:GLN:HA	1.74	0.53
1:I:20:ILE:HD12	1:K:46:SER:HB3	1.90	0.52
1:E:191:VAL:HG13	1:E:192:PRO:HA	1.92	0.52
1:M:8:ILE:HG21	1:N:16:ARG:HG2	1.90	0.52
1:T:40:VAL:O	1:T:44:ILE:HG12	2.10	0.52
1:M:59:ASP:OD1	1:M:87:ASP:HB2	2.10	0.52
1:A:122:ILE:HD12	1:A:169:THR:HG22	1.92	0.51
1:C:143:ILE:HA	1:C:146:THR:HG23	1.92	0.51
1:A:115:LEU:CD2	1:A:190:MET:HE3	2.40	0.51
1:F:113:PHE:HB3	1:F:190:MET:HG3	1.91	0.51
1:C:98:SER:HB2	1:C:123:HIS:CE1	2.45	0.51
1:C:24:LEU:HD23	1:C:31[A]:MET:HE1	1.92	0.51
1:I:4:ILE:N	1:I:5:PRO:CD	2.73	0.51
1:E:33:GLY:HA3	1:G:42:ASN:ND2	2.26	0.51
1:M:120:VAL:HG11	1:M:185:LEU:HD13	1.92	0.51
1:M:176:THR:HG22	1:M:179:GLU:H	1.76	0.51
1:K:18:TYR:OH	1:K:26:LYS:HE2	2.11	0.51
1:S:93:ILE:O	1:S:93:ILE:HG13	2.10	0.51
1:E:133:ALA:HB3	1:M:124:GLN:HE21	1.76	0.51
1:I:51:LEU:HD13	1:I:60:ILE:HG12	1.93	0.51
1:V:124:GLN:HE22	1:I:134:THR:HG23	1.76	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:154:LEU:HB3	1:I:165:ILE:HD13	1.94	0.50
1:E:176:THR:HB	1:E:179:GLU:OE1	2.12	0.50
1:E:28:ARG:HG2	1:E:51:LEU:HD22	1.94	0.50
1:M:9:GLU:O	1:M:15:GLU:HA	2.12	0.50
1:T:28:ARG:HG2	1:T:51:LEU:HD22	1.93	0.50
1:F:133:ALA:HB3	1:N:124:GLN:HE21	1.76	0.50
1:V:79:ASP:HB3	1:A:115:LEU:HD13	1.94	0.50
1:M:23:ARG:HH21	1:M:26:LYS:HB3	1.77	0.49
1:V:176:THR:HG22	1:V:179:GLU:H	1.77	0.49
1:A:134:THR:HG23	1:K:124:GLN:NE2	2.27	0.49
1:I:72:THR:HG21	1:T:95[A]:MET:HE2	1.93	0.49
1:L:171:ARG:HH22	1:M:132:GLN:HE22	1.56	0.49
1:I:125:PRO:HD2	1:I:147:ARG:HG2	1.94	0.49
1:I:33:GLY:HA2	1:I:65:ASN:O	2.12	0.49
1:V:142:HIS:HE1	1:A:119:GLU:OE2	1.96	0.49
1:C:28:ARG:HG2	1:C:51:LEU:HD22	1.95	0.49
1:I:8:ILE:HD11	1:K:17:ALA:O	2.13	0.49
1:L:115:LEU:HD21	1:L:190:MET:CE	2.42	0.49
1:M:93:ILE:HG22	1:M:115:LEU:CD1	2.43	0.48
1:E:125:PRO:HD2	1:E:147:ARG:HG2	1.96	0.48
1:S:171:ARG:NH2	1:T:132:GLN:HE21	2.11	0.48
1:T:23:ARG:HD3	1:T:23:ARG:C	2.33	0.48
1:V:41:ALA:HB2	1:V:73:ALA:HB1	1.95	0.48
1:K:150:LEU:O	1:K:154:LEU:HB2	2.14	0.48
1:M:115:LEU:HD13	1:N:79:ASP:HB3	1.95	0.48
1:S:23:ARG:HD3	1:S:23:ARG:O	2.13	0.48
1:V:117:ASN:HB2	1:B:79:ASP:OD1	2.13	0.48
1:A:115:LEU:HD21	1:A:190:MET:CE	2.44	0.48
1:A:134:THR:HG23	1:K:124:GLN:HE22	1.78	0.48
1:B:115:LEU:HD23	1:B:190:MET:HB2	1.95	0.48
