



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:25 AM UTC

PDB ID : 5CBF / pdb\_00005cbf  
Title : Structural and Functional Characterization of a Calcium-activated Cation channel from Tsukamurella Paurometabola  
Authors : Dhakshnamoorthy, B.; Rohaim, A.; Rui, H.; Blachowicz, L.; Roux, B.  
Deposited on : 2015-06-30  
Resolution : 3.61 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 : 4-5-2 with Phenix2.0  
Xtrriage (Phenix) : 2.0  
EDS : 3.0  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

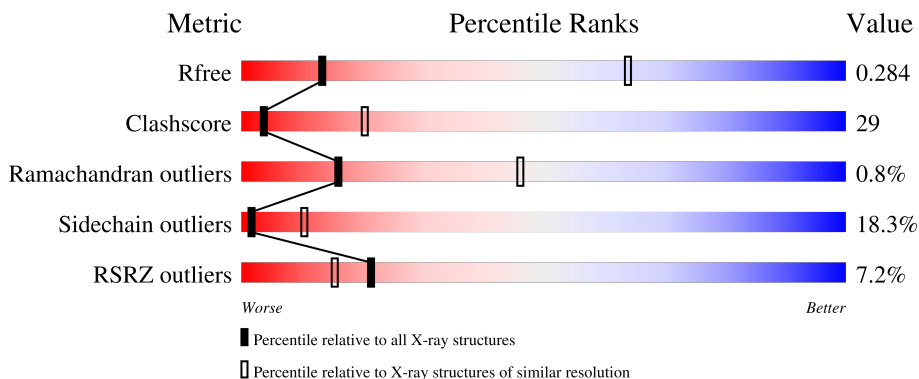
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.61 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	1062 (3.72-3.52)
Clashscore	190562	1092 (3.72-3.52)
Ramachandran outliers	187476	1057 (3.72-3.52)
Sidechain outliers	187428	1055 (3.72-3.52)
RSRZ outliers	180081	1060 (3.72-3.52)

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	129	 2% 35% 33% 11% 21%
1	B	129	 2% 33% 36% 10% 21%
1	C	129	 9% 41% 26% 12% 21%
1	D	129	 5% 36% 31% 11% 21%
1	E	129	 4% 37% 28% 13% 21%

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Mol	Chain	Length	Quality of chain
1	F	129	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into five segments: 11% red, 33% green, 35% yellow, 10% orange, and 21% grey. The segments are stacked from left to right in the order: red, green, yellow, orange, grey.</p>

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 4623 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 Ion transport 2 domain protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	102	770	510	123	135	2	0	0	0
1	B	102	770	510	123	135	2	0	0	0
1	C	102	770	510	123	135	2	0	0	0
1	D	102	770	510	123	135	2	0	0	0
1	E	102	770	510	123	135	2	0	0	0
1	F	102	770	510	123	135	2	0	0	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	124	HIS	-	expression tag	UNP D5UM26
A	125	HIS	-	expression tag	UNP D5UM26
A	126	HIS	-	expression tag	UNP D5UM26
A	127	HIS	-	expression tag	UNP D5UM26
A	128	HIS	-	expression tag	UNP D5UM26
A	129	HIS	-	expression tag	UNP D5UM26
B	124	HIS	-	expression tag	UNP D5UM26
B	125	HIS	-	expression tag	UNP D5UM26
B	126	HIS	-	expression tag	UNP D5UM26
B	127	HIS	-	expression tag	UNP D5UM26
B	128	HIS	-	expression tag	UNP D5UM26
B	129	HIS	-	expression tag	UNP D5UM26
C	124	HIS	-	expression tag	UNP D5UM26
C	125	HIS	-	expression tag	UNP D5UM26
C	126	HIS	-	expression tag	UNP D5UM26
C	127	HIS	-	expression tag	UNP D5UM26
C	128	HIS	-	expression tag	UNP D5UM26

