



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2023 – 12:38 pm GMT

PDB ID : 8C7E  
Title : Serendipitous cyanase structure from Serratia-like contamination  
Authors : Bloch, Y.; Skladanowska, K.; Savvides, S.N.  
Deposited on : 2023-01-14  
Resolution : 2.80 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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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 : 2.32.1  
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) : 2.32.1

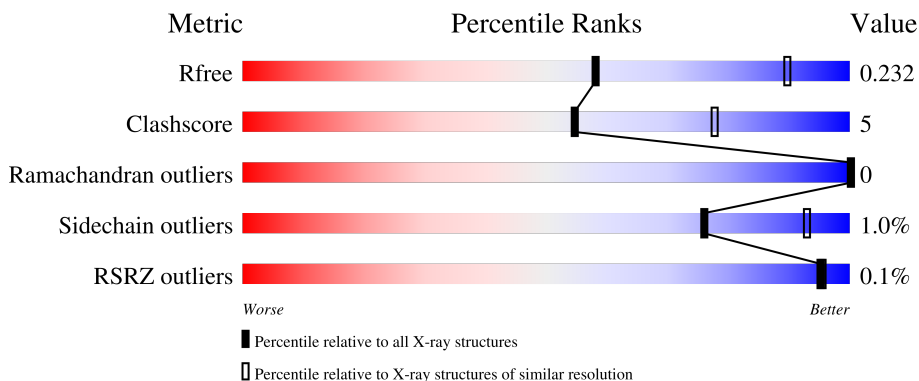
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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  $\geq 3$ , 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	155	88% (green), 12% (yellow)
1	B	155	81% (green), 19% (yellow), 0% (orange), 0% (red)
1	C	155	88% (green), 11% (yellow), 0% (orange), 0% (red)
1	D	155	90% (green), 10% (yellow), 0% (orange), 0% (red)
1	E	155	86% (green), 13% (yellow), 0% (orange), 0% (red)

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Mol	Chain	Length	Quality of chain
1	F	155	 87% 12%
1	G	155	 89% 11%
1	H	155	 85% 15%
1	I	155	 90% 10%
1	J	155	 92% 8%
1	K	155	 91% 9%
1	L	155	 % 83% 16%
1	M	155	 87% 12%
1	N	155	 83% 17%
1	O	155	 82% 18%
1	P	155	 86% 14%
1	Q	155	 80% 19%
1	R	155	 89% 11%
1	S	155	 % 86% 13%
1	T	155	 87% 13%

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 23339 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 Cyanate hydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	155	1164	748	198	216	2	0	0	0
1	B	155	1164	748	198	216	2	0	0	0
1	C	155	1164	748	198	216	2	0	0	0
1	D	155	1164	748	198	216	2	0	0	0
1	E	155	1164	748	198	216	2	0	0	0
1	F	155	1164	748	198	216	2	0	0	0
1	G	155	1164	748	198	216	2	0	0	0
1	H	155	1164	748	198	216	2	0	0	0
1	I	155	1164	748	198	216	2	0	0	0
1	J	155	1168	750	198	218	2	0	0	0
1	K	155	1164	748	198	216	2	0	0	0
1	L	155	1164	748	198	216	2	0	0	0
1	M	155	1168	750	198	218	2	0	0	0
1	N	155	1164	748	198	216	2	0	0	0
1	O	155	1168	750	198	218	2	0	0	0
1	P	155	1164	748	198	216	2	0	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	Q	155	1164	748	198	216	2	0	0	0
1	R	155	1164	748	198	216	2	0	0	0
1	S	155	1168	750	198	218	2	0	0	0
1	T	155	1164	748	198	216	2	0	0	0

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	34	GLN	GLU	conflict	UNP A0A3E2EMR4
A	73	LEU	GLN	conflict	UNP A0A3E2EMR4
B	34	GLN	GLU	conflict	UNP A0A3E2EMR4
B	73	LEU	GLN	conflict	UNP A0A3E2EMR4
C	34	GLN	GLU	conflict	UNP A0A3E2EMR4
C	73	LEU	GLN	conflict	UNP A0A3E2EMR4
D	34	GLN	GLU	conflict	UNP A0A3E2EMR4
D	73	LEU	GLN	conflict	UNP A0A3E2EMR4
E	34	GLN	GLU	conflict	UNP A0A3E2EMR4
E	73	LEU	GLN	conflict	UNP A0A3E2EMR4
F	34	GLN	GLU	conflict	UNP A0A3E2EMR4
F	73	LEU	GLN	conflict	UNP A0A3E2EMR4
G	34	GLN	GLU	conflict	UNP A0A3E2EMR4
G	73	LEU	GLN	conflict	UNP A0A3E2EMR4
H	34	GLN	GLU	conflict	UNP A0A3E2EMR4
H	73	LEU	GLN	conflict	UNP A0A3E2EMR4
I	34	GLN	GLU	conflict	UNP A0A3E2EMR4
I	73	LEU	GLN	conflict	UNP A0A3E2EMR4
J	34	GLN	GLU	conflict	UNP A0A3E2EMR4
J	73	LEU	GLN	conflict	UNP A0A3E2EMR4
K	34	GLN	GLU	conflict	UNP A0A3E2EMR4
K	73	LEU	GLN	conflict	UNP A0A3E2EMR4
L	34	GLN	GLU	conflict	UNP A0A3E2EMR4
L	73	LEU	GLN	conflict	UNP A0A3E2EMR4
M	34	GLN	GLU	conflict	UNP A0A3E2EMR4
M	73	LEU	GLN	conflict	UNP A0A3E2EMR4
N	34	GLN	GLU	conflict	UNP A0A3E2EMR4
N	73	LEU	GLN	conflict	UNP A0A3E2EMR4
O	34	GLN	GLU	conflict	UNP A0A3E2EMR4
O	73	LEU	GLN	conflict	UNP A0A3E2EMR4
P	34	GLN	GLU	conflict	UNP A0A3E2EMR4