1:E:123:HIS:C	1:E:123:HIS:ND1	2.66	0.48
1:G:89:GLN:HG2	1:G:111:LYS:HB3	1.96	0.48
1:S:154:LEU:HB3	1:S:165:ILE:HD13	1.95	0.48
1:G:153:ILE:O	1:G:157:ARG:HG3	2.14	0.47
1:A:162:ILE:HG22	1:A:163:GLU:OE1	2.13	0.47
1:B:98:SER:CB	1:B:123:HIS:CE1	2.90	0.47
1:G:191:VAL:CG2	1:G:192:PRO:HA	2.31	0.47
1:I:23:ARG:NH2	1:I:27:ASP:OD1	2.47	0.47
1:K:23:ARG:NH1	1:L:50:PHE:HE1	2.13	0.47
1:M:75:PHE:CZ	1:M:149:LYS:HD3	2.49	0.47
1:S:98:SER:HB2	1:S:123:HIS:HE1	1.74	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:20:ILE:HA	1:B:20:ILE:HD12	1.71	0.47
1:N:99:MET:CE	1:N:99:MET:HA	2.44	0.47
1:S:166:GLN:HE21	1:S:166:GLN:HA	1.80	0.47
1:S:8:ILE:HA	1:S:16:ARG:O	2.15	0.47
1:T:98:SER:H	1:T:121:MET:HE2	1.80	0.47
1:I:41:ALA:HB2	1:I:73:ALA:HB1	1.97	0.47
1:C:109:LYS:NZ	1:C:110:GLY:H	2.13	0.47
1:N:143:ILE:HA	1:N:146:THR:CG2	2.46	0.46
1:A:107:GLY:HA3	1:A:112:ARG:HG2	1.98	0.46
1:F:98:SER:CB	1:F:123:HIS:CE1	2.91	0.46
1:G:84:ILE:HD12	1:G:86:PRO:HG2	1.97	0.46
1:A:152:ARG:CA	1:A:162:ILE:HD11	2.36	0.46
1:M:107:GLY:O	1:M:112:ARG:HD3	2.16	0.46
1:M:33:GLY:HA2	1:M:65:ASN:O	2.16	0.46
1:B:163:GLU:CD	1:B:163:GLU:N	2.68	0.46
1:M:91:ILE:HG23	1:M:190:MET:CE	2.46	0.46
1:N:98:SER:HB2	1:N:99:MET:H	1.59	0.46
1:L:119:GLU:OE2	1:M:142:HIS:HE1	1.98	0.46
1:B:124:GLN:HE21	1:T:133:ALA:HB3	1.80	0.46
1:S:65:ASN:HB2	1:S:93:ILE:HG13	1.98	0.46
1:I:123:HIS:C	1:I:123:HIS:HD1	2.19	0.46
1:M:123:HIS:ND1	1:M:123:HIS:C	2.69	0.46
1:C:92:CYS:SG	1:C:118:ALA:HB1	2.56	0.46
1:M:121:MET:HE3	1:M:123:HIS:HB3	1.97	0.46
1:K:119:GLU:OE2	1:L:142:HIS:HE1	2.00	0.45
1:A:124:GLN:NE2	1:K:133:ALA:HB3	2.30	0.45
1:A:165:ILE:O	1:A:169:THR:HG23	2.16	0.45
1:B:124:GLN:NE2	1:T:134:THR:HG23	2.31	0.45
1:E:124:GLN:NE2	1:M:133:ALA:HB3	2.30	0.45
1:A:191:VAL:HG22	1:A:192:PRO:HD2	1.99	0.45
1:S:125:PRO:HD2	1:S:147:ARG:HG2	1.98	0.45
1:V:20:ILE:HG13	1:V:20:ILE:H	1.60	0.45
1:K:115:LEU:HD21	1:K:190:MET:HE2	1.98	0.45
1:A:42:ASN:HD21	1:G:31:MET:HB3	1.82	0.45
1:L:154:LEU:HD23	1:L:154:LEU:HA	1.86	0.45
1:K:95[A]:MET:HB3	1:L:72:THR:HG21	1.99	0.45
1:V:35:GLN:HG3	1:V:68:GLY:O	2.16	0.45
1:K:41:ALA:HA	1:K:77:ILE:HD11	1.98	0.45
1:L:8:ILE:HG12	1:L:17:ALA:HA	1.99	0.45
1:F:124:GLN:HE21	1:N:133:ALA:HB3	1.81	0.44