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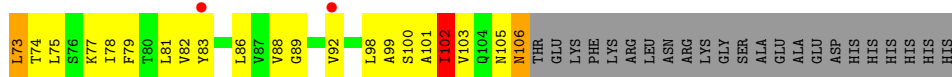
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Chain	Residue	Modelled	Actual	Comment	Reference
C	129	HIS	-	expression tag	UNP D5UM26
D	124	HIS	-	expression tag	UNP D5UM26
D	125	HIS	-	expression tag	UNP D5UM26
D	126	HIS	-	expression tag	UNP D5UM26
D	127	HIS	-	expression tag	UNP D5UM26
D	128	HIS	-	expression tag	UNP D5UM26
D	129	HIS	-	expression tag	UNP D5UM26
E	124	HIS	-	expression tag	UNP D5UM26
E	125	HIS	-	expression tag	UNP D5UM26
E	126	HIS	-	expression tag	UNP D5UM26
E	127	HIS	-	expression tag	UNP D5UM26
E	128	HIS	-	expression tag	UNP D5UM26
E	129	HIS	-	expression tag	UNP D5UM26
F	124	HIS	-	expression tag	UNP D5UM26
F	125	HIS	-	expression tag	UNP D5UM26
F	126	HIS	-	expression tag	UNP D5UM26
F	127	HIS	-	expression tag	UNP D5UM26
F	128	HIS	-	expression tag	UNP D5UM26
F	129	HIS	-	expression tag	UNP D5UM26

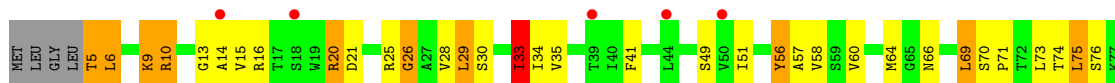
- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Ca 1 1	0	0
2	B	1	Total Ca 1 1	0	0
2	E	1	Total Ca 1 1	0	0

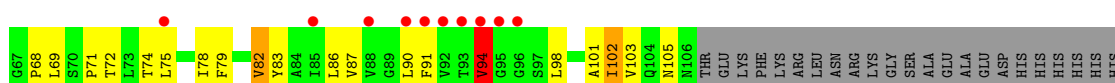
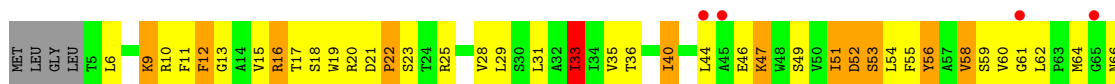




• Molecule 1: Ion transport 2 domain protein



• Molecule 1: Ion transport 2 domain protein



## 4 Data and refinement statistics i

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	116.05Å 116.05Å 132.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 3.61 50.00 – 3.61	Depositor EDS
% Data completeness (in resolution range)	99.1 (50.00-3.61) 99.1 (50.00-3.61)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.61 (at 3.57Å)	Xtrriage
Refinement program	REFMAC 5.8.0103	Depositor
R, $R_{free}$	0.239 , 0.279 0.241 , 0.284	Depositor DCC
$R_{free}$ test set	489 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	122.1	Xtrriage
Anisotropy	0.022	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.25 , 40.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.25$	Xtrriage
Estimated twinning fraction	0.206 for -h,k,-l	Xtrriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	4623	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	114.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.99% 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: CA

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	1.05	0/787	1.64	14/1075 (1.3%)
1	B	1.05	1/787 (0.1%)	1.62	13/1075 (1.2%)
1	C	1.04	1/787 (0.1%)	1.64	10/1075 (0.9%)
1	D	1.08	1/787 (0.1%)	1.62	9/1075 (0.8%)
1	E	1.06	2/787 (0.3%)	1.66	15/1075 (1.4%)
1	F	1.12	1/787 (0.1%)	1.60	9/1075 (0.8%)
All	All	1.07	6/4722 (0.1%)	1.63	70/6450 (1.1%)

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	A	0	1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	51	ILE	C-O	7.49	1.32	1.24
1	F	51	ILE	C-O	7.33	1.32	1.24
1	B	21	ASP	CA-C	5.32	1.58	1.53
1	E	89	GLY	C-O	5.21	1.30	1.23
1	D	32	ALA	C-O	5.12	1.29	1.24
1	E	26	GLY	C-O	-5.07	1.18	1.24