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Chain	Residue	Modelled	Actual	Comment	Reference
P	73	LEU	GLN	conflict	UNP A0A3E2EMR4
Q	34	GLN	GLU	conflict	UNP A0A3E2EMR4
Q	73	LEU	GLN	conflict	UNP A0A3E2EMR4
R	34	GLN	GLU	conflict	UNP A0A3E2EMR4
R	73	LEU	GLN	conflict	UNP A0A3E2EMR4
S	34	GLN	GLU	conflict	UNP A0A3E2EMR4
S	73	LEU	GLN	conflict	UNP A0A3E2EMR4
T	34	GLN	GLU	conflict	UNP A0A3E2EMR4
T	73	LEU	GLN	conflict	UNP A0A3E2EMR4

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Cl 2 2	0	0
2	C	2	Total Cl 2 2	0	0
2	D	1	Total Cl 1 1	0	0
2	E	1	Total Cl 1 1	0	0
2	F	1	Total Cl 1 1	0	0
2	G	1	Total Cl 1 1	0	0
2	I	1	Total Cl 1 1	0	0
2	J	1	Total Cl 1 1	0	0
2	K	1	Total Cl 1 1	0	0
2	L	1	Total Cl 1 1	0	0
2	M	2	Total Cl 2 2	0	0
2	N	1	Total Cl 1 1	0	0
2	O	2	Total Cl 2 2	0	0
2	P	1	Total Cl 1 1	0	0
2	R	1	Total Cl 1 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	S	1	Total Cl 1 1	0	0

- Molecule 3 is water.

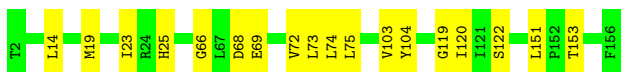
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O 1 1	0	0
3	B	1	Total O 1 1	0	0
3	C	2	Total O 2 2	0	0
3	D	2	Total O 2 2	0	0
3	E	1	Total O 1 1	0	0
3	H	2	Total O 2 2	0	0
3	I	2	Total O 2 2	0	0
3	J	2	Total O 2 2	0	0
3	K	4	Total O 4 4	0	0
3	L	2	Total O 2 2	0	0
3	M	1	Total O 1 1	0	0
3	O	1	Total O 1 1	0	0
3	Q	1	Total O 1 1	0	0
3	T	1	Total O 1 1	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: Cyanate hydratase

Chain A:  88% 12%




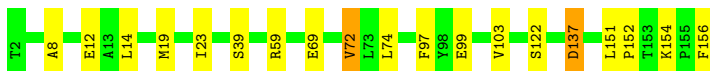
- Molecule 1: Cyanate hydratase

Chain B:  81% 19%




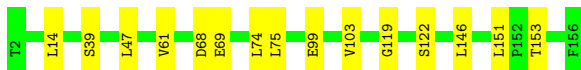
- Molecule 1: Cyanate hydratase

Chain C:  88% 11%




- Molecule 1: Cyanate hydratase

Chain D:  90% 10%



- Molecule 1: Cyanate hydratase

Chain E:  86% 13%



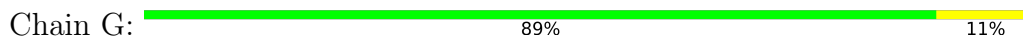
- Molecule 1: Cyanate hydratase

Chain F:  87% 12%

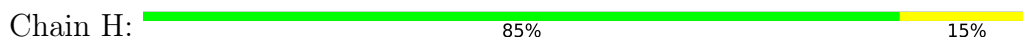




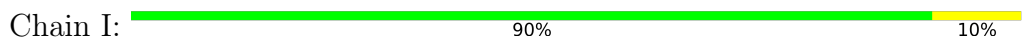
- Molecule 1: Cyanate hydratase



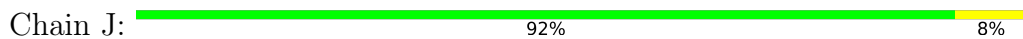
- Molecule 1: Cyanate hydratase



- Molecule 1: Cyanate hydratase



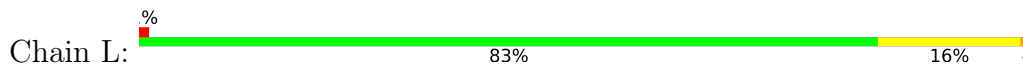
- Molecule 1: Cyanate hydratase



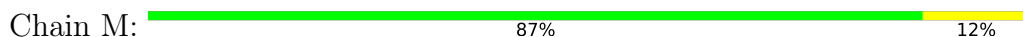
- Molecule 1: Cyanate hydratase



- Molecule 1: Cyanate hydratase

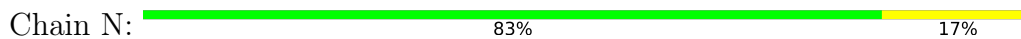


- Molecule 1: Cyanate hydratase

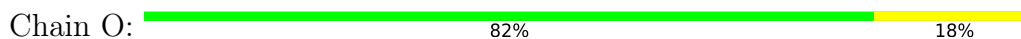




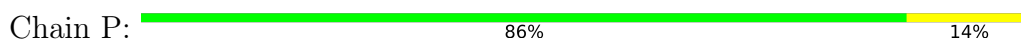
- Molecule 1: Cyanate hydratase



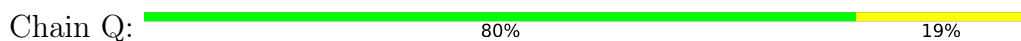
- Molecule 1: Cyanate hydratase



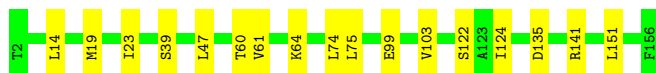
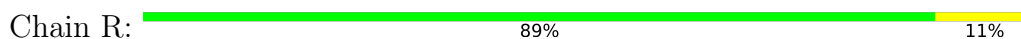
- Molecule 1: Cyanate hydratase



