



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 08:37 PM EDT

PDB ID : 7CTO / pdb_00007cto
Title : Staphylococcus aureus MsrB
Authors : Kim, H.J.
Deposited on : 2020-08-19
Resolution : 3.47 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:

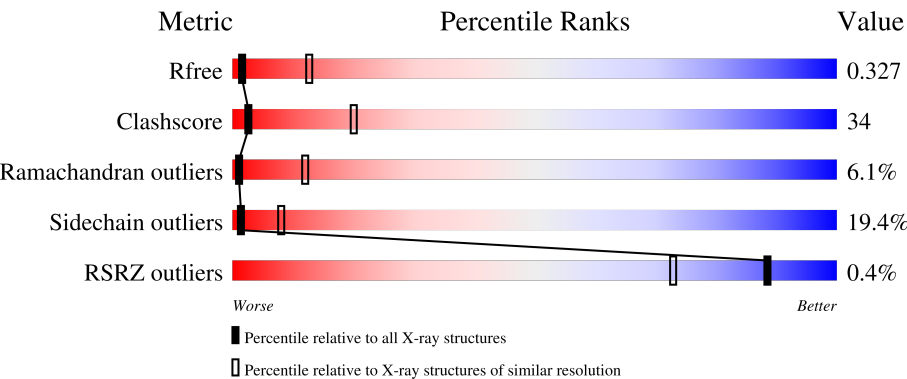
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
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 3.47 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	1083 (3.52-3.44)
Clashscore	190562	1139 (3.52-3.44)
Ramachandran outliers	187476	1111 (3.52-3.44)
Sidechain outliers	187428	1112 (3.52-3.44)
RSRZ outliers	180081	1082 (3.52-3.44)

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	159	35%37%11%•16%
1	B	159	% 33%42%10%•13%
1	C	159	34%38%13%14%
1	D	159	42%32%12%14%
1	E	159	31%41%13%•14%

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Mol	Chain	Length	Quality of chain
1	F	159	% 41% 31% 12% • 15%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 6564 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 Peptide methionine sulfoxide reductase MsrB.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	134	Total	C	N	O	S	0	0	0
			1076	684	172	216	4			
1	B	139	Total	C	N	O	S	0	0	0
			1121	713	179	225	4			
1	C	136	Total	C	N	O	S	0	0	0
			1093	694	175	220	4			
1	D	136	Total	C	N	O	S	0	0	0
			1095	695	174	222	4			
1	E	136	Total	C	N	O	S	0	0	0
			1095	697	174	220	4			
1	F	135	Total	C	N	O	S	0	0	0
			1084	688	173	219	4			

There are 102 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-16	MET	-	initiating methionine	UNP W8TTH3
A	-15	HIS	-	expression tag	UNP W8TTH3
A	-14	HIS	-	expression tag	UNP W8TTH3
A	-13	HIS	-	expression tag	UNP W8TTH3
A	-12	HIS	-	expression tag	UNP W8TTH3
A	-11	HIS	-	expression tag	UNP W8TTH3
A	-10	HIS	-	expression tag	UNP W8TTH3
A	-9	GLU	-	expression tag	UNP W8TTH3
A	-8	ASN	-	expression tag	UNP W8TTH3
A	-7	LEU	-	expression tag	UNP W8TTH3
A	-6	TYR	-	expression tag	UNP W8TTH3
A	-5	PHE	-	expression tag	UNP W8TTH3
A	-4	GLN	-	expression tag	UNP W8TTH3
A	-3	GLY	-	expression tag	UNP W8TTH3
A	-2	ALA	-	expression tag	UNP W8TTH3
A	-1	ALA	-	expression tag	UNP W8TTH3
A	0	SER	-	expression tag	UNP W8TTH3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	MET	-	initiating methionine	UNP W8TTH3
B	-15	HIS	-	expression tag	UNP W8TTH3
B	-14	HIS	-	expression tag	UNP W8TTH3
B	-13	HIS	-	expression tag	UNP W8TTH3
B	-12	HIS	-	expression tag	UNP W8TTH3
B	-11	HIS	-	expression tag	UNP W8TTH3
B	-10	HIS	-	expression tag	UNP W8TTH3
B	-9	GLU	-	expression tag	UNP W8TTH3
B	-8	ASN	-	expression tag	UNP W8TTH3
B	-7	LEU	-	expression tag	UNP W8TTH3
B	-6	TYR	-	expression tag	UNP W8TTH3
B	-5	PHE	-	expression tag	UNP W8TTH3
B	-4	GLN	-	expression tag	UNP W8TTH3
B	-3	GLY	-	expression tag	UNP W8TTH3
B	-2	ALA	-	expression tag	UNP W8TTH3
B	-1	ALA	-	expression tag	UNP W8TTH3
B	0	SER	-	expression tag	UNP W8TTH3
C	-16	MET	-	initiating methionine	UNP W8TTH3
C	-15	HIS	-	expression tag	UNP W8TTH3
C	-14	HIS	-	expression tag	UNP W8TTH3
C	-13	HIS	-	expression tag	UNP W8TTH3
C	-12	HIS	-	expression tag	UNP W8TTH3
C	-11	HIS	-	expression tag	UNP W8TTH3
C	-10	HIS	-	expression tag	UNP W8TTH3
C	-9	GLU	-	expression tag	UNP W8TTH3
C	-8	ASN	-	expression tag	UNP W8TTH3
C	-7	LEU	-	expression tag	UNP W8TTH3
C	-6	TYR	-	expression tag	UNP W8TTH3
C	-5	PHE	-	expression tag	UNP W8TTH3
C	-4	GLN	-	expression tag	UNP W8TTH3
C	-3	GLY	-	expression tag	UNP W8TTH3
C	-2	ALA	-	expression tag	UNP W8TTH3
C	-1	ALA	-	expression tag	UNP W8TTH3
C	0	SER	-	expression tag	UNP W8TTH3
D	-16	MET	-	initiating methionine	UNP W8TTH3
D	-15	HIS	-	expression tag	UNP W8TTH3
D	-14	HIS	-	expression tag	UNP W8TTH3
D	-13	HIS	-	expression tag	UNP W8TTH3
D	-12	HIS	-	expression tag	UNP W8TTH3
D	-11	HIS	-	expression tag	UNP W8TTH3
D	-10	HIS	-	expression tag	UNP W8TTH3
D	-9	GLU	-	expression tag	UNP W8TTH3

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-8	ASN	-	expression tag	UNP W8TTH3
D	-7	LEU	-	expression tag	UNP W8TTH3
D	-6	TYR	-	expression tag	UNP W8TTH3
D	-5	PHE	-	expression tag	UNP W8TTH3
D	-4	GLN	-	expression tag	UNP W8TTH3
D	-3	GLY	-	expression tag	UNP W8TTH3
D	-2	ALA	-	expression tag	UNP W8TTH3
D	-1	ALA	-	expression tag	UNP W8TTH3
D	0	SER	-	expression tag	UNP W8TTH3
E	-16	MET	-	initiating methionine	UNP W8TTH3
E	-15	HIS	-	expression tag	UNP W8TTH3
E	-14	HIS	-	expression tag	UNP W8TTH3
E	-13	HIS	-	expression tag	UNP W8TTH3
E	-12	HIS	-	expression tag	UNP W8TTH3
E	-11	HIS	-	expression tag	UNP W8TTH3
E	-10	HIS	-	expression tag	UNP W8TTH3
E	-9	GLU	-	expression tag	UNP W8TTH3
E	-8	ASN	-	expression tag	UNP W8TTH3
E	-7	LEU	-	expression tag	UNP W8TTH3
E	-6	TYR	-	expression tag	UNP W8TTH3
E	-5	PHE	-	expression tag	UNP W8TTH3
E	-4	GLN	-	expression tag	UNP W8TTH3
E	-3	GLY	-	expression tag	UNP W8TTH3
E	-2	ALA	-	expression tag	UNP W8TTH3
E	-1	ALA	-	expression tag	UNP W8TTH3
E	0	SER	-	expression tag	UNP W8TTH3
F	-16	MET	-	initiating methionine	UNP W8TTH3
F	-15	HIS	-	expression tag	UNP W8TTH3
F	-14	HIS	-	expression tag	UNP W8TTH3
F	-13	HIS	-	expression tag	UNP W8TTH3
F	-12	HIS	-	expression tag	UNP W8TTH3
F	-11	HIS	-	expression tag	UNP W8TTH3
F	-10	HIS	-	expression tag	UNP W8TTH3
F	-9	GLU	-	expression tag	UNP W8TTH3
F	-8	ASN	-	expression tag	UNP W8TTH3
F	-7	LEU	-	expression tag	UNP W8TTH3
F	-6	TYR	-	expression tag	UNP W8TTH3
F	-5	PHE	-	expression tag	UNP W8TTH3
F	-4	GLN	-	expression tag	UNP W8TTH3
F	-3	GLY	-	expression tag	UNP W8TTH3
F	-2	ALA	-	expression tag	UNP W8TTH3
F	-1	ALA	-	expression tag	UNP W8TTH3