All (70) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	20	ARG	N-CA-C	-8.32	93.34	107.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	92	VAL	O-C-N	8.02	129.77	121.91
1	C	40	ILE	N-CA-CB	7.70	119.55	110.55
1	D	70	SER	CA-C-N	6.91	128.47	119.84
1	D	70	SER	C-N-CA	6.91	128.47	119.84
1	D	13	GLY	N-CA-C	6.80	129.30	113.18
1	F	23	SER	N-CA-C	6.73	118.30	110.97
1	D	44	LEU	N-CA-C	6.68	118.44	111.03
1	A	56	TYR	N-CA-CB	6.63	119.62	110.01
1	A	102	ILE	N-CA-CB	6.58	118.25	110.55
1	D	102	ILE	N-CA-CB	6.42	118.79	110.57
1	A	92	VAL	N-CA-C	6.36	116.84	110.23
1	F	33	ILE	O-C-N	6.29	128.07	121.91
1	B	21	ASP	CA-C-N	6.25	127.66	119.84
1	B	21	ASP	C-N-CA	6.25	127.66	119.84
1	E	26	GLY	N-CA-C	-6.20	107.85	114.67
1	D	21	ASP	N-CA-C	6.17	119.27	109.64
1	F	102	ILE	N-CA-CB	5.99	117.56	110.55
1	F	83	TYR	CB-CA-C	5.94	120.66	110.79
1	E	58	VAL	CB-CA-C	5.82	120.00	112.14
1	A	70	SER	CA-C-N	5.79	125.81	119.90
1	A	70	SER	C-N-CA	5.79	125.81	119.90
1	A	68	PRO	N-CA-C	-5.78	106.63	113.86
1	F	56	TYR	N-CA-CB	5.71	118.52	110.12
1	E	21	ASP	N-CA-C	5.68	119.55	107.91
1	D	83	TYR	CB-CA-C	5.66	120.47	110.85
1	E	56	TYR	N-CA-CB	5.64	118.41	110.12
1	F	12	PHE	N-CA-C	-5.61	98.11	107.99
1	E	33	ILE	N-CA-CB	5.60	117.10	110.55
1	E	58	VAL	CA-C-N	5.59	128.33	120.28
1	E	58	VAL	C-N-CA	5.59	128.33	120.28
1	B	92	VAL	O-C-N	5.59	127.53	121.94
1	A	21	ASP	CA-C-N	5.58	126.82	119.84
1	A	21	ASP	C-N-CA	5.58	126.82	119.84
1	B	98	LEU	O-C-N	5.57	128.03	122.12
1	A	31	LEU	O-C-N	5.56	127.80	122.07
1	C	10	ARG	N-CA-CB	5.54	118.01	109.98
1	B	67	GLY	CA-C-N	5.54	124.88	118.85
1	B	67	GLY	C-N-CA	5.54	124.88	118.85
1	E	33	ILE	O-C-N	5.53	127.33	121.91
1	E	57	ALA	O-C-N	5.48	127.72	122.07
1	A	101	ALA	CA-C-O	-5.48	114.74	120.55
1	C	21	ASP	N-CA-C	5.45	120.71	109.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	75	LEU	O-C-N	5.44	127.67	122.07
1	D	56	TYR	N-CA-CB	5.40	118.06	110.12
1	B	57	ALA	O-C-N	5.39	127.62	122.07
1	A	85	ILE	CA-C-O	-5.35	115.39	120.95
1	C	100	SER	O-C-N	5.30	127.53	122.07
1	B	56	TYR	N-CA-CB	5.29	117.89	110.12
1	C	93	THR	CA-C-O	-5.29	114.95	120.55
1	C	19	TRP	CA-C-N	5.28	129.84	121.72
1	C	19	TRP	C-N-CA	5.28	129.84	121.72
1	B	26	GLY	N-CA-C	-5.23	107.83	114.16
1	B	35	VAL	O-C-N	5.23	126.94	121.87
1	B	33	ILE	CB-CA-C	5.22	119.19	112.14
1	A	86	LEU	N-CA-C	5.15	116.70	111.14
1	D	79	PHE	O-C-N	5.14	127.57	122.12
1	B	80	THR	O-C-N	5.14	127.36	122.07
1	C	42	TYR	N-CA-CB	5.14	117.67	110.12
1	F	94	VAL	CA-C-N	5.13	125.63	119.94
1	F	94	VAL	C-N-CA	5.13	125.63	119.94
1	C	83	TYR	O-C-N	5.12	127.35	122.07
1	B	86	LEU	N-CA-C	5.12	116.55	111.07
1	E	21	ASP	CA-C-N	5.07	126.18	119.84
1	E	21	ASP	C-N-CA	5.07	126.18	119.84
1	F	68	PRO	N-CA-C	-5.06	109.42	114.68
1	E	95	GLY	CA-C-N	5.03	125.53	119.94
1	E	95	GLY	C-N-CA	5.03	125.53	119.94
1	A	6	LEU	CA-C-N	5.01	129.16	121.19
1	A	6	LEU	C-N-CA	5.01	129.16	121.19