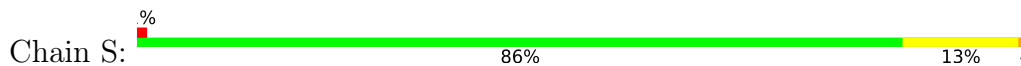
- Molecule 1: Cyanate hydratase



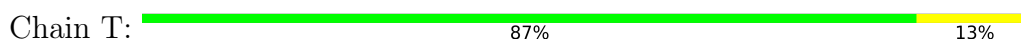
- Molecule 1: Cyanate hydratase



- Molecule 1: Cyanate hydratase



- Molecule 1: Cyanate hydratase





## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.12Å 90.12Å 736.36Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	78.05 – 2.80 78.05 – 2.80	Depositor EDS
% Data completeness (in resolution range)	96.6 (78.05-2.80) 96.6 (78.05-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.38 (at 2.82Å)	Xtrriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, $R_{free}$	0.202 , 0.236 0.199 , 0.232	Depositor DCC
$R_{free}$ test set	3981 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	58.1	Xtrriage
Anisotropy	0.401	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 40.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	0.078 for h,-h-k,-l	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	23339	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	60.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.33% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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 [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.27	0/1188	0.50	0/1617
1	B	0.29	0/1188	0.51	0/1617
1	C	0.26	0/1188	0.50	0/1617
1	D	0.27	0/1188	0.50	0/1617
1	E	0.30	0/1188	0.54	1/1617 (0.1%)
1	F	0.26	0/1188	0.50	0/1617
1	G	0.26	0/1188	0.50	0/1617
1	H	0.28	0/1188	0.56	2/1617 (0.1%)
1	I	0.26	0/1188	0.49	0/1617
1	J	0.26	0/1192	0.49	0/1622
1	K	0.27	0/1188	0.50	0/1617
1	L	0.28	0/1188	0.53	1/1617 (0.1%)
1	M	0.27	0/1192	0.54	0/1622
1	N	0.26	0/1188	0.49	0/1617
1	O	0.28	0/1192	0.51	0/1622
1	P	0.26	0/1188	0.49	0/1617
1	Q	0.26	0/1188	0.52	0/1617
1	R	0.26	0/1188	0.50	0/1617
1	S	0.26	0/1192	0.49	0/1622
1	T	0.27	0/1188	0.51	0/1617
All	All	0.27	0/23776	0.51	4/32360 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	137	ASP	CB-CG-OD1	7.18	124.76	118.30
1	E	137	ASP	CB-CG-OD1	6.01	123.71	118.30
1	L	137	ASP	CB-CG-OD1	5.84	123.56	118.30
1	H	137	ASP	CB-CG-OD2	-5.08	113.73	118.30

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	1164	0	1187	15	0
1	B	1164	0	1187	23	0
1	C	1164	0	1187	12	0
1	D	1164	0	1187	10	0
1	E	1164	0	1187	15	0
1	F	1164	0	1187	15	0
1	G	1164	0	1187	12	0
1	H	1164	0	1187	17	0
1	I	1164	0	1187	13	0
1	J	1168	0	1191	9	0
1	K	1164	0	1187	12	0
1	L	1164	0	1187	17	0
1	M	1168	0	1191	15	0
1	N	1164	0	1187	17	0
1	O	1168	0	1191	20	0
1	P	1164	0	1187	17	0
1	Q	1164	0	1187	22	0
1	R	1164	0	1187	13	0
1	S	1168	0	1191	14	0
1	T	1164	0	1187	15	0
2	A	2	0	0	0	0
2	C	2	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
2	M	2	0	0	0	0
2	N	1	0	0	0	0
2	O	2	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	P	1	0	0	0	0
2	R	1	0	0	0	0
2	S	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	E	1	0	0	0	0
3	H	2	0	0	0	0
3	I	2	0	0	0	0
3	J	2	0	0	0	0
3	K	4	0	0	0	0
3	L	2	0	0	0	0
3	M	1	0	0	0	0
3	O	1	0	0	0	0
3	Q	1	0	0	0	0
3	T	1	0	0	0	0
All	All	23339	0	23756	246	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 5.