1:B:109:LYS:HE3	1:B:187:ASP:OD1	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:23:ARG:HH11	1:L:50:PHE:HE1	1.65	0.44
1:N:98:SER:HB3	1:N:123:HIS:HE1	1.74	0.44
1:F:28:ARG:HG2	1:F:51:LEU:HD22	1.99	0.44
1:L:39:ASN:HD22	1:L:39:ASN:HA	1.67	0.44
1:A:115:LEU:CD2	1:A:190:MET:CE	2.95	0.44
1:N:173:ASN:HD22	1:N:174:PHE:N	2.16	0.44
1:B:105:ALA:HB2	1:B:185:LEU:HD23	2.00	0.44
1:I:123:HIS:C	1:I:123:HIS:ND1	2.70	0.44
1:S:98:SER:OG	1:S:99:MET:N	2.50	0.44
1:C:24:LEU:HD23	1:C:31[A]:MET:CE	2.48	0.44
1:I:173:ASN:HD22	1:I:173:ASN:C	2.20	0.44
1:G:24:LEU:HD12	1:G:24:LEU:HA	1.88	0.44
1:N:16:ARG:HB3	1:N:18:TYR:CE1	2.53	0.44
1:N:41:ALA:O	1:N:45:VAL:HG23	2.17	0.44
1:N:6:THR:O	1:S:22:SER:HB3	2.17	0.43
1:A:4:ILE:HA	1:A:5:PRO:HD3	1.79	0.43
1:C:124:GLN:HE22	1:S:134:THR:HG23	1.81	0.43
1:B:129:ALA:O	1:T:128:GLY:HA2	2.18	0.43
1:F:133:ALA:HB3	1:N:124:GLN:NE2	2.33	0.43
1:I:16:ARG:HB3	1:I:18:TYR:CE1	2.54	0.43
1:B:98:SER:H	1:B:121:MET:HE2	1.83	0.43
1:I:35:GLN:HG2	1:I:68:GLY:O	2.19	0.43
1:K:171:ARG:HH22	1:L:132:GLN:HE22	1.64	0.43
1:A:8:ILE:HG12	1:A:17:ALA:CB	2.49	0.43
1:E:175:LEU:HD12	1:E:179:GLU:HB3	2.00	0.43
1:G:85:LYS:N	1:G:86:PRO:HD2	2.33	0.43
1:L:171:ARG:NH2	1:M:132:GLN:HE22	2.11	0.43
1:F:56:SER:HA	1:F:86:PRO:HG3	2.01	0.43
1:L:122:ILE:HD12	1:L:169:THR:HG22	2.01	0.43
1:E:98:SER:HB2	1:E:123:HIS:CE1	2.54	0.43
1:G:124:GLN:NE2	1:L:133:ALA:HB3	2.32	0.43
1:A:176:THR:OG1	1:A:179:GLU:HG3	2.18	0.42
1:A:24:LEU:HD12	1:A:24:LEU:HA	1.91	0.42
1:C:20:ILE:HG13	1:C:20:ILE:H	1.65	0.42
1:A:45:VAL:HG12	1:G:31:MET:HE1	2.00	0.42
1:K:62:LEU:HD23	1:K:90:THR:HG22	2.01	0.42
1:V:92:CYS:HB2	1:V:104:LEU:HD13	2.02	0.42
1:A:19:ASP:HB3	1:G:6:THR:HG22	2.01	0.42
1:A:93:ILE:HG22	1:A:115:LEU:CD1	2.47	0.42
1:E:98:SER:HB2	1:E:123:HIS:NE2	2.34	0.42
1:G:98:SER:CB	1:G:123:HIS:HE1	2.31	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:132:GLN:HE21	1:A:171:ARG:HH22	1.66	0.42
1:S:23:ARG:CD	1:S:23:ARG:C	2.88	0.42
1:T:99:MET:CE	1:T:99:MET:HA	2.49	0.42
1:C:163:GLU:HG2	1:C:167:LYS:NZ	2.34	0.42
1:I:4:ILE:N	1:I:5:PRO:HD3	2.35	0.42
1:V:142:HIS:CE1	1:A:119:GLU:OE2	2.73	0.42
1:L:125:PRO:HD2	1:L:147:ARG:HG2	2.02	0.42