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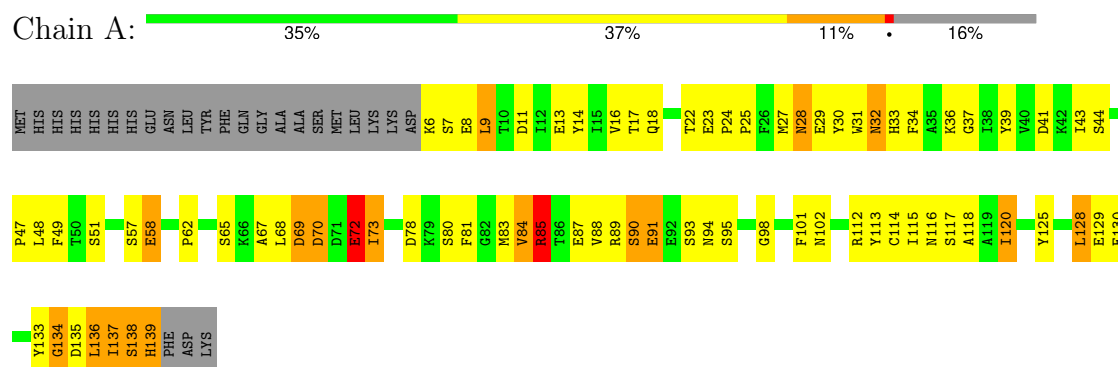
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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	SER	-	expression tag	UNP W8TTH3

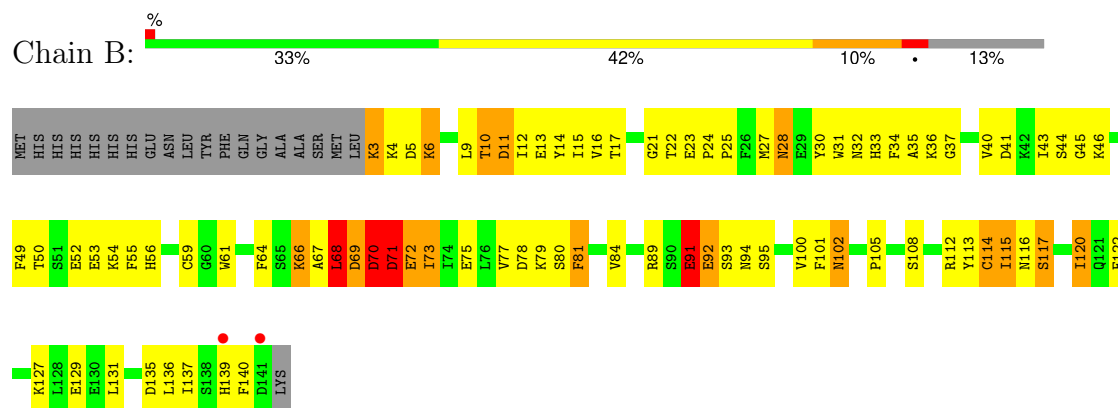
3 Residue-property plots [i](#)

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.

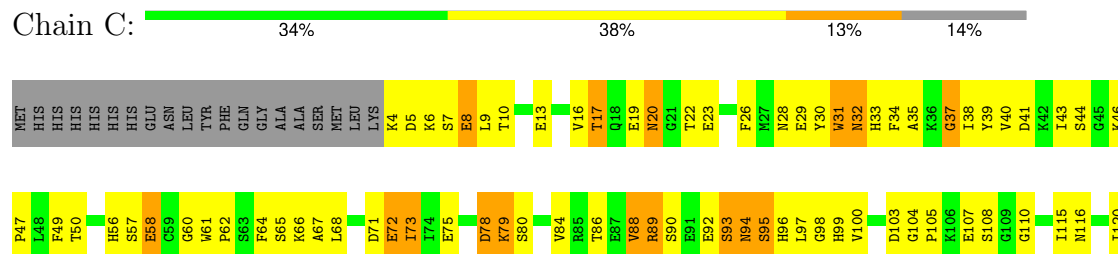
• Molecule 1: Peptide methionine sulfoxide reductase MsrB

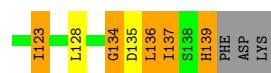


• Molecule 1: Peptide methionine sulfoxide reductase MsrB

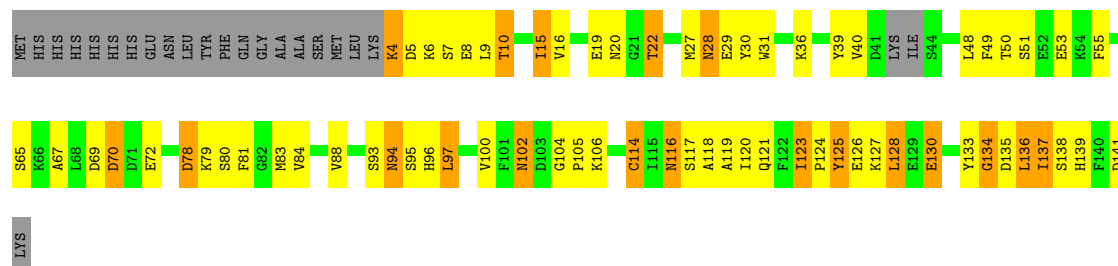


• Molecule 1: Peptide methionine sulfoxide reductase MsrB

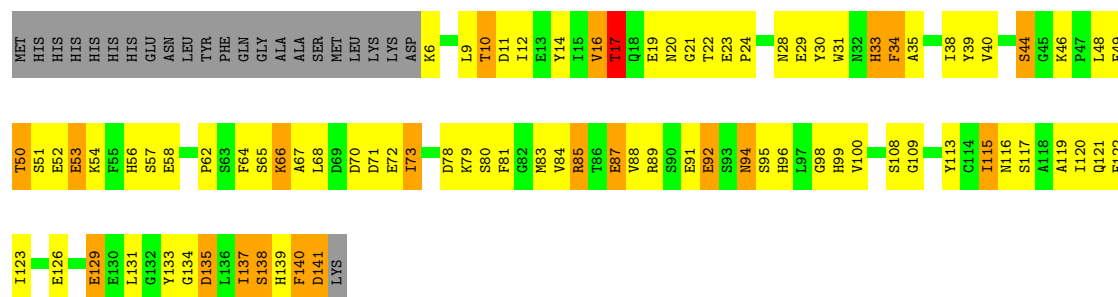




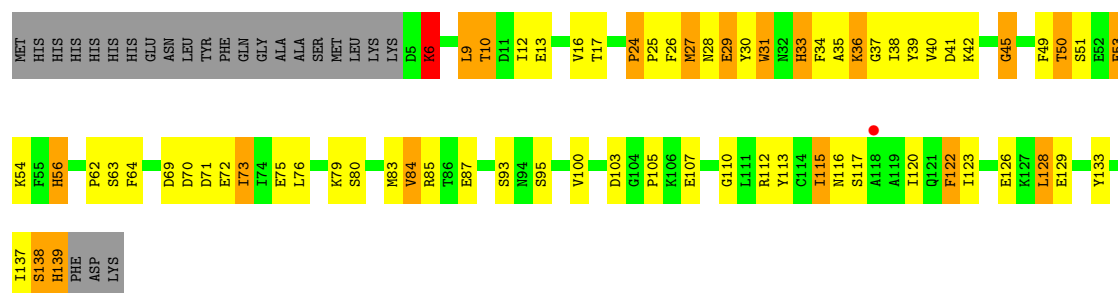
- Molecule 1: Peptide methionine sulfoxide reductase MsrB