All (246) 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:B:55:GLU:OE2	1:B:59:ARG:NH2	2.08	0.85
1:C:152:PRO:HB2	1:C:154:LYS:HE3	1.58	0.84
1:E:119:GLY:HA2	1:E:153:THR:HG23	1.58	0.83
1:Q:55:GLU:HG3	1:Q:56:PRO:HD3	1.61	0.82
1:F:152:PRO:HB2	1:F:154:LYS:HE3	1.61	0.81
1:M:151:LEU:HB3	1:P:124:ILE:HD11	1.61	0.79
1:O:151:LEU:HD21	1:S:151:LEU:HD11	1.64	0.79
1:E:153:THR:HG22	1:H:124:ILE:HG23	1.68	0.75
1:F:55:GLU:HG3	1:F:56:PRO:HD3	1.67	0.75
1:D:151:LEU:HD21	1:G:151:LEU:HD11	1.69	0.73
1:A:69:GLU:HG3	1:O:69:GLU:HB3	1.73	0.71
1:B:151:LEU:HD21	1:J:151:LEU:HD21	1.73	0.70
1:F:14:LEU:HD22	1:F:74:LEU:HB3	1.74	0.69
1:K:151:LEU:HD21	1:R:151:LEU:HD21	1.75	0.69
1:M:151:LEU:HD11	1:P:151:LEU:HD21	1.76	0.68
1:A:68:ASP:HA	1:K:24:ARG:HD3	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:28:THR:HG22	1:I:30:GLU:H	1.60	0.67
1:C:14:LEU:HD22	1:C:74:LEU:HB3	1.77	0.66
1:L:14:LEU:HD22	1:L:74:LEU:HB3	1.76	0.65
1:A:66:GLY:HA3	1:K:25:HIS:CE1	2.32	0.65
1:N:156:PHE:OXT	1:T:39:SER:OG	2.13	0.65
1:K:135:ASP:HB2	1:K:141:ARG:HG3	1.78	0.64
1:S:14:LEU:HD21	1:S:75:LEU:HG	1.79	0.64
1:O:14:LEU:HD21	1:O:75:LEU:HG	1.80	0.64
1:L:47:LEU:HD21	1:L:61:VAL:HG11	1.80	0.64
1:G:14:LEU:HD22	1:G:74:LEU:HB3	1.80	0.63
1:N:151:LEU:HD21	1:T:151:LEU:HD21	1.81	0.62
1:C:59:ARG:HG2	1:C:72:VAL:HG11	1.81	0.61
1:L:135:ASP:HB2	1:L:141:ARG:HG3	1.83	0.61
1:T:135:ASP:HB2	1:T:141:ARG:HG3	1.82	0.61
1:N:83:SER:HG	1:P:106:SER:HG	1.50	0.60
1:K:156:PHE:OXT	1:R:39:SER:OG	2.19	0.59
1:P:5:LEU:HA	1:Q:19:MET:HE1	1.82	0.59
1:L:122:SER:HB2	1:L:151:LEU:HD13	1.85	0.59
1:Q:47:LEU:HD21	1:Q:61:VAL:HG11	1.85	0.59
1:I:14:LEU:HD22	1:I:74:LEU:HB3	1.84	0.59
1:A:19:MET:O	1:A:23:ILE:HG12	2.02	0.58
1:E:59:ARG:NH2	1:E:69:GLU:OE2	2.34	0.58
1:O:31:ALA:O	1:O:34:GLN:HG2	2.03	0.58
1:S:73:LEU:HD21	1:T:23:ILE:HG21	1.85	0.58
1:O:152:PRO:HB2	1:O:154:LYS:HE3	1.85	0.58
1:A:14:LEU:HD21	1:A:75:LEU:HG	1.85	0.58
1:H:31:ALA:O	1:H:34:GLN:HG2	2.04	0.58
1:H:68:ASP:OD1	1:H:69:GLU:N	2.32	0.58
1:E:156:PHE:OXT	1:H:39:SER:OG	2.19	0.57
1:L:69:GLU:HA	1:L:72:VAL:HG23	1.85	0.57
1:P:14:LEU:HD22	1:P:74:LEU:HB3	1.85	0.57
1:B:31:ALA:O	1:B:34:GLN:HG2	2.05	0.57
1:O:55:GLU:HG2	1:O:59:ARG:NH1	2.19	0.57
1:O:122:SER:HB2	1:O:151:LEU:HD13	1.87	0.57
1:I:30:GLU:OE2	1:I:34:GLN:NE2	2.37	0.56
1:C:19:MET:O	1:C:23:ILE:HG12	2.05	0.56
1:R:47:LEU:HD21	1:R:61:VAL:HG11	1.88	0.56
1:F:127:LYS:HB2	1:F:147:ASP:HB3	1.88	0.56
1:J:55:GLU:OE2	1:J:59:ARG:HG3	2.06	0.56
1:P:60:THR:HG22	1:P:64:LYS:HE3	1.87	0.56
1:R:19:MET:O	1:R:23:ILE:HG12	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:30:GLU:O	1:N:34:GLN:HG3	2.06	0.55
1:L:59:ARG:HG2	1:L:72:VAL:HG11	1.88	0.55
1:B:72:VAL:HG23	1:B:76:GLN:HE21	1.70	0.54
1:L:99:GLU:O	1:L:103:VAL:HG23	2.07	0.54
1:L:151:LEU:HD21	1:Q:151:LEU:HD21	1.88	0.54
1:B:14:LEU:HD21	1:B:75:LEU:HG	1.90	0.53
1:G:99:GLU:O	1:G:103:VAL:HG23	2.07	0.53
1:H:135:ASP:OD1	1:H:141:ARG:HD2	2.08	0.53
1:M:14:LEU:HD22	1:M:74:LEU:HB3	1.90	0.53
1:S:99:GLU:O	1:S:103:VAL:HG23	2.08	0.53
1:B:69:GLU:N	1:B:69:GLU:OE1	2.41	0.53
1:I:122:SER:HB2	1:I:151:LEU:HD13	1.90	0.53
1:P:31:ALA:O	1:P:34:GLN:HG2	2.09	0.53
1:E:151:LEU:HD21	1:H:151:LEU:HD11	1.89	0.53
1:G:135:ASP:HB2	1:G:141:ARG:HG3	1.91	0.53
1:Q:127:LYS:HB2	1:Q:147:ASP:HB3	1.91	0.52
1:D:47:LEU:HD21	1:D:61:VAL:HG11	1.92	0.52
1:N:127:LYS:HB2	1:N:147:ASP:HB3	1.90	0.52
1:D:14:LEU:HD22	1:D:74:LEU:HB3	1.92	0.52
1:C:137:ASP:OD1	1:C:137:ASP:N	2.32	0.52
1:F:55:GLU:HG3	1:F:56:PRO:CD	2.38	0.52
1:A:73:LEU:HD21	1:E:23:ILE:HG21	1.91	0.52
1:K:122:SER:HB2	1:K:151:LEU:HD13	1.92	0.52
1:H:130:ILE:HD12	1:H:144:ILE:HG13	1.92	0.51
1:Q:68:ASP:OD1	1:Q:69:GLU:N	2.43	0.51
1:D:68:ASP:OD1	1:D:69:GLU:N	2.37	0.51
1:I:6:HIS:ND1	1:J:19:MET:HG3	2.26	0.51
1:I:99:GLU:O	1:I:103:VAL:HG23	2.10	0.51
1:A:122:SER:HB2	1:A:151:LEU:HD13	1.93	0.51
1:B:99:GLU:O	1:B:103:VAL:HG23	2.11	0.51
1:C:99:GLU:O	1:C:103:VAL:HG23	2.11	0.51
1:H:135:ASP:HB2	1:H:141:ARG:HG3	1.92	0.51
1:R:99:GLU:O	1:R:103:VAL:HG23	2.11	0.51