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	101	ALA	Mainchain

## 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	770	0	808	53	0
1	B	770	0	808	51	0
1	C	770	0	808	47	0
1	D	770	0	808	63	0
1	E	770	0	808	48	0
1	F	770	0	808	53	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	E	1	0	0	0	0
All	All	4623	0	4848	275	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 29.

All (275) 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:A:10:ARG:NH2	1:A:15:VAL:HG21	1.25	1.45
1:C:40:ILE:O	1:C:44:LEU:HD12	1.40	1.19
1:A:10:ARG:NH2	1:A:15:VAL:CG2	2.07	1.18
1:F:90:LEU:O	1:F:94:VAL:HG23	1.44	1.15
1:A:20:ARG:HB3	1:A:25:ARG:NH2	1.63	1.10
1:F:40:ILE:O	1:F:44:LEU:HD12	1.55	1.06
1:F:22:PRO:HD2	1:F:25:ARG:HG3	1.35	1.05
1:C:40:ILE:O	1:C:44:LEU:CD1	2.07	1.02
1:F:40:ILE:O	1:F:44:LEU:CD1	2.08	1.02
1:E:90:LEU:O	1:E:94:VAL:HG23	1.58	1.01
1:F:20:ARG:HH22	1:F:105:ASN:ND2	1.61	0.98
1:F:19:TRP:NE1	1:F:21:ASP:OD2	1.97	0.98
1:A:20:ARG:HB3	1:A:25:ARG:HH22	1.28	0.96
1:F:17:THR:HG22	1:F:20:ARG:HH21	1.30	0.96
1:F:22:PRO:HD2	1:F:25:ARG:CG	1.97	0.93
1:C:40:ILE:HG22	1:C:44:LEU:HD11	1.51	0.92
1:D:54:LEU:HD23	1:E:81:LEU:HD13	1.56	0.88
1:D:12:PHE:O	1:D:16:ARG:HD2	1.73	0.88
1:A:10:ARG:HH22	1:A:15:VAL:HG21	1.06	0.87
1:C:46:GLU:C	1:C:47:LYS:HG2	1.99	0.86
1:D:17:THR:HA	1:D:20:ARG:NH1	1.90	0.86
1:D:103:VAL:O	1:D:106:ASN:ND2	2.09	0.86
1:F:90:LEU:O	1:F:94:VAL:CG2	2.25	0.85
1:D:67:GLY:HA2	1:E:70:SER:OG	1.78	0.83
1:F:31:LEU:O	1:F:35:VAL:HG23	1.78	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:10:ARG:HH11	1:C:14:ALA:HB3	1.44	0.82
1:F:19:TRP:CE2	1:F:21:ASP:OD2	2.33	0.81
1:D:54:LEU:HD23	1:E:81:LEU:CD1	2.10	0.81
1:D:99:ALA:HB2	1:E:89:GLY:HA3	1.62	0.80
1:F:20:ARG:HB2	1:F:25:ARG:HH22	1.44	0.80
1:A:90:LEU:O	1:A:94:VAL:HG23	1.83	0.79
1:C:51:ILE:HD12	1:C:51:ILE:H	1.46	0.79
1:B:64:MET:HA	1:D:60:VAL:HA	1.63	0.79
1:F:9:LYS:O	1:F:13:GLY:HA3	1.83	0.78
1:A:10:ARG:CZ	1:A:15:VAL:HG21	2.13	0.77