- Molecule 1: Peptide methionine sulfoxide reductase MsrB



- Molecule 1: Peptide methionine sulfoxide reductase MsrB



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	114.83Å 122.19Å 73.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	83.68 – 3.47 83.68 – 3.47	Depositor EDS
% Data completeness (in resolution range)	99.8 (83.68-3.47) 99.9 (83.68-3.47)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.35 (at 3.49Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.216 , 0.323 0.230 , 0.327	Depositor DCC
R_{free} test set	686 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	111.2	Xtriage
Anisotropy	0.059	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 110.3	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	0.92	EDS
Total number of atoms	6564	wwPDB-VP
Average B, all atoms (Å ²)	120.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.81% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.15	0/1104	1.46	6/1491 (0.4%)
1	B	1.02	0/1150	1.57	8/1551 (0.5%)
1	C	1.09	2/1121 (0.2%)	1.45	4/1513 (0.3%)
1	D	1.06	0/1123	1.51	4/1515 (0.3%)
1	E	1.08	0/1124	1.49	3/1518 (0.2%)
1	F	1.06	0/1112	1.46	2/1502 (0.1%)
All	All	1.08	2/6734 (0.0%)	1.49	27/9090 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	123	ILE	N-CA	5.12	1.49	1.45
1	C	37	GLY	C-O	5.09	1.27	1.23

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	64	PHE	CA-C-N	6.42	129.21	120.54
1	B	64	PHE	C-N-CA	6.42	129.21	120.54
1	F	122	PHE	CA-C-O	-6.29	114.28	121.19
1	E	16	VAL	CA-C-N	5.95	128.17	120.44
1	E	16	VAL	C-N-CA	5.95	128.17	120.44
1	E	17	THR	CA-CB-OG1	-5.89	100.76	109.60
1	F	24	PRO	N-CA-CB	5.82	106.45	103.19
1	D	124	PRO	N-CA-C	5.76	120.12	111.14
1	A	36	LYS	CA-C-N	5.74	128.92	121.23
1	A	36	LYS	C-N-CA	5.74	128.92	121.23
1	A	69	ASP	CA-CB-CG	5.59	118.19	112.60
1	B	40	VAL	CA-C-O	-5.51	117.32	121.68
1	B	91	GLU	N-CA-C	-5.51	104.92	111.03
1	D	27	MET	CA-C-N	5.38	131.82	121.54
1	D	27	MET	C-N-CA	5.38	131.82	121.54
1	C	32	ASN	CA-C-N	5.38	129.81	121.26

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	32	ASN	C-N-CA	5.38	129.81	121.26
1	B	54	LYS	CA-C-O	-5.32	115.44	121.72
1	B	105	PRO	N-CA-C	-5.26	102.86	110.80
1	A	14	TYR	CB-CA-C	5.21	119.46	110.86
1	C	128	LEU	CA-C-N	5.06	127.06	120.28
1	C	128	LEU	C-N-CA	5.06	127.06	120.28
1	B	70	ASP	CA-C-N	5.04	131.16	121.54
1	B	70	ASP	C-N-CA	5.04	131.16	121.54
1	A	32	ASN	CA-C-N	5.03	128.24	121.05
1	A	32	ASN	C-N-CA	5.03	128.24	121.05
1	D	22	THR	CA-CB-OG1	-5.01	102.09	109.60

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	1076	0	1008	70	0
1	B	1121	0	1051	93	0
1	C	1093	0	1025	76	0
1	D	1095	0	1013	70	0
1	E	1095	0	1021	76	0
1	F	1084	0	1012	58	0
All	All	6564	0	6130	431	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 34.

All (431) 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:E:88:VAL:HG12	1:E:98:GLY:C	1.80	1.07
1:C:49:PHE:CE1	1:C:67:ALA:HB2	1.89	1.06
1:B:46:LYS:HZ3	1:B:72:GLU:HB3	1.17	1.02