1:B:69:GLU:HA	1:B:72:VAL:HG13	1.93	0.50
1:F:30:GLU:O	1:F:34:GLN:HG3	2.12	0.50
1:M:135:ASP:OD2	1:M:137:ASP:HB3	2.10	0.50
1:Q:55:GLU:HG3	1:Q:56:PRO:CD	2.37	0.50
1:I:14:LEU:HD21	1:I:75:LEU:HG	1.94	0.50
1:B:122:SER:HB2	1:B:151:LEU:HD13	1.94	0.50
1:A:151:LEU:HD21	1:I:151:LEU:HD21	1.94	0.49
1:I:19:MET:O	1:I:23:ILE:HG12	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:O:55:GLU:HG2	1:O:59:ARG:HH12	1.77	0.49
1:S:59:ARG:HG2	1:S:72:VAL:HG11	1.94	0.49
1:L:31:ALA:O	1:L:34:GLN:HG2	2.12	0.49
1:E:25:HIS:CE1	1:O:66:GLY:HA3	2.47	0.49
1:M:126:PHE:HE1	1:M:146:LEU:HB3	1.77	0.49
1:R:122:SER:HB2	1:R:151:LEU:HD13	1.93	0.49
1:M:156:PHE:OXT	1:P:39:SER:OG	2.30	0.49
1:Q:14:LEU:HD22	1:Q:74:LEU:HB3	1.93	0.49
1:L:14:LEU:HD21	1:L:75:LEU:HG	1.94	0.48
1:M:68:ASP:OD1	1:M:69:GLU:N	2.37	0.48
1:M:99:GLU:O	1:M:103:VAL:HG23	2.13	0.48
1:B:152:PRO:HB2	1:B:154:LYS:HD3	1.95	0.48
1:Q:99:GLU:O	1:Q:103:VAL:HG23	2.13	0.48
1:F:47:LEU:HD21	1:F:61:VAL:HG11	1.95	0.48
1:N:14:LEU:HD22	1:N:74:LEU:HB3	1.95	0.48
1:E:39:SER:OG	1:H:156:PHE:OXT	2.25	0.48
1:G:4:SER:HB2	1:G:78:ILE:HB	1.95	0.48
1:N:99:GLU:O	1:N:103:VAL:HG23	2.12	0.48
1:S:119:GLY:HA2	1:S:153:THR:OG1	2.13	0.48
1:T:99:GLU:O	1:T:103:VAL:HG23	2.13	0.48
1:N:14:LEU:HD21	1:N:75:LEU:HG	1.95	0.48
1:L:135:ASP:OD2	1:L:141:ARG:NH1	2.37	0.47
1:N:120:ILE:HG12	1:T:124:ILE:HG13	1.96	0.47
1:A:19:MET:HG3	1:B:6:HIS:ND1	2.29	0.47
1:C:122:SER:HB2	1:C:151:LEU:HD13	1.95	0.47
1:H:14:LEU:HD21	1:H:75:LEU:HG	1.96	0.47
1:H:99:GLU:O	1:H:103:VAL:HG23	2.14	0.47
1:M:135:ASP:OD1	1:M:141:ARG:HD2	2.13	0.47
1:Q:6:HIS:ND1	1:R:19:MET:HG3	2.29	0.47
1:J:122:SER:HB2	1:J:151:LEU:HD13	1.97	0.47
1:P:57:ALA:O	1:P:61:VAL:HG22	2.15	0.47
1:N:48:LEU:HD13	1:P:106:SER:HB3	1.96	0.47
1:T:14:LEU:HD22	1:T:74:LEU:HB3	1.96	0.47
1:N:44:THR:HB	1:P:107:THR:HG23	1.97	0.46
1:R:60:THR:HG22	1:R:64:LYS:HE3	1.97	0.46
1:M:135:ASP:HB2	1:M:141:ARG:HG3	1.97	0.46
1:T:122:SER:HB2	1:T:151:LEU:HD13	1.97	0.46
1:A:103:VAL:HG23	1:A:104:TYR:CD2	2.51	0.46
1:Q:14:LEU:HD21	1:Q:75:LEU:HG	1.98	0.46
1:T:119:GLY:HA2	1:T:153:THR:OG1	2.15	0.46
1:B:56:PRO:O	1:B:60:THR:HG23	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:39:SER:OG	1:F:156:PHE:OXT	2.27	0.46
1:R:14:LEU:HD21	1:R:75:LEU:HG	1.98	0.46
1:C:156:PHE:OXT	1:F:39:SER:OG	2.30	0.46
1:D:14:LEU:HD21	1:D:75:LEU:HG	1.98	0.46
1:P:14:LEU:HD21	1:P:75:LEU:HG	1.97	0.46
1:R:14:LEU:HD22	1:R:74:LEU:HB3	1.98	0.46
1:K:99:GLU:O	1:K:103:VAL:HG23	2.15	0.46
1:E:99:GLU:O	1:E:103:VAL:HG23	2.16	0.46
1:F:119:GLY:HA2	1:F:153:THR:OG1	2.16	0.45
1:H:126:PHE:HE1	1:H:146:LEU:HB3	1.81	0.45
1:B:95:TYR:CE1	1:B:98:TYR:HD2	2.35	0.45
1:N:135:ASP:HB2	1:N:141:ARG:HG3	1.99	0.45
1:T:68:ASP:O	1:T:72:VAL:HG23	2.16	0.45
1:E:155:PRO:HB3	1:H:42:PHE:CE2	2.51	0.45
1:L:19:MET:O	1:L:23:ILE:HG12	2.16	0.45
1:Q:28:THR:O	1:Q:32:LEU:HD13	2.17	0.45
1:S:11:ARG:NH2	1:S:46:ALA:O	2.47	0.45
1:C:69:GLU:HA	1:C:72:VAL:HG23	1.97	0.45
1:G:119:GLY:HA2	1:G:153:THR:OG1	2.17	0.45
1:N:20:ALA:HB1	1:N:24:ARG:NH1	2.32	0.45
1:T:15:THR:HG23	1:T:48:LEU:HA	1.99	0.45
1:H:119:GLY:HA2	1:H:153:THR:OG1	2.17	0.45
1:Q:122:SER:HB2	1:Q:151:LEU:HD13	1.97	0.45
1:D:119:GLY:HA2	1:D:153:THR:OG1	2.17	0.45
1:Q:133:VAL:HG13	1:Q:141:ARG:HB2	1.99	0.45
1:B:69:GLU:H	1:B:69:GLU:CD	2.21	0.44
1:C:8:ALA:O	1:C:12:GLU:HG2	2.16	0.44
1:M:119:GLY:HA2	1:M:153:THR:OG1	2.18	0.44
1:A:14:LEU:HD22	1:A:74:LEU:HB3	2.00	0.44
1:R:135:ASP:HB2	1:R:141:ARG:HG3	1.98	0.44
1:O:149:LYS:HD3	1:T:135:ASP:OD2	2.18	0.44
1:Q:19:MET:O	1:Q:23:ILE:HG13	2.18	0.44
1:Q:119:GLY:HA2	1:Q:153:THR:OG1	2.18	0.44
1:M:127:LYS:HB2	1:M:147:ASP:HB3	2.00	0.44
1:O:135:ASP:HB2	1:O:141:ARG:HG3	2.00	0.44
1:G:56:PRO:O	1:G:60:THR:HG23	2.18	0.44
1:O:22:LYS:HE3	1:O:29:PHE:CE2	2.53	0.44
1:B:19:MET:O	1:B:23:ILE:HG13	2.18	0.43
1:F:73:LEU:HD21	1:G:23:ILE:HG22	2.00	0.43
1:K:119:GLY:HA2	1:K:153:THR:OG1	2.18	0.43
1:N:6:HIS:ND1	1:O:19:MET:HG3	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:69:GLU:H	1:S:69:GLU:CD	2.21	0.43