1:S:122:ILE:HG13	1:S:173:ASN:HB3	2.01	0.42
1:N:96:ALA:O	1:N:101:SER:HB3	2.19	0.42
1:I:113:PHE:HB3	1:I:190:MET:HG3	2.02	0.42
1:C:59:ASP:OD1	1:C:87:ASP:HB2	2.20	0.41
1:K:115:LEU:HD21	1:K:190:MET:CE	2.50	0.41
1:T:75:PHE:CZ	1:T:149:LYS:HD3	2.55	0.41
1:B:43:SER:O	1:B:47:GLN:HG3	2.19	0.41
1:N:10:THR:OG1	1:S:16:ARG:NH2	2.48	0.41
1:N:85:LYS:N	1:N:86:PRO:CD	2.84	0.41
1:E:119:GLU:OE2	1:G:142:HIS:HE1	2.03	0.41
1:E:23:ARG:HH21	1:E:26:LYS:HD2	1.85	0.41
1:M:117:ASN:O	1:N:149:LYS:NZ	2.53	0.41
1:T:150:LEU:O	1:T:154:LEU:HB2	2.20	0.41
1:V:129:ALA:O	1:I:128:GLY:HA2	2.19	0.41
1:V:73:ALA:O	1:V:76:ALA:HB3	2.21	0.41
1:N:181:LYS:HA	1:N:186:ILE:HG13	2.03	0.41
1:B:109:LYS:HD2	1:B:109:LYS:HA	1.98	0.41
1:C:31[A]:MET:HG2	1:F:42:ASN:OD1	2.21	0.41
1:F:4:ILE:HA	1:F:5:PRO:HD3	1.71	0.41
1:I:115:LEU:HD13	1:K:79:ASP:HB3	2.02	0.41
1:K:154:LEU:HA	1:K:154:LEU:HD23	1.92	0.41
1:V:13:ARG:HG2	1:A:14:GLY:HA2	2.03	0.41
1:A:82:GLN:HA	1:A:82:GLN:NE2	2.36	0.41
1:I:109:LYS:HD2	1:I:109:LYS:HA	1.76	0.41
1:I:112:ARG:HE	1:I:112:ARG:HB2	1.59	0.41
1:E:171:ARG:HH22	1:G:132:GLN:NE2	2.19	0.41
1:N:154:LEU:HB3	1:N:165:ILE:HD13	2.03	0.41
1:A:75:PHE:CZ	1:A:149:LYS:HD3	2.57	0.40
1:C:11:THR:O	1:C:11:THR:HG23	2.21	0.40
1:E:154:LEU:HD23	1:E:154:LEU:HA	1.92	0.40
1:F:9:GLU:O	1:F:15:GLU:HA	2.21	0.40
1:T:191:VAL:HG22	1:T:192:PRO:HA	2.03	0.40
1:A:40:VAL:O	1:A:44:ILE:HG12	2.21	0.40
1:B:65:ASN:HB2	1:B:93:ILE:HG13	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:116:PRO:HD3	1:C:190:MET:O	2.21	0.40
1:I:57:GLU:OE1	1:I:85:LYS:NZ	2.49	0.40
1:T:30:ILE:HD11	1:T:51:LEU:HD12	2.03	0.40
1:V:143:ILE:HA	1:V:146:THR:HG23	2.03	0.40
1:S:105:ALA:HB2	1:S:185:LEU:HD23	2.02	0.40
1:L:171:ARG:HH21	1:M:135:GLU:HG3	1.85	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	189/200 (94%)	182 (96%)	6 (3%)	1 (0%)	29	54
1	B	179/200 (90%)	173 (97%)	6 (3%)	0	100	100
1	C	183/200 (92%)	178 (97%)	4 (2%)	1 (0%)	29	54
1	E	183/200 (92%)	181 (99%)	2 (1%)	0	100	100
1	F	182/200 (91%)	177 (97%)	4 (2%)	1 (0%)	29	54
1	G	177/200 (88%)	169 (96%)	7 (4%)	1 (1%)	25	50
1	I	187/200 (94%)	181 (97%)	6 (3%)	0	100	100
1	K	183/200 (92%)	175 (96%)	7 (4%)	1 (0%)	29	54
1	L	180/200 (90%)	172 (96%)	8 (4%)	0	100	100