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:88:VAL:HG12	1:E:98:GLY:O	1.68	0.92
1:C:89:ARG:NH1	1:C:94:ASN:HD21	1.68	0.91
1:E:23:GLU:HG2	1:E:116:ASN:OD1	1.69	0.91
1:B:46:LYS:NZ	1:B:72:GLU:HB3	1.88	0.89
1:F:36:LYS:H	1:F:36:LYS:HD2	1.39	0.87
1:B:30:TYR:HB2	1:B:117:SER:HB3	1.57	0.86
1:D:20:ASN:HA	1:D:96:HIS:CD2	2.13	0.84
1:F:39:TYR:HD2	1:F:120:ILE:HG21	1.42	0.82
1:D:50:THR:HG23	1:D:53:GLU:HG2	1.61	0.80
1:E:54:LYS:HG3	1:E:64:PHE:CE2	2.17	0.79
1:C:32:ASN:O	1:C:34:PHE:CD1	2.35	0.79
1:F:36:LYS:HD2	1:F:36:LYS:N	1.97	0.79
1:A:72:GLU:O	1:A:91:GLU:CD	2.26	0.79
1:B:69:ASP:OD1	1:B:69:ASP:N	2.15	0.77
1:D:102:ASN:H	1:D:102:ASN:HD22	1.29	0.77
1:C:32:ASN:O	1:C:34:PHE:HD1	1.66	0.76
1:E:78:ASP:HB2	1:E:87:GLU:HB3	1.67	0.76
1:B:28:ASN:HB3	1:B:117:SER:OG	1.86	0.76
1:F:41:ASP:OD2	1:F:93:SER:HB3	1.84	0.76
1:B:66:LYS:HA	1:B:113:TYR:CE1	2.22	0.75
1:D:94:ASN:CG	1:D:94:ASN:O	2.30	0.75
1:D:127:LYS:HD3	1:D:130:GLU:OE1	1.88	0.74
1:F:39:TYR:CD2	1:F:120:ILE:HG21	2.22	0.74
1:F:33:HIS:CE1	1:F:35:ALA:HB3	2.22	0.73
1:D:48:LEU:HD13	1:D:97:LEU:HD11	1.68	0.73
1:D:102:ASN:N	1:D:102:ASN:ND2	2.37	0.73
1:D:123:ILE:O	1:D:128:LEU:HD11	1.89	0.73
1:E:20:ASN:HA	1:E:96:HIS:CD2	2.24	0.73
1:E:46:LYS:HD2	1:E:92:GLU:HG2	1.70	0.72
1:D:127:LYS:O	1:D:130:GLU:HB2	1.89	0.72
1:D:102:ASN:HD22	1:D:102:ASN:N	1.88	0.71
1:B:68:LEU:HD12	1:B:68:LEU:C	2.16	0.71
1:D:50:THR:CG2	1:D:53:GLU:HG2	2.22	0.70
1:D:39:TYR:HB3	1:D:120:ILE:HG23	1.73	0.70
1:D:30:TYR:HD2	1:D:117:SER:HB2	1.55	0.69
1:A:91:GLU:CD	1:A:91:GLU:N	2.51	0.69
1:F:17:THR:HB	1:F:95:SER:HB2	1.73	0.69
1:A:90:SER:O	1:A:94:ASN:HA	1.92	0.68
1:D:30:TYR:CD2	1:D:117:SER:HB2	2.28	0.68
1:E:66:LYS:HG3	1:E:67:ALA:O	1.93	0.68
1:D:118:ALA:C	1:D:120:ILE:H	2.01	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:LYS:HD3	1:C:84:VAL:HG22	1.76	0.68
1:D:28:ASN:O	1:D:30:TYR:N	2.25	0.67
1:B:28:ASN:H	1:B:28:ASN:ND2	1.91	0.67
1:C:46:LYS:NZ	1:C:72:GLU:OE2	2.27	0.67
1:E:49:PHE:CE1	1:E:67:ALA:HA	2.29	0.67
1:D:116:ASN:H	1:D:116:ASN:HD22	1.41	0.67
1:A:128:LEU:HD23	1:A:133:TYR:HB2	1.77	0.67
1:A:134:GLY:O	1:A:137:ILE:HG12	1.95	0.67
1:F:39:TYR:CE2	1:F:122:PHE:HB2	2.30	0.67
1:B:72:GLU:OE1	1:B:72:GLU:HA	1.94	0.67
1:D:28:ASN:HB3	1:D:117:SER:OG	1.95	0.66
1:D:78:ASP:OD1	1:D:78:ASP:C	2.38	0.66
1:D:88:VAL:O	1:D:97:LEU:HB2	1.95	0.66
1:A:22:THR:O	1:E:24:PRO:HB3	1.96	0.66
1:C:40:VAL:HG12	1:C:47:PRO:HA	1.77	0.66
1:D:48:LEU:CD1	1:D:97:LEU:HD11	2.27	0.66
1:A:113:TYR:O	1:A:115:ILE:HG12	1.97	0.65
1:B:72:GLU:O	1:B:91:GLU:HB3	1.97	0.65
1:F:64:PHE:O	1:F:113:TYR:HD1	1.79	0.64
1:D:116:ASN:H	1:D:116:ASN:ND2	1.95	0.64
1:C:10:THR:OG1	1:C:13:GLU:HG3	1.97	0.64
1:F:17:THR:HB	1:F:95:SER:CB	2.28	0.64
1:E:54:LYS:CG	1:E:64:PHE:CE2	2.81	0.63
1:D:40:VAL:HG22	1:D:121:GLN:O	1.98	0.63
1:D:55:PHE:CZ	1:D:105:PRO:CD	2.81	0.63
1:C:89:ARG:NH1	1:C:94:ASN:ND2	2.44	0.63
1:E:23:GLU:CG	1:E:116:ASN:OD1	2.46	0.63
1:C:9:LEU:HD11	1:C:43:ILE:HD11	1.81	0.63
1:D:16:VAL:O	1:D:96:HIS:HB3	1.99	0.63
1:B:17:THR:HB	1:B:95:SER:HB2	1.80	0.63
1:C:17:THR:HG22	1:C:95:SER:HB3	1.80	0.63
1:C:34:PHE:CE1	1:C:56:HIS:CE1	2.87	0.62
1:D:48:LEU:HD13	1:D:97:LEU:CD1	2.29	0.62
1:B:33:HIS:NE2	1:B:122:PHE:CD1	2.66	0.62
1:E:28:ASN:HD21	1:E:62:PRO:HD3	1.65	0.62
1:F:79:LYS:HA	1:F:83:MET:O	1.99	0.62
1:B:28:ASN:CB	1:B:117:SER:OG	2.47	0.62
1:B:37:GLY:O	1:B:140:PHE:HZ	1.83	0.62
1:A:28:ASN:HD21	1:A:62:PRO:HD3	1.66	0.61
1:B:49:PHE:CE1	1:B:67:ALA:HA	2.35	0.61
1:E:28:ASN:ND2	1:E:62:PRO:HD3	2.16	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:30:TYR:HA	1:C:33:HIS:HB3	1.81	0.61
1:E:48:LEU:C	1:E:49:PHE:CD2	2.78	0.61
1:C:94:ASN:ND2	1:C:94:ASN:O	2.32	0.60
1:D:128:LEU:HD13	1:D:128:LEU:H	1.66	0.60
1:A:102:ASN:HD21	1:C:84:VAL:H	1.50	0.60
1:D:127:LYS:CD	1:D:130:GLU:OE1	2.49	0.60
1:E:85:ARG:HH12	1:E:99:HIS:CE1	2.20	0.60
1:D:134:GLY:O	1:D:136:LEU:N	2.35	0.60
1:B:49:PHE:HE2	1:B:115:ILE:HD13	1.65	0.60
1:E:20:ASN:HA	1:E:96:HIS:CG	2.37	0.59
1:C:73:ILE:HD11	1:C:97:LEU:HD12	1.83	0.59
1:B:112:ARG:NH1	1:B:114:CYS:SG	2.76	0.59
1:E:88:VAL:CG1	1:E:98:GLY:C	2.67	0.59
1:A:128:LEU:C	1:A:130:GLU:H	2.11	0.59
1:B:35:ALA:HB3	1:B:122:PHE:CZ	2.37	0.59
1:A:133:TYR:O	1:A:135:ASP:N	2.35	0.59
1:B:84:VAL:HG11	1:E:100:VAL:HG13	1.84	0.59
1:F:30:TYR:CD2	1:F:117:SER:HB2	2.37	0.59
1:F:16:VAL:HG13	1:F:116:ASN:HD22	1.68	0.59
1:C:33:HIS:HD2	1:C:39:TYR:OH	1.86	0.58
1:F:6:LYS:O	1:F:6:LYS:HG2	2.02	0.58
1:B:25:PRO:C	1:B:27:MET:H	2.10	0.58
1:C:28:ASN:HD21	1:C:62:PRO:HD3	1.69	0.58
1:B:17:THR:CB	1:B:95:SER:HB2	2.34	0.58
1:B:66:LYS:HA	1:B:113:TYR:HE1	1.68	0.58
1:B:66:LYS:C	1:B:68:LEU:H	2.12	0.58
1:D:20:ASN:HA	1:D:96:HIS:CG	2.39	0.58
1:A:90:SER:HA	1:A:91:GLU:OE2	2.04	0.58
1:D:100:VAL:HG13	1:F:84:VAL:HG11	1.85	0.58
1:E:54:LYS:HG3	1:E:64:PHE:CD2	2.38	0.58
1:A:9:LEU:HD21	1:A:43:ILE:CD1	2.34	0.58
1:E:39:TYR:HA	1:E:121:GLN:O	2.03	0.58
1:A:72:GLU:O	1:A:91:GLU:OE2	2.22	0.57
1:F:39:TYR:HD2	1:F:120:ILE:CG2	2.17	0.57
1:B:89:ARG:NH2	1:B:94:ASN:HD21	2.02	0.57
1:F:56:HIS:ND1	1:F:56:HIS:N	2.53	0.57
1:A:78:ASP:OD1	1:A:80:SER:HB2	2.04	0.57
1:B:9:LEU:O	1:B:10:THR:C	2.47	0.57
1:B:46:LYS:HZ3	1:B:72:GLU:CB	2.04	0.57
1:A:84:VAL:O	1:A:85:ARG:HB2	2.04	0.57
1:F:37:GLY:HA3	1:F:122:PHE:HE1	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:49:PHE:HB3	1:C:64:PHE:CD2	2.40	0.56
1:C:88:VAL:N	1:C:98:GLY:O	2.38	0.56
1:E:34:PHE:HD1	1:E:34:PHE:N	2.04	0.56
1:B:25:PRO:HG3	1:B:61:TRP:CZ2	2.40	0.56
1:D:104:GLY:O	1:D:105:PRO:C	2.47	0.56
1:D:118:ALA:O	1:D:120:ILE:N	2.38	0.56
1:D:127:LYS:O	1:D:130:GLU:N	2.39	0.56
1:A:41:ASP:OD2	1:A:93:SER:HB3	2.06	0.56
1:B:73:ILE:HD13	1:B:73:ILE:O	2.06	0.56
1:E:50:THR:OG1	1:E:51:SER:N	2.38	0.56
1:E:71:ASP:O	1:E:91:GLU:OE1	2.24	0.56
1:B:67:ALA:O	1:B:69:ASP:N	2.39	0.55
1:C:44:SER:HB3	1:C:93:SER:HB2	1.87	0.55
1:D:4:LYS:N	1:D:7:SER:OG	2.39	0.55
1:F:84:VAL:O	1:F:84:VAL:HG12	2.06	0.55
1:B:41:ASP:OD2	1:B:93:SER:HB3	2.07	0.55
1:D:128:LEU:H	1:D:128:LEU:CD1	2.19	0.55
1:C:46:LYS:HD3	1:C:92:GLU:HB3	1.89	0.55
1:B:9:LEU:O	1:B:11:ASP:N	2.40	0.55
1:E:34:PHE:N	1:E:34:PHE:CD1	2.74	0.55
1:E:88:VAL:HG13	1:E:88:VAL:O	2.06	0.55