1:A:119:GLY:HA2	1:A:153:THR:OG1	2.19	0.43
1:L:126:PHE:HE1	1:L:146:LEU:HB3	1.82	0.43
1:M:106:SER:HB3	1:Q:48:LEU:HD13	1.99	0.43
1:Q:32:LEU:HD23	1:Q:47:LEU:HD13	2.00	0.43
1:F:14:LEU:HD21	1:F:75:LEU:HG	2.00	0.43
1:O:119:GLY:HA2	1:O:153:THR:OG1	2.18	0.43
1:O:126:PHE:HE1	1:O:146:LEU:HB3	1.82	0.43
1:M:59:ARG:NH2	1:M:69:GLU:OE2	2.45	0.43
1:P:135:ASP:HB2	1:P:141:ARG:HG3	2.01	0.43
1:A:120:ILE:HG12	1:I:124:ILE:HG13	2.00	0.43
1:B:69:GLU:HG3	1:K:69:GLU:HG3	2.01	0.42
1:E:149:LYS:HD3	1:I:135:ASP:OD2	2.19	0.42
1:P:2:THR:HG23	1:P:78:ILE:HD11	2.01	0.42
1:H:6:HIS:ND1	1:I:19:MET:HG3	2.34	0.42
1:P:34:GLN:O	1:P:64:LYS:NZ	2.52	0.42
1:S:11:ARG:HD2	1:S:11:ARG:HA	1.89	0.42
1:J:18:ILE:HD13	1:J:47:LEU:HD22	2.02	0.42
1:P:103:VAL:HG23	1:P:104:TYR:CD2	2.54	0.42
1:N:126:PHE:HE1	1:N:146:LEU:HB3	1.84	0.42
1:D:99:GLU:O	1:D:103:VAL:HG23	2.19	0.42
1:P:122:SER:HB2	1:P:151:LEU:HD13	2.01	0.42
1:F:74:LEU:HD12	1:F:74:LEU:HA	1.93	0.42
1:K:96:ARG:NH2	2:O:201:CL:CL	2.82	0.42
1:G:74:LEU:HD12	1:G:74:LEU:HA	1.88	0.42
1:I:34:GLN:O	1:I:64:LYS:NZ	2.44	0.42
1:L:42:PHE:CE2	1:Q:155:PRO:HB3	2.55	0.42
1:S:89:PRO:HG3	1:S:94:ILE:HG22	2.02	0.42
1:O:2:THR:OG1	1:O:3:HIS:N	2.53	0.42
1:O:55:GLU:HG3	1:O:76:GLN:NE2	2.35	0.41
1:J:119:GLY:HA2	1:J:153:THR:OG1	2.19	0.41
1:S:47:LEU:HD21	1:S:61:VAL:HG11	2.01	0.41
1:A:23:ILE:CD1	1:B:3:HIS:HB3	2.50	0.41
1:D:146:LEU:HB2	1:H:144:ILE:HB	2.02	0.41
1:T:14:LEU:HD21	1:T:75:LEU:HG	2.03	0.41
1:R:124:ILE:H	1:R:124:ILE:HG13	1.70	0.41
1:C:97:PHE:CD2	1:G:108:LEU:HD13	2.55	0.41
1:E:42:PHE:CE2	1:H:155:PRO:HB3	2.55	0.41
1:F:73:LEU:HD21	1:G:23:ILE:CG2	2.51	0.41
1:J:129:ASP:OD1	1:J:130:ILE:N	2.54	0.41
1:B:74:LEU:HD12	1:B:74:LEU:HA	1.90	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:122:SER:HB2	1:D:151:LEU:HD13	2.03	0.41
1:D:39:SER:OG	1:G:156:PHE:OXT	2.27	0.41
1:M:106:SER:OG	1:Q:83:SER:HB2	2.21	0.41
1:N:122:SER:HB2	1:N:151:LEU:HD13	2.03	0.41
1:B:129:ASP:OD1	1:B:130:ILE:N	2.54	0.41
1:F:69:GLU:HG2	1:F:70:GLU:N	2.35	0.41
1:K:151:LEU:CD2	1:R:151:LEU:HD21	2.49	0.41
1:L:74:LEU:HD12	1:L:74:LEU:HA	1.86	0.41
1:Q:43:VAL:O	1:Q:47:LEU:HG	2.20	0.41
1:B:55:GLU:HG2	1:B:76:GLN:NE2	2.36	0.41
1:E:154:LYS:HA	1:E:154:LYS:HD2	1.82	0.41
1:A:25:HIS:CE1	1:K:66:GLY:HA3	2.56	0.40
1:E:122:SER:HB2	1:E:151:LEU:HD13	2.02	0.40
1:O:155:PRO:HB3	1:S:42:PHE:CE2	2.56	0.40
1:L:156:PHE:OXT	1:Q:39:SER:OG	2.39	0.40
1:O:42:PHE:CE2	1:S:155:PRO:HB3	2.57	0.40
1:O:120:ILE:HG12	1:S:124:ILE:HG13	2.03	0.40
1:B:39:SER:OG	1:J:156:PHE:OXT	2.28	0.40
1:E:135:ASP:HB2	1:E:141:ARG:HG3	2.03	0.40
1:N:20:ALA:HB1	1:N:24:ARG:HH12	1.86	0.40
1:B:89:PRO:HG3	1:B:94:ILE:HG22	2.04	0.40
1:L:126:PHE:HA	1:L:147:ASP:O	2.22	0.40
1:T:22:LYS:HE3	1:T:29:PHE:CE2	2.55	0.40
1:B:156:PHE:OXT	1:J:39:SER:OG	2.29	0.40
1:T:22:LYS:HE3	1:T:29:PHE:CD2	2.56	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	153/155 (99%)	151 (99%)	2 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	153/155 (99%)	150 (98%)	3 (2%)	0	100	100
1	C	153/155 (99%)	150 (98%)	3 (2%)	0	100	100
1	D	153/155 (99%)	151 (99%)	2 (1%)	0	100	100
1	E	153/155 (99%)	150 (98%)	3 (2%)	0	100	100
1	F	153/155 (99%)	150 (98%)	3 (2%)	0	100	100
1	G	153/155 (99%)	149 (97%)	4 (3%)	0	100	100
1	H	153/155 (99%)	151 (99%)	2 (1%)	0	100	100
1	I	153/155 (99%)	151 (99%)	2 (1%)	0	100	100
1	J	153/155 (99%)	149 (97%)	4 (3%)	0	100	100
1	K	153/155 (99%)	149 (97%)	4 (3%)	0	100	100
1	L	153/155 (99%)	149 (97%)	4 (3%)	0	100	100
1	M	153/155 (99%)	150 (98%)	3 (2%)	0	100	100
1	N	153/155 (99%)	151 (99%)	2 (1%)	0	100	100
1	O	153/155 (99%)	150 (98%)	3 (2%)	0	100	100
1	P	153/155 (99%)	150 (98%)	3 (2%)	0	100	100
1	Q	153/155 (99%)	149 (97%)	4 (3%)	0	100	100
1	R	153/155 (99%)	150 (98%)	3 (2%)	0	100	100
1	S	153/155 (99%)	149 (97%)	4 (3%)	0	100	100
1	T	153/155 (99%)	150 (98%)	3 (2%)	0	100	100
All	All	3060/3100 (99%)	2999 (98%)	61 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	120/122 (98%)	119 (99%)	1 (1%)	81	94
1	B	120/122 (98%)	119 (99%)	1 (1%)	81	94