1	M	182/200 (91%)	174 (96%)	7 (4%)	1 (0%)	29	54
1	N	182/200 (91%)	172 (94%)	10 (6%)	0	100	100
1	S	182/200 (91%)	177 (97%)	5 (3%)	0	100	100
1	T	183/200 (92%)	179 (98%)	4 (2%)	0	100	100
1	V	189/200 (94%)	185 (98%)	4 (2%)	0	100	100
All	All	2561/2800 (92%)	2475 (97%)	80 (3%)	6 (0%)	47	73

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	55	ASP
1	G	5	PRO
1	M	56	SER
1	A	98	SER
1	F	94	GLY
1	K	94	GLY

5.3.2 Protein sidechains [i](#)

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	157/165 (95%)	143 (91%)	14 (9%)	9 22
1	B	150/165 (91%)	136 (91%)	14 (9%)	9 21
1	C	154/165 (93%)	135 (88%)	19 (12%)	4 11
1	E	154/165 (93%)	138 (90%)	16 (10%)	7 16
1	F	153/165 (93%)	136 (89%)	17 (11%)	6 14
1	G	149/165 (90%)	136 (91%)	13 (9%)	10 23
1	I	155/165 (94%)	139 (90%)	16 (10%)	7 16
1	K	153/165 (93%)	137 (90%)	16 (10%)	7 16
1	L	151/165 (92%)	137 (91%)	14 (9%)	9 21
1	M	153/165 (93%)	138 (90%)	15 (10%)	8 18
1	N	153/165 (93%)	136 (89%)	17 (11%)	6 14
1	S	153/165 (93%)	136 (89%)	17 (11%)	6 14
1	T	153/165 (93%)	133 (87%)	20 (13%)	4 10
1	V	157/165 (95%)	143 (91%)	14 (9%)	9 22
All	All	2145/2310 (93%)	1923 (90%)	222 (10%)	7 16

All (222) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	V	13	ARG
1	V	20	ILE
1	V	23	ARG
1	V	24	LEU
1	V	56	SER
1	V	72	THR
1	V	123	HIS
1	V	145	LYS
1	V	146	THR
1	V	147	ARG
1	V	155	SER
1	V	157	ARG
1	V	173	ASN
1	V	176	THR
1	A	19	ASP
1	A	23	ARG
1	A	24	LEU
1	A	32	LEU
1	A	42	ASN
1	A	72	THR
1	A	123	HIS
1	A	146	THR
1	A	147	ARG
1	A	157	ARG
1	A	162	ILE
1	A	166	GLN
1	A	171	ARG
1	A	173	ASN
1	B	20	ILE
1	B	23	ARG
1	B	24	LEU
1	B	57	GLU
1	B	72	THR
1	B	89	GLN
1	B	109	LYS
1	B	123	HIS
1	B	130	GLN
1	B	146	THR
1	B	147	ARG
1	B	166	GLN
1	B	173	ASN
1	B	191	VAL
1	C	10	THR

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Mol	Chain	Res	Type
1	C	20	ILE
1	C	23	ARG
1	C	24	LEU
1	C	35	GLN
1	C	72	THR
1	C	93	ILE
1	C	98	SER
1	C	109	LYS
1	C	123	HIS
1	C	145	LYS
1	C	146	THR
1	C	147	ARG
1	C	160	GLN
1	C	166	GLN
1	C	171	ARG
1	C	173	ASN
1	C	175	LEU
1	C	176	THR
1	E	9	GLU
1	E	10	THR
1	E	16	ARG
1	E	22	SER
1	E	23	ARG
1	E	24	LEU
1	E	99	MET
1	E	109	LYS
1	E	123	HIS
1	E	130	GLN
1	E	146	THR
1	E	147	ARG
1	E	154	LEU
1	E	157	ARG