1:C:49:PHE:CD1	1:C:67:ALA:HB2	2.39	0.55
1:D:55:PHE:CZ	1:D:105:PRO:HD3	2.40	0.55
1:F:54:LYS:HB2	1:F:64:PHE:CE2	2.42	0.55
1:E:50:THR:CG2	1:E:68:LEU:HD13	2.38	0.54
1:E:17:THR:HG22	1:E:95:SER:HB2	1.88	0.54
1:A:94:ASN:O	1:A:94:ASN:ND2	2.40	0.54
1:A:81:PHE:O	1:A:83:MET:HG3	2.08	0.54
1:B:3:LYS:O	1:B:6:LYS:N	2.40	0.54
1:B:81:PHE:CD1	1:B:81:PHE:N	2.75	0.54
1:A:88:VAL:N	1:A:98:GLY:O	2.41	0.53
1:D:15:ILE:O	1:D:19:GLU:N	2.39	0.53
1:E:88:VAL:HG13	1:E:98:GLY:H	1.73	0.53
1:C:49:PHE:HB3	1:C:64:PHE:CE2	2.43	0.53
1:E:66:LYS:HA	1:E:113:TYR:CE1	2.42	0.53
1:F:62:PRO:HD2	1:F:115:ILE:O	2.08	0.53
1:D:117:SER:O	1:D:120:ILE:HB	2.08	0.53
1:A:28:ASN:HD22	1:A:117:SER:CB	2.21	0.53
1:E:38:ILE:O	1:E:122:PHE:HA	2.08	0.53
1:C:9:LEU:HD21	1:C:43:ILE:CD1	2.39	0.52
1:F:80:SER:HB2	1:F:85:ARG:HH21	1.72	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:93:SER:OG	1:C:95:SER:OG	2.27	0.52
1:A:13:GLU:HA	1:A:118:ALA:HB1	1.92	0.52
1:D:118:ALA:C	1:D:120:ILE:N	2.67	0.52
1:B:35:ALA:CB	1:B:122:PHE:CZ	2.92	0.52
1:B:139:HIS:H	1:B:139:HIS:CD2	2.27	0.52
1:A:84:VAL:HG22	1:C:86:THR:HG21	1.90	0.52
1:A:80:SER:HB3	1:A:81:PHE:CD2	2.45	0.52
1:B:37:GLY:C	1:B:140:PHE:HZ	2.18	0.52
1:B:102:ASN:N	1:B:102:ASN:ND2	2.57	0.52
1:C:90:SER:O	1:C:94:ASN:HA	2.10	0.52
1:F:13:GLU:CD	1:F:42:LYS:HE2	2.35	0.52
1:F:37:GLY:HA3	1:F:122:PHE:CE1	2.45	0.52
1:D:128:LEU:CD1	1:D:128:LEU:N	2.73	0.51
1:E:85:ARG:NH1	1:E:99:HIS:CE1	2.78	0.51
1:A:90:SER:O	1:A:94:ASN:CA	2.58	0.51
1:A:90:SER:O	1:A:94:ASN:N	2.43	0.51
1:F:28:ASN:HB2	1:F:117:SER:OG	2.10	0.51
1:B:10:THR:OG1	1:B:13:GLU:HB2	2.11	0.51
1:F:24:PRO:HB2	1:F:27:MET:HG3	1.92	0.51
1:A:84:VAL:HG13	1:A:85:ARG:N	2.26	0.51
1:D:93:SER:O	1:D:93:SER:OG	2.19	0.51
1:B:5:ASP:CG	1:B:6:LYS:N	2.69	0.50
1:D:36:LYS:HA	1:D:51:SER:CB	2.41	0.50
1:F:105:PRO:C	1:F:107:GLU:H	2.18	0.50
1:D:30:TYR:HB2	1:D:117:SER:CB	2.41	0.50
1:E:29:GLU:HG2	1:E:30:TYR:N	2.26	0.50
1:E:80:SER:HB3	1:E:81:PHE:CD2	2.47	0.50
1:E:30:TYR:OH	1:E:121:GLN:NE2	2.44	0.50
1:B:15:ILE:HG22	1:B:21:GLY:HA3	1.94	0.50
1:C:105:PRO:HD2	1:C:108:SER:OG	2.11	0.50
1:D:114:CYS:O	1:D:114:CYS:SG	2.70	0.50
1:D:39:TYR:HB3	1:D:120:ILE:CG2	2.42	0.50
1:E:141:ASP:OD2	1:E:141:ASP:N	2.44	0.50
1:F:70:ASP:HA	1:F:73:ILE:HD13	1.93	0.50
1:C:49:PHE:CZ	1:C:67:ALA:HB2	2.44	0.50
1:A:125:TYR:HA	1:A:128:LEU:HD12	1.94	0.49
1:B:68:LEU:C	1:B:68:LEU:CD1	2.85	0.49
1:D:20:ASN:OD1	1:D:96:HIS:HB2	2.12	0.49
1:E:123:ILE:HD13	1:E:133:TYR:CE1	2.47	0.49
1:F:69:ASP:HB3	1:F:72:GLU:HG2	1.95	0.49
1:C:72:GLU:C	1:C:72:GLU:CD	2.80	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:128:LEU:O	1:D:133:TYR:HB2	2.12	0.49
1:E:57:SER:O	1:E:58:GLU:C	2.56	0.49
1:F:9:LEU:HB2	1:F:13:GLU:OE2	2.13	0.49
1:F:30:TYR:HA	1:F:33:HIS:HB3	1.95	0.49
1:A:44:SER:HB3	1:A:93:SER:HB3	1.94	0.49
1:C:20:ASN:HA	1:C:96:HIS:CD2	2.48	0.49
1:B:17:THR:HG22	1:B:95:SER:HB2	1.95	0.49
1:D:116:ASN:ND2	1:D:116:ASN:N	2.61	0.49
1:A:31:TRP:C	1:A:31:TRP:CD2	2.90	0.49
1:B:80:SER:HB3	1:B:81:PHE:CE1	2.48	0.49
1:C:61:TRP:O	1:C:62:PRO:C	2.56	0.49
1:F:137:ILE:O	1:F:139:HIS:N	2.46	0.49
1:D:20:ASN:CA	1:D:96:HIS:CD2	2.91	0.48
1:F:36:LYS:N	1:F:36:LYS:CD	2.73	0.48
1:F:45:GLY:O	1:F:133:TYR:OH	2.30	0.48
1:B:80:SER:HB3	1:B:81:PHE:CZ	2.48	0.48
1:C:4:LYS:HA	1:C:8:GLU:OE2	2.14	0.48
1:C:32:ASN:O	1:C:34:PHE:CE1	2.66	0.48
1:E:94:ASN:ND2	1:E:94:ASN:O	2.46	0.48
1:B:13:GLU:O	1:B:17:THR:OG1	2.28	0.48
1:A:94:ASN:O	1:A:94:ASN:CG	2.55	0.48
1:B:3:LYS:HA	1:B:3:LYS:HE2	1.95	0.48
1:C:72:GLU:O	1:C:72:GLU:OE1	2.31	0.48
1:A:30:TYR:HA	1:A:33:HIS:HB3	1.96	0.48
1:B:46:LYS:NZ	1:B:72:GLU:CB	2.71	0.47
1:E:137:ILE:O	1:E:139:HIS:N	2.47	0.47
1:A:25:PRO:HD2	1:E:22:THR:HG21	1.96	0.47
1:B:44:SER:HB3	1:B:93:SER:HB2	1.96	0.47
1:D:79:LYS:HA	1:D:83:MET:O	2.15	0.47
1:A:133:TYR:O	1:A:134:GLY:C	2.57	0.47
1:B:3:LYS:O	1:B:5:ASP:N	2.48	0.47
1:C:33:HIS:NE2	1:C:35:ALA:HB3	2.30	0.47
1:E:67:ALA:CB	1:E:73:ILE:HD11	2.45	0.47
1:F:25:PRO:HB2	1:F:26:PHE:CD2	2.49	0.47
1:A:73:ILE:HA	1:A:91:GLU:OE2	2.14	0.47
1:B:23:GLU:OE1	1:B:117:SER:N	2.48	0.47
1:A:31:TRP:O	1:A:31:TRP:CE3	2.68	0.47
1:B:10:THR:O	1:B:11:ASP:HB2	2.14	0.47
1:B:11:ASP:O	1:B:14:TYR:N	2.47	0.47
1:C:89:ARG:HA	1:C:97:LEU:HG	1.97	0.47
1:C:104:GLY:H	1:C:110:GLY:C	2.23	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:28:ASN:OD1	1:B:61:TRP:HD1	1.98	0.47
1:C:6:LYS:C	1:C:8:GLU:H	2.22	0.47
1:E:19:GLU:O	1:E:20:ASN:C	2.57	0.47
1:E:30:TYR:HB2	1:E:117:SER:HB2	1.96	0.47
1:A:84:VAL:HG21	1:C:100:VAL:HG13	1.96	0.47
1:E:94:ASN:O	1:E:94:ASN:CG	2.57	0.47
1:A:78:ASP:OD1	1:A:80:SER:CB	2.63	0.46
1:A:125:TYR:HA	1:A:128:LEU:CD1	2.46	0.46
1:B:100:VAL:HG13	1:E:84:VAL:HG21	1.97	0.46
1:C:73:ILE:HD13	1:C:73:ILE:HA	1.62	0.46
1:B:23:GLU:HG2	1:B:116:ASN:OD1	2.14	0.46
1:C:19:GLU:C	1:C:20:ASN:HD22	2.23	0.46
1:F:33:HIS:CD2	1:F:39:TYR:OH	2.68	0.46
1:E:78:ASP:OD1	1:E:80:SER:CB	2.64	0.46
1:E:20:ASN:OD1	1:E:89:ARG:NH1	2.48	0.46
1:A:27:MET:HE2	1:E:21:GLY:HA2	1.97	0.46
1:A:91:GLU:CD	1:A:91:GLU:H	2.22	0.46
1:B:102:ASN:N	1:B:102:ASN:HD22	2.14	0.46
1:B:139:HIS:CD2	1:B:139:HIS:N	2.83	0.46
1:C:41:ASP:HB3	1:C:44:SER:OG	2.16	0.46
1:A:8:GLU:O	1:A:8:GLU:HG3	2.16	0.46
1:B:17:THR:CG2	1:B:95:SER:HB2	2.46	0.46
1:B:115:ILE:HD11	1:B:120:ILE:HG23	1.97	0.46
1:F:69:ASP:HB3	1:F:72:GLU:CG	2.46	0.46
1:B:25:PRO:C	1:B:27:MET:N	2.71	0.46
1:B:66:LYS:C	1:B:68:LEU:N	2.72	0.46
1:A:37:GLY:O	1:A:51:SER:HB3	2.15	0.46
1:C:93:SER:HG	1:C:95:SER:HG	1.59	0.45
1:B:37:GLY:O	1:B:140:PHE:CZ	2.68	0.45
1:C:89:ARG:HH11	1:C:94:ASN:HD21	1.54	0.45
1:A:41:ASP:HB2	1:A:48:LEU:HD11	1.97	0.45
1:B:66:LYS:HA	1:B:113:TYR:CD1	2.50	0.45
1:C:17:THR:CG2	1:C:95:SER:HB3	2.46	0.45
1:C:29:GLU:HG2	1:C:30:TYR:H	1.82	0.45
1:A:49:PHE:CD1	1:A:67:ALA:HA	2.51	0.45
1:A:72:GLU:C	1:A:91:GLU:CG	2.90	0.45
1:B:35:ALA:HB3	1:B:122:PHE:CE1	2.50	0.45
1:C:38:ILE:HG12	1:C:123:ILE:HB	1.99	0.45
1:D:70:ASP:OD1	1:D:70:ASP:N	2.49	0.45
1:F:10:THR:CG2	1:F:12:ILE:HG22	2.46	0.45
1:A:135:ASP:O	1:A:137:ILE:N	2.49	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:31:TRP:O	1:B:31:TRP:CE3	2.69	0.45
1:E:49:PHE:HA	1:E:68:LEU:HB2	1.99	0.45
1:B:41:ASP:OD2	1:B:93:SER:CB	2.64	0.45