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	120/122 (98%)	118 (98%)	2 (2%)	60	87
1	D	120/122 (98%)	120 (100%)	0	100	100
1	E	120/122 (98%)	117 (98%)	3 (2%)	47	80
1	F	120/122 (98%)	119 (99%)	1 (1%)	81	94
1	G	120/122 (98%)	119 (99%)	1 (1%)	81	94
1	H	120/122 (98%)	120 (100%)	0	100	100
1	I	120/122 (98%)	120 (100%)	0	100	100
1	J	121/122 (99%)	121 (100%)	0	100	100
1	K	120/122 (98%)	120 (100%)	0	100	100
1	L	120/122 (98%)	117 (98%)	3 (2%)	47	80
1	M	121/122 (99%)	119 (98%)	2 (2%)	60	87
1	N	120/122 (98%)	117 (98%)	3 (2%)	47	80
1	O	121/122 (99%)	121 (100%)	0	100	100
1	P	120/122 (98%)	119 (99%)	1 (1%)	81	94
1	Q	120/122 (98%)	116 (97%)	4 (3%)	38	72
1	R	120/122 (98%)	120 (100%)	0	100	100
1	S	121/122 (99%)	119 (98%)	2 (2%)	60	87
1	T	120/122 (98%)	120 (100%)	0	100	100
All	All	2404/2440 (98%)	2380 (99%)	24 (1%)	76	93