1:E:34:ILE:HG22	1:E:83:TYR:CE1	2.20	0.76
1:F:87:VAL:HA	1:F:90:LEU:HD12	1.67	0.76
1:A:14:ALA:O	1:A:18:SER:HB2	1.86	0.75
1:B:22:PRO:HD2	1:B:25:ARG:HB2	1.69	0.75
1:A:51:ILE:H	1:A:51:ILE:HD12	1.52	0.74
1:A:16:ARG:HD2	1:A:17:THR:HG23	1.69	0.73
1:E:16:ARG:HH22	1:E:102:ILE:HG13	1.53	0.73
1:F:82:VAL:O	1:F:86:LEU:HD12	1.88	0.73
1:C:54:LEU:O	1:C:58:VAL:HG22	1.90	0.72
1:B:55:PHE:HA	1:D:81:LEU:HD21	1.70	0.72
1:C:101:ALA:O	1:C:105:ASN:HB2	1.89	0.72
1:D:46:GLU:OE2	1:D:71:PRO:HA	1.90	0.72
1:C:10:ARG:CZ	1:C:15:VAL:HG23	2.20	0.71
1:E:49:SER:HB2	1:F:51:ILE:HD13	1.71	0.71
1:B:46:GLU:C	1:B:47:LYS:HG2	2.16	0.71
1:A:46:GLU:C	1:A:47:LYS:HG2	2.15	0.71
1:B:46:GLU:O	1:B:47:LYS:HG2	1.89	0.71
1:C:44:LEU:HD12	1:C:44:LEU:H	1.56	0.70
1:C:46:GLU:O	1:C:47:LYS:HG2	1.92	0.70
1:B:34:ILE:CG2	1:B:83:TYR:HE1	2.04	0.69
1:A:10:ARG:CZ	1:A:15:VAL:CG2	2.68	0.69
1:A:66:ASN:ND2	1:A:69:LEU:O	2.26	0.69
1:A:46:GLU:OE2	1:A:71:PRO:HA	1.93	0.69
1:B:64:MET:HE1	1:D:81:LEU:HD23	1.74	0.68
1:D:55:PHE:HA	1:E:81:LEU:HD21	1.76	0.68
1:D:17:THR:HA	1:D:20:ARG:HH12	1.58	0.67
1:A:5:THR:O	1:A:8:PHE:CE2	2.47	0.67
1:B:20:ARG:HB3	1:B:25:ARG:HH22	1.60	0.66
1:D:51:ILE:HD12	1:D:51:ILE:H	1.59	0.66
1:F:31:LEU:HD22	1:F:94:VAL:HG21	1.78	0.66
1:F:40:ILE:O	1:F:44:LEU:HD13	1.96	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:44:LEU:HD12	1:F:44:LEU:H	1.61	0.65
1:C:10:ARG:HH11	1:C:14:ALA:CB	2.09	0.65
1:E:66:ASN:HB2	1:E:69:LEU:O	1.97	0.65
1:B:49:SER:HB2	1:C:51:ILE:HD13	1.79	0.64
1:F:29:LEU:O	1:F:33:ILE:HG22	1.98	0.64
1:E:94:VAL:O	1:E:98:LEU:HB2	1.98	0.64
1:A:98:LEU:O	1:A:102:ILE:HG12	1.98	0.64
1:F:51:ILE:H	1:F:51:ILE:HD12	1.63	0.63
1:A:20:ARG:O	1:A:25:ARG:HD2	1.98	0.63
1:F:44:LEU:HD12	1:F:44:LEU:N	2.12	0.63
1:C:10:ARG:NE	1:C:15:VAL:HG23	2.14	0.62
1:E:16:ARG:HH22	1:E:102:ILE:CG1	2.12	0.62
1:C:73:LEU:O	1:C:77:LYS:HG3	1.99	0.62
1:A:30:SER:O	1:A:34:ILE:HG12	1.99	0.62
1:D:64:MET:HA	1:E:60:VAL:HA	1.81	0.62