1:D:4:LYS:N	1:D:7:SER:HG	2.14	0.45
1:F:29:GLU:HG2	1:F:30:TYR:CE1	2.51	0.45
1:E:20:ASN:OD1	1:E:20:ASN:N	2.47	0.45
1:B:49:PHE:CE2	1:B:115:ILE:HD13	2.50	0.45
1:B:129:GLU:OE1	1:B:137:ILE:CD1	2.64	0.45
1:D:55:PHE:CZ	1:D:105:PRO:HD2	2.51	0.45
1:E:10:THR:OG1	1:E:12:ILE:HG22	2.15	0.45
1:D:31:TRP:O	1:D:31:TRP:CE3	2.70	0.45
1:F:128:LEU:HD12	1:F:128:LEU:HA	1.81	0.45
1:C:9:LEU:HD21	1:C:43:ILE:HD11	1.99	0.44
1:C:47:PRO:HG2	1:C:136:LEU:HD11	1.99	0.44
1:C:137:ILE:HG22	1:C:137:ILE:O	2.16	0.44
1:D:137:ILE:HD12	1:D:138:SER:N	2.33	0.44
1:E:88:VAL:CG1	1:E:98:GLY:H	2.30	0.44
1:A:84:VAL:O	1:A:85:ARG:CB	2.65	0.44
1:B:27:MET:HB3	1:B:27:MET:HE3	1.70	0.44
1:E:50:THR:C	1:E:53:GLU:OE1	2.61	0.44
1:E:129:GLU:HG3	1:E:134:GLY:HA3	1.98	0.44
1:C:19:GLU:C	1:C:20:ASN:ND2	2.75	0.44
1:E:50:THR:HG22	1:E:68:LEU:HD13	1.98	0.44
1:C:4:LYS:N	1:C:8:GLU:HB3	2.33	0.44
1:F:54:LYS:HA	1:F:63:SER:O	2.18	0.44
1:A:84:VAL:CG2	1:C:86:THR:HG21	2.47	0.44
1:D:49:PHE:CD1	1:D:67:ALA:HA	2.52	0.44
1:E:30:TYR:HD2	1:E:117:SER:HB2	1.82	0.44
1:E:46:LYS:HD2	1:E:92:GLU:CG	2.46	0.44
1:E:53:GLU:CD	1:E:53:GLU:H	2.26	0.44
1:C:72:GLU:OE1	1:C:90:SER:HA	2.19	0.43
1:D:9:LEU:C	1:D:10:THR:O	2.60	0.43
1:E:17:THR:HG23	1:E:119:ALA:HB2	2.00	0.43
1:E:28:ASN:HB2	1:E:117:SER:OG	2.18	0.43
1:F:100:VAL:HA	1:F:112:ARG:O	2.18	0.43
1:A:17:THR:HB	1:A:18:GLN:HE21	1.83	0.43
1:A:49:PHE:CE1	1:A:67:ALA:HA	2.53	0.43
1:B:139:HIS:H	1:B:139:HIS:HD2	1.66	0.43
1:A:23:GLU:CD	1:A:117:SER:H	2.25	0.43
1:E:78:ASP:HB2	1:E:87:GLU:CB	2.45	0.43
1:F:75:GLU:HA	1:F:87:GLU:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:SER:O	1:A:139:HIS:C	2.61	0.43
1:B:43:ILE:HB	1:B:93:SER:OG	2.19	0.43
1:C:115:ILE:HG22	1:C:116:ASN:O	2.17	0.43
1:C:17:THR:HG22	1:C:95:SER:CB	2.46	0.43
1:B:33:HIS:CD2	1:B:122:PHE:CD1	3.07	0.43
1:B:66:LYS:O	1:B:68:LEU:N	2.51	0.43
1:E:54:LYS:HD2	1:E:64:PHE:CZ	2.53	0.43
1:F:54:LYS:HB2	1:F:64:PHE:CZ	2.54	0.43
1:A:24:PRO:HB3	1:E:22:THR:O	2.19	0.43
1:B:33:HIS:CE1	1:B:35:ALA:CB	3.02	0.43
1:B:5:ASP:CG	1:B:6:LYS:H	2.27	0.43
1:C:23:GLU:OE1	1:C:116:ASN:HA	2.19	0.43
1:C:57:SER:O	1:C:58:GLU:HB2	2.19	0.43
1:D:30:TYR:HB2	1:D:117:SER:HB3	2.01	0.43
1:E:31:TRP:CD2	1:E:31:TRP:C	2.97	0.43
1:C:26:PHE:HB3	1:D:81:PHE:CZ	2.54	0.43
1:B:78:ASP:OD2	1:B:78:ASP:C	2.62	0.42
1:D:93:SER:O	1:D:94:ASN:C	2.62	0.42
1:E:62:PRO:HD2	1:E:115:ILE:O	2.19	0.42
1:F:17:THR:CB	1:F:95:SER:HB2	2.44	0.42
1:A:41:ASP:OD2	1:A:93:SER:CB	2.68	0.42
1:A:57:SER:O	1:A:58:GLU:C	2.62	0.42
1:C:44:SER:C	1:C:46:LYS:N	2.77	0.42
1:D:4:LYS:O	1:D:8:GLU:HB3	2.19	0.42
1:B:36:LYS:N	1:B:36:LYS:HD3	2.34	0.42
1:F:30:TYR:CD1	1:F:33:HIS:CD2	3.08	0.42
1:F:49:PHE:HB3	1:F:64:PHE:CD2	2.54	0.42
1:C:26:PHE:CE1	1:C:60:GLY:HA2	2.53	0.42
1:D:30:TYR:HB2	1:D:117:SER:HB2	1.99	0.42
1:D:69:ASP:HB3	1:D:72:GLU:CG	2.50	0.42
1:F:31:TRP:CD2	1:F:31:TRP:C	2.97	0.42
1:F:50:THR:HG23	1:F:53:GLU:CG	2.49	0.42
1:F:84:VAL:O	1:F:84:VAL:CG1	2.67	0.42
1:A:39:TYR:HB3	1:A:120:ILE:CG2	2.49	0.42
1:C:33:HIS:CD2	1:C:39:TYR:OH	2.69	0.42
1:C:134:GLY:C	1:C:136:LEU:N	2.78	0.42
1:A:39:TYR:HB3	1:A:120:ILE:HG21	2.01	0.42
1:A:47:PRO:HG2	1:A:136:LEU:HD11	2.02	0.42
1:B:23:GLU:HB2	1:B:24:PRO:HD2	2.02	0.42
1:B:28:ASN:ND2	1:B:28:ASN:N	2.63	0.42
1:C:33:HIS:CE1	1:C:35:ALA:HB3	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:40:VAL:HG13	1:E:123:ILE:HD12	2.02	0.42
1:A:49:PHE:HA	1:A:68:LEU:HB2	2.01	0.42
1:C:20:ASN:ND2	1:C:20:ASN:N	2.67	0.42
1:A:73:ILE:HD13	1:A:73:ILE:H	1.85	0.42
1:D:69:ASP:OD1	1:D:70:ASP:N	2.53	0.42
1:D:81:PHE:C	1:D:83:MET:H	2.28	0.42
1:C:5:ASP:O	1:C:9:LEU:N	2.44	0.42
1:C:89:ARG:CB	1:C:95:SER:O	2.68	0.42
1:E:79:LYS:CD	1:E:84:VAL:HG12	2.50	0.42
1:F:137:ILE:O	1:F:138:SER:C	2.63	0.42
1:B:70:ASP:OD2	1:B:70:ASP:N	2.53	0.41
1:A:16:VAL:HG12	1:A:17:THR:N	2.35	0.41
1:A:69:ASP:O	1:A:70:ASP:C	2.63	0.41
1:B:53:GLU:HA	1:B:53:GLU:OE2	2.19	0.41
1:C:44:SER:HB2	1:C:46:LYS:HB2	2.01	0.41
1:F:25:PRO:HB2	1:F:26:PHE:HD2	1.85	0.41
1:F:33:HIS:HE1	1:F:35:ALA:HB3	1.77	0.41
1:A:17:THR:HB	1:A:95:SER:HB2	2.03	0.41
1:B:33:HIS:NE2	1:B:35:ALA:HB3	2.35	0.41
1:E:33:HIS:NE2	1:E:35:ALA:HB3	2.35	0.41
1:F:39:TYR:CD2	1:F:122:PHE:HB2	2.56	0.41
1:A:87:GLU:OE2	1:A:89:ARG:NE	2.49	0.41
1:B:33:HIS:CE1	1:B:35:ALA:HB3	2.54	0.41
1:D:136:LEU:HD12	1:D:136:LEU:HA	1.96	0.41
1:B:101:PHE:CD2	1:B:112:ARG:NH2	2.88	0.41
1:E:140:PHE:O	1:E:141:ASP:C	2.63	0.41
1:F:16:VAL:HG13	1:F:116:ASN:ND2	2.36	0.41
1:A:28:ASN:ND2	1:A:117:SER:OG	2.44	0.41
1:C:139:HIS:CD2	1:C:139:HIS:N	2.86	0.41
1:F:37:GLY:HA3	1:F:123:ILE:O	2.20	0.41
1:B:70:ASP:HB2	1:B:71:ASP:H	1.45	0.41
1:C:78:ASP:OD2	1:C:78:ASP:C	2.63	0.41
1:E:48:LEU:O	1:E:49:PHE:CD2	2.73	0.41
1:F:9:LEU:HD13	1:F:9:LEU:N	2.36	0.41
1:A:32:ASN:O	1:A:34:PHE:HD1	2.04	0.41
1:B:32:ASN:HD21	1:B:56:HIS:CE1	2.39	0.41
1:B:68:LEU:HD12	1:B:68:LEU:O	2.21	0.41
1:C:37:GLY:HA3	1:C:123:ILE:O	2.21	0.41
1:C:66:LYS:O	1:C:66:LYS:HG3	2.20	0.41
1:D:36:LYS:HA	1:D:51:SER:OG	2.20	0.41
1:E:126:GLU:CD	1:E:126:GLU:H	2.29	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:125:TYR:C	1:D:125:TYR:CD1	2.97	0.41
1:E:44:SER:OG	1:E:92:GLU:HG3	2.21	0.41
1:E:50:THR:HG1	1:E:53:GLU:CD	2.28	0.41
1:B:127:LYS:O	1:B:131:LEU:HD12	2.21	0.40
1:D:4:LYS:HG3	1:D:7:SER:OG	2.21	0.40
1:E:134:GLY:O	1:E:135:ASP:C	2.64	0.40
1:B:34:PHE:HE2	1:B:55:PHE:HA	1.87	0.40
1:F:34:PHE:HA	1:F:54:LYS:HZ2	1.86	0.40
1:A:73:ILE:HD13	1:A:73:ILE:N	2.36	0.40
1:A:89:ARG:NH2	1:A:94:ASN:O	2.51	0.40
1:B:77:VAL:HG12	1:B:78:ASP:N	2.35	0.40
1:C:28:ASN:HD21	1:C:31:TRP:HB3	1.86	0.40
1:E:33:HIS:C	1:E:34:PHE:CD1	2.99	0.40
1:B:115:ILE:HD11	1:B:120:ILE:CG2	2.51	0.40
1:A:101:PHE:O	1:A:112:ARG:N	2.51	0.40
1:B:44:SER:HB2	1:B:92:GLU:HG3	2.03	0.40
1:B:49:PHE:CD1	1:B:67:ALA:HA	2.56	0.40
1:C:31:TRP:C	1:C:31:TRP:CE3	3.00	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	132/159 (83%)	101 (76%)	23 (17%)	8 (6%)	1	12
1	B	137/159 (86%)	103 (75%)	26 (19%)	8 (6%)	1	12
1	C	134/159 (84%)	92 (69%)	35 (26%)	7 (5%)	1	14
1	D	132/159 (83%)	99 (75%)	21 (16%)	12 (9%)	0	7
1	E	134/159 (84%)	99 (74%)	28 (21%)	7 (5%)	1	14
1	F	133/159 (84%)	107 (80%)	19 (14%)	7 (5%)	1	14