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	VAL
1	B	72	VAL
1	C	72	VAL
1	C	137	ASP
1	E	4	SER
1	E	140	GLU
1	E	154	LYS
1	F	55	GLU
1	G	154	LYS
1	L	16	ASP
1	L	72	VAL
1	L	73	LEU
1	M	38	LEU

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Mol	Chain	Res	Type
1	M	137	ASP
1	N	2	THR
1	N	16	ASP
1	N	80	LEU
1	P	61	VAL
1	Q	19	MET
1	Q	25	HIS
1	Q	27	LEU
1	Q	55	GLU
1	S	72	VAL
1	S	74	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	76	GLN
1	E	25	HIS
1	T	76	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 20 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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	155/155 (100%)	-0.17	0 100 100	42, 54, 75, 129	0
1	B	155/155 (100%)	-0.06	0 100 100	44, 62, 96, 118	0
1	C	155/155 (100%)	-0.08	0 100 100	42, 56, 85, 128	0
1	D	155/155 (100%)	-0.13	0 100 100	43, 59, 81, 111	0
1	E	155/155 (100%)	-0.15	0 100 100	41, 53, 74, 105	0
1	F	155/155 (100%)	-0.20	0 100 100	41, 56, 74, 89	0
1	G	155/155 (100%)	-0.13	0 100 100	41, 55, 81, 119	0
1	H	155/155 (100%)	-0.15	0 100 100	40, 56, 79, 138	0
1	I	155/155 (100%)	-0.18	0 100 100	41, 54, 77, 140	0
1	J	155/155 (100%)	-0.21	0 100 100	40, 54, 75, 98	0
1	K	155/155 (100%)	-0.17	0 100 100	42, 55, 77, 112	0
1	L	155/155 (100%)	-0.08	1 (0%) 89 86	44, 61, 85, 111	0
1	M	155/155 (100%)	0.01	0 100 100	51, 69, 91, 132	0
1	N	155/155 (100%)	-0.06	0 100 100	46, 61, 87, 124	0
1	O	155/155 (100%)	-0.13	0 100 100	43, 56, 85, 138	0
1	P	155/155 (100%)	-0.09	0 100 100	47, 62, 80, 93	0
1	Q	155/155 (100%)	0.05	0 100 100	52, 68, 92, 127	0
1	R	155/155 (100%)	-0.12	0 100 100	46, 60, 85, 121	0
1	S	155/155 (100%)	-0.18	1 (0%) 89 86	44, 55, 77, 139	0
1	T	155/155 (100%)	-0.14	0 100 100	41, 55, 78, 117	0
All	All	3100/3100 (100%)	-0.12	2 (0%) 95 95	40, 58, 85, 140	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	S	137	ASP	2.1
1	L	2	THR	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 monosaccharides 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, 95<sup>th</sup> 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.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	CL	O	201	1/1	0.96	0.08	52,52,52,52	0
2	CL	P	201	1/1	0.96	0.15	56,56,56,56	0
2	CL	D	201	1/1	0.97	0.08	49,49,49,49	0
2	CL	M	202	1/1	0.97	0.07	61,61,61,61	0
2	CL	R	201	1/1	0.97	0.08	55,55,55,55	0
2	CL	J	201	1/1	0.98	0.12	52,52,52,52	0
2	CL	L	201	1/1	0.98	0.15	65,65,65,65	0
2	CL	M	201	1/1	0.98	0.18	63,63,63,63	0
2	CL	C	201	1/1	0.98	0.07	54,54,54,54	0
2	CL	N	201	1/1	0.98	0.15	47,47,47,47	0
2	CL	A	201	1/1	0.98	0.12	55,55,55,55	0
2	CL	O	202	1/1	0.98	0.08	52,52,52,52	0
2	CL	F	201	1/1	0.98	0.08	54,54,54,54	0
2	CL	G	201	1/1	0.98	0.13	63,63,63,63	0
2	CL	S	201	1/1	0.98	0.09	48,48,48,48	0
2	CL	K	201	1/1	0.99	0.07	49,49,49,49	0
2	CL	C	202	1/1	0.99	0.11	55,55,55,55	0
2	CL	A	202	1/1	0.99	0.12	50,50,50,50	0
2	CL	I	201	1/1	0.99	0.12	53,53,53,53	0
2	CL	E	201	1/1	0.99	0.09	57,57,57,57	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.