1	E	166	GLN
1	E	176	THR
1	F	20	ILE
1	F	23	ARG
1	F	24	LEU
1	F	35	GLN
1	F	52	GLN
1	F	72	THR
1	F	109	LYS
1	F	123	HIS

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Mol	Chain	Res	Type
1	F	145	LYS
1	F	147	ARG
1	F	154	LEU
1	F	157	ARG
1	F	173	ASN
1	F	175	LEU
1	F	176	THR
1	F	182	GLU
1	F	191	VAL
1	G	6	THR
1	G	18	TYR
1	G	24	LEU
1	G	72	THR
1	G	123	HIS
1	G	143	ILE
1	G	146	THR
1	G	147	ARG
1	G	157	ARG
1	G	160	GLN
1	G	166	GLN
1	G	173	ASN
1	G	175	LEU
1	I	15	GLU
1	I	19	ASP
1	I	20	ILE
1	I	23	ARG
1	I	24	LEU
1	I	72	THR
1	I	109	LYS
1	I	123	HIS
1	I	124	GLN
1	I	146	THR
1	I	147	ARG
1	I	154	LEU
1	I	166	GLN
1	I	171	ARG
1	I	173	ASN
1	I	181	LYS
1	K	8	ILE
1	K	15	GLU
1	K	19	ASP
1	K	20	ILE

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Mol	Chain	Res	Type
1	K	23	ARG
1	K	95[A]	MET
1	K	95[B]	MET
1	K	109	LYS
1	K	123	HIS
1	K	146	THR
1	K	147	ARG
1	K	154	LEU
1	K	166	GLN
1	K	171	ARG
1	K	175	LEU
1	K	182	GLU
1	L	9	GLU
1	L	19	ASP
1	L	27	ASP
1	L	35	GLN
1	L	39	ASN
1	L	72	THR
1	L	99	MET
1	L	123	HIS
1	L	147	ARG
1	L	157	ARG
1	L	162	ILE
1	L	166	GLN
1	L	173	ASN
1	L	191	VAL
1	M	15	GLU
1	M	20	ILE
1	M	23	ARG
1	M	35	GLN
1	M	52	GLN
1	M	72	THR
1	M	123	HIS
1	M	146	THR
1	M	147	ARG
1	M	166	GLN
1	M	171	ARG
1	M	173	ASN
1	M	176	THR
1	M	188	GLU
1	M	191	VAL
1	N	19	ASP

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Mol	Chain	Res	Type
1	N	20	ILE
1	N	23	ARG
1	N	56	SER
1	N	72	THR
1	N	98	SER
1	N	99	MET
1	N	109	LYS
1	N	123	HIS
1	N	146	THR
1	N	147	ARG
1	N	149	LYS
1	N	157	ARG
1	N	166	GLN
1	N	171	ARG
1	N	173	ASN
1	N	181	LYS
1	S	9	GLU
1	S	20	ILE
1	S	23	ARG
1	S	24	LEU
1	S	31[A]	MET
1	S	31[B]	MET
1	S	57	GLU
1	S	72	THR
1	S	123	HIS
1	S	146	THR
1	S	147	ARG
1	S	166	GLN
1	S	171	ARG
1	S	173	ASN
1	S	175	LEU
1	S	181	LYS
1	S	191	VAL
1	T	23	ARG
1	T	24	LEU
1	T	35	GLN
1	T	42	ASN
1	T	72	THR
1	T	95[A]	MET
1	T	95[B]	MET
1	T	109	LYS
1	T	123	HIS

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Mol	Chain	Res	Type
1	T	146	THR
1	T	147	ARG
1	T	154	LEU
1	T	157	ARG
1	T	162	ILE
1	T	166	GLN
1	T	171	ARG
1	T	173	ASN
1	T	175	LEU
1	T	176	THR
1	T	191	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (136) such sidechains are listed below:

Mol	Chain	Res	Type