1:A:73:LEU:O	1:A:77:LYS:HG3	1.99	0.62
1:F:20:ARG:HB2	1:F:25:ARG:NH2	2.14	0.62
1:A:51:ILE:HG12	1:B:74:THR:HG23	1.82	0.61
1:D:22:PRO:HD2	1:D:25:ARG:HB2	1.82	0.61
1:B:46:GLU:OE2	1:B:71:PRO:HA	2.00	0.61
1:B:34:ILE:CG2	1:B:83:TYR:CE1	2.84	0.61
1:D:54:LEU:CD2	1:E:81:LEU:HD13	2.30	0.61
1:B:34:ILE:HG23	1:B:83:TYR:HE1	1.64	0.60
1:F:20:ARG:HH22	1:F:105:ASN:HD21	1.45	0.60
1:F:46:GLU:OE2	1:F:72:THR:N	2.31	0.60
1:B:17:THR:CG2	1:B:101:ALA:HB3	2.31	0.60
1:C:20:ARG:HE	1:C:101:ALA:HB2	1.67	0.60
1:D:42:TYR:CE2	1:D:56:TYR:HD2	2.19	0.60
1:D:103:VAL:HA	1:D:106:ASN:HD21	1.66	0.60
1:C:36:THR:O	1:C:40:ILE:HG13	2.01	0.60
1:D:66:ASN:ND2	1:D:69:LEU:O	2.32	0.59
1:B:20:ARG:O	1:B:25:ARG:NH2	2.34	0.59
1:D:10:ARG:HH12	1:D:15:VAL:HB	1.67	0.59
1:A:20:ARG:CB	1:A:25:ARG:HH22	2.09	0.59
1:F:49:SER:O	1:F:53:SER:OG	2.19	0.59
1:E:34:ILE:HG22	1:E:83:TYR:HE1	1.66	0.59
1:B:34:ILE:HG23	1:B:83:TYR:CE1	2.38	0.58
1:A:14:ALA:O	1:A:18:SER:CB	2.52	0.58
1:F:46:GLU:OE2	1:F:71:PRO:HA	2.03	0.58
1:C:51:ILE:H	1:C:51:ILE:CD1	2.14	0.58
1:F:20:ARG:NE	1:F:101:ALA:HB1	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:LYS:O	1:B:13:GLY:N	2.36	0.58
1:B:17:THR:HG22	1:B:101:ALA:HB3	1.85	0.58
1:D:36:THR:O	1:D:40:ILE:HG12	2.04	0.57
1:D:73:LEU:O	1:D:77:LYS:HG3	2.04	0.57
1:C:42:TYR:OH	1:C:80:THR:HG23	2.05	0.57
1:F:22:PRO:HD2	1:F:25:ARG:HG2	1.86	0.57
1:A:16:ARG:O	1:A:20:ARG:NH2	2.37	0.57
1:C:19:TRP:HD1	1:C:21:ASP:OD1	1.88	0.57
1:F:20:ARG:O	1:F:25:ARG:NH2	2.38	0.56
1:A:17:THR:HG22	1:A:105:ASN:ND2	2.20	0.56
1:E:51:ILE:HD12	1:E:51:ILE:H	1.70	0.56
1:F:36:THR:O	1:F:40:ILE:HD12	2.06	0.55
1:C:9:LYS:O	1:C:13:GLY:HA3	2.05	0.55
1:D:102:ILE:HD13	1:E:90:LEU:HD21	1.88	0.55
1:A:16:ARG:HH11	1:A:17:THR:HG23	1.72	0.55
1:F:20:ARG:O	1:F:22:PRO:HD3	2.06	0.55
1:F:20:ARG:HH22	1:F:105:ASN:CG	2.14	0.55
1:A:99:ALA:HB2	1:B:89:GLY:HA3	1.88	0.55
1:C:51:ILE:HD12	1:C:51:ILE:N	2.20	0.54
1:F:74:THR:O	1:F:78:ILE:HG13	2.07	0.54
1:B:54:LEU:HD23	1:D:81