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	802/954 (84%)	601 (75%)	152 (19%)	49 (6%)	1	12

All (49) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	7	SER
1	A	136	LEU
1	B	4	LYS
1	B	10	THR
1	B	11	ASP
1	B	68	LEU
1	B	71	ASP
1	D	28	ASN
1	D	29	GLU
1	D	125	TYR
1	D	134	GLY
1	D	135	ASP
1	E	135	ASP
1	E	138	SER
1	E	140	PHE
1	F	27	MET
1	F	138	SER
1	A	58	GLU
1	A	85	ARG
1	A	134	GLY
1	B	12	ILE
1	B	92	GLU
1	C	72	GLU
1	C	93	SER
1	C	134	GLY
1	C	135	ASP
1	D	10	THR
1	D	94	ASN
1	D	119	ALA
1	D	139	HIS
1	E	33	HIS
1	F	45	GLY
1	F	110	GLY
1	A	129	GLU
1	C	58	GLU
1	F	6	LYS
1	F	33	HIS

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Mol	Chain	Res	Type
1	C	103	ASP
1	D	15	ILE
1	D	80	SER
1	D	130	GLU
1	E	70	ASP
1	E	109	GLY
1	F	51	SER
1	A	114	CYS
1	A	72	GLU
1	C	16	VAL
1	E	16	VAL
1	B	45	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	119/141 (84%)	100 (84%)	19 (16%)	2	14
1	B	124/141 (88%)	97 (78%)	27 (22%)	1	6
1	C	121/141 (86%)	96 (79%)	25 (21%)	1	6
1	D	121/141 (86%)	101 (84%)	20 (16%)	2	13
1	E	121/141 (86%)	92 (76%)	29 (24%)	1	4
1	F	120/141 (85%)	99 (82%)	21 (18%)	2	11
All	All	726/846 (86%)	585 (81%)	141 (19%)	1	8