1	V	82	GLN
1	V	89	GLN
1	V	117	ASN
1	V	123	HIS
1	V	124	GLN
1	V	132	GLN
1	V	142	HIS
1	V	151	ASN
1	V	160	GLN
1	V	166	GLN
1	V	173	ASN
1	A	39	ASN
1	A	42	ASN
1	A	82	GLN
1	A	89	GLN
1	A	117	ASN
1	A	124	GLN
1	A	132	GLN
1	A	142	HIS
1	A	151	ASN
1	A	160	GLN
1	A	166	GLN
1	A	173	ASN
1	B	39	ASN
1	B	42	ASN
1	B	89	GLN
1	B	123	HIS

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Mol	Chain	Res	Type
1	B	124	GLN
1	B	132	GLN
1	B	142	HIS
1	B	151	ASN
1	B	160	GLN
1	B	166	GLN
1	B	173	ASN
1	C	82	GLN
1	C	89	GLN
1	C	123	HIS
1	C	124	GLN
1	C	132	GLN
1	C	142	HIS
1	C	151	ASN
1	C	160	GLN
1	C	166	GLN
1	C	173	ASN
1	E	35	GLN
1	E	39	ASN
1	E	82	GLN
1	E	89	GLN
1	E	124	GLN
1	E	151	ASN
1	E	166	GLN
1	E	173	ASN
1	F	39	ASN
1	F	89	GLN
1	F	117	ASN
1	F	123	HIS
1	F	124	GLN
1	F	132	GLN
1	F	151	ASN
1	F	166	GLN
1	F	173	ASN
1	G	42	ASN
1	G	82	GLN
1	G	89	GLN
1	G	123	HIS
1	G	124	GLN
1	G	132	GLN
1	G	141	ASN
1	G	142	HIS

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Mol	Chain	Res	Type
1	G	151	ASN
1	G	160	GLN
1	G	166	GLN
1	G	173	ASN
1	I	89	GLN
1	I	124	GLN
1	I	132	GLN
1	I	142	HIS
1	I	151	ASN
1	I	160	GLN
1	I	166	GLN
1	I	173	ASN
1	K	52	GLN
1	K	82	GLN
1	K	89	GLN
1	K	123	HIS
1	K	124	GLN
1	K	132	GLN
1	K	151	ASN
1	K	160	GLN
1	K	166	GLN
1	K	173	ASN
1	L	39	ASN
1	L	89	GLN
1	L	117	ASN
1	L	124	GLN
1	L	132	GLN
1	L	142	HIS
1	L	160	GLN
1	L	166	GLN
1	L	173	ASN
1	M	42	ASN
1	M	82	GLN
1	M	89	GLN
1	M	117	ASN
1	M	124	GLN
1	M	132	GLN
1	M	142	HIS
1	M	160	GLN
1	M	166	GLN
1	M	173	ASN
1	N	39	ASN

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Mol	Chain	Res	Type
1	N	89	GLN
1	N	124	GLN
1	N	132	GLN
1	N	142	HIS
1	N	151	ASN
1	N	160	GLN
1	N	166	GLN
1	N	173	ASN
1	S	39	ASN
1	S	89	GLN
1	S	117	ASN
1	S	123	HIS
1	S	124	GLN
1	S	142	HIS
1	S	151	ASN
1	S	160	GLN
1	S	166	GLN
1	S	173	ASN
1	T	89	GLN
1	T	124	GLN
1	T	132	GLN
1	T	151	ASN
1	T	160	GLN
1	T	166	GLN
1	T	173	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates 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	189/200 (94%)	0.05	4 (2%) 63 65	43, 54, 68, 84	0
1	B	182/200 (91%)	-0.07	1 (0%) 91 92	46, 54, 64, 72	0
1	C	185/200 (92%)	0.03	3 (1%) 72 74	47, 56, 67, 78	0
1	E	185/200 (92%)	0.17	9 (4%) 29 28	54, 61, 73, 81	0
1	F	185/200 (92%)	0.18	7 (3%) 40 39	53, 62, 73, 85	0
1	G	180/200 (90%)	0.15	4 (2%) 62 63	49, 60, 69, 75	0
1	I	189/200 (94%)	0.10	5 (2%) 56 57	50, 60, 77, 93	0
1	K	186/200 (93%)	0.19	7 (3%) 40 39	51, 63, 81, 88	0
1	L	184/200 (92%)	0.42	11 (5%) 21 20	58, 65, 84, 103	0
1	M	185/200 (92%)	0.38	10 (5%) 25 24	57, 63, 74, 83	0
1	N	185/200 (92%)	0.21	6 (3%) 47 48	50, 60, 73, 85	0
1	S	184/200 (92%)	0.11	5 (2%) 54 55	49, 55, 75, 91	0
1	T	186/200 (93%)	0.07	7 (3%) 40 39	48, 57, 79, 89	0
1	V	189/200 (94%)	0.06	5 (2%) 56 57	44, 55, 68, 83	0
All	All	2594/2800 (92%)	0.15	84 (3%) 47 48	43, 59, 75, 103	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	10	THR	7.2
1	S	17	ALA	6.5
1	N	15	GLU	5.5
1	K	14	GLY	4.9
1	L	8	ILE	4.7
1	L	16	ARG	4.6
1	S	16	ARG	4.5
1	L	9	GLU	4.5

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Mol	Chain	Res	Type	RSRZ
1	I	15	GLU	4.3
1	A	13	ARG	4.0
1	K	15	GLU	4.0
1	E	16	ARG	4.0
1	M	15	GLU	4.0
1	L	162	ILE	3.8
1	S	9	GLU	3.8
1	E	109	LYS	3.7
1	N	16	ARG	3.6
1	T	16	ARG	3.6
1	A	57	GLU	3.6
1	S	15	GLU	3.4
1	M	55	ASP	3.4
1	T	10	THR	3.4
1	G	188	GLU	3.4
1	M	57	GLU	3.3
1	V	191	VAL	3.3
1	M	9	GLU	3.3
1	M	56	SER	3.2
1	E	17	ALA	3.2
1	T	15	GLU	3.2
1	M	16	ARG	3.1
1	L	17	ALA	3.1
1	V	15	GLU	3.1
1	K	10	THR	3.1
1	E	15	GLU	3.1
1	L	188	GLU	3.0
1	F	9	GLU	3.0
1	I	13	ARG	3.0
1	E	143	ILE	2.9
1	K	16	ARG	2.9
1	F	168	ASP	2.9
1	F	163	GLU	2.8
1	F	57	GLU	2.8
1	G	110	GLY	2.8
1	V	190	MET	2.8
1	L	158	THR	2.8
1	I	57	GLU	2.7
1	F	10	THR	2.7
1	F	85	LYS	2.7
1	M	182	GLU	2.6
1	N	17	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	K	143	ILE	2.5
1	T	9	GLU	2.5
1	I	14	GLY	2.5
1	F	15	GLU	2.5
1	E	182	GLU	2.5
1	S	23	ARG	2.5
1	C	188	GLU	2.4
1	K	31	MET	2.4
1	L	4	ILE	2.4
1	K	128	GLY	2.4
1	E	130	GLN	2.3
1	G	182	GLU	2.3
1	I	163	GLU	2.3
1	M	27	ASP	2.3
1	M	192	PRO	2.3
1	C	16	ARG	2.3
1	L	169	THR	2.3
1	A	36	ILE	2.2
1	T	17	ALA	2.2
1	N	9	GLU	2.2
1	L	168	ASP	2.2
1	T	19	ASP	2.2
1	N	57	GLU	2.1
1	A	130	GLN	2.1
1	V	57	GLU	2.1
1	E	4	ILE	2.1
1	C	9	GLU	2.1
1	E	190	MET	2.1
1	G	169	THR	2.1
1	V	141	ASN	2.1
1	N	10	THR	2.1
1	T	190	MET	2.1
1	B	191	VAL	2.1
1	M	188	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.