All (141) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	9	LEU
1	A	11	ASP
1	A	28	ASN
1	A	29	GLU
1	A	65	SER

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Mol	Chain	Res	Type
1	A	70	ASP
1	A	72	GLU
1	A	73	ILE
1	A	84	VAL
1	A	85	ARG
1	A	90	SER
1	A	91	GLU
1	A	116	ASN
1	A	120	ILE
1	A	128	LEU
1	A	137	ILE
1	A	138	SER
1	A	139	HIS
1	B	3	LYS
1	B	6	LYS
1	B	16	VAL
1	B	22	THR
1	B	28	ASN
1	B	50	THR
1	B	52	GLU
1	B	59	CYS
1	B	66	LYS
1	B	68	LEU
1	B	69	ASP
1	B	70	ASP
1	B	71	ASP
1	B	72	GLU
1	B	73	ILE
1	B	75	GLU
1	B	79	LYS
1	B	81	PHE
1	B	91	GLU
1	B	102	ASN
1	B	108	SER
1	B	114	CYS
1	B	115	ILE
1	B	117	SER
1	B	120	ILE
1	B	135	ASP
1	B	136	LEU
1	C	7	SER
1	C	8	GLU

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Mol	Chain	Res	Type
1	C	17	THR
1	C	20	ASN
1	C	22	THR
1	C	31	TRP
1	C	50	THR
1	C	65	SER
1	C	68	LEU
1	C	71	ASP
1	C	73	ILE
1	C	75	GLU
1	C	78	ASP
1	C	79	LYS
1	C	80	SER
1	C	88	VAL
1	C	89	ARG
1	C	94	ASN
1	C	95	SER
1	C	99	HIS
1	C	107	GLU
1	C	120	ILE
1	C	136	LEU
1	C	137	ILE
1	C	139	HIS
1	D	4	LYS
1	D	5	ASP
1	D	6	LYS
1	D	22	THR
1	D	65	SER
1	D	70	ASP
1	D	78	ASP
1	D	84	VAL
1	D	95	SER
1	D	97	LEU
1	D	102	ASN
1	D	106	LYS
1	D	114	CYS
1	D	116	ASN
1	D	123	ILE
1	D	126	GLU
1	D	128	LEU
1	D	136	LEU
1	D	137	ILE

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Mol	Chain	Res	Type
1	D	141	ASP
1	E	6	LYS
1	E	9	LEU
1	E	10	THR
1	E	11	ASP
1	E	14	TYR
1	E	17	THR
1	E	34	PHE
1	E	44	SER
1	E	50	THR
1	E	52	GLU
1	E	53	GLU
1	E	56	HIS
1	E	65	SER
1	E	66	LYS
1	E	72	GLU
1	E	73	ILE
1	E	83	MET
1	E	85	ARG
1	E	87	GLU
1	E	92	GLU
1	E	94	ASN
1	E	108	SER
1	E	115	ILE
1	E	120	ILE
1	E	129	GLU
1	E	131	LEU
1	E	137	ILE
1	E	138	SER
1	E	141	ASP
1	F	6	LYS
1	F	9	LEU
1	F	10	THR
1	F	29	GLU
1	F	31	TRP
1	F	36	LYS
1	F	38	ILE
1	F	40	VAL
1	F	50	THR
1	F	53	GLU
1	F	56	HIS
1	F	71	ASP

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Mol	Chain	Res	Type
1	F	73	ILE
1	F	76	LEU
1	F	84	VAL
1	F	103	ASP
1	F	115	ILE
1	F	126	GLU
1	F	128	LEU
1	F	129	GLU
1	F	139	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	28	ASN
1	A	33	HIS
1	A	94	ASN
1	A	102	ASN
1	B	28	ASN
1	B	32	ASN
1	B	56	HIS
1	B	94	ASN
1	B	139	HIS
1	C	20	ASN
1	C	28	ASN
1	C	33	HIS
1	C	94	ASN
1	C	102	ASN
1	C	121	GLN
1	C	139	HIS
1	D	96	HIS
1	D	99	HIS
1	D	116	ASN
1	E	33	HIS
1	E	94	ASN
1	E	121	GLN
1	F	32	ASN
1	F	33	HIS
1	F	94	ASN
1	F	99	HIS
1	F	116	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	134/159 (84%)	-0.38	0 100 100	71, 115, 152, 187	0
1	B	139/159 (87%)	-0.32	2 (1%) 73 49	64, 97, 141, 178	0
1	C	136/159 (85%)	-0.48	0 100 100	71, 107, 145, 173	0
1	D	136/159 (85%)	-0.25	0 100 100	70, 129, 177, 236	0
1	E	136/159 (85%)	-0.39	0 100 100	80, 137, 173, 187	0
1	F	135/159 (84%)	-0.36	1 (0%) 84 63	79, 126, 171, 205	0
All	All	816/954 (85%)	-0.36	3 (0%) 88 73	64, 116, 169, 236	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	139	HIS	2.6
1	B	141	ASP	2.5
1	F	118	ALA	2.2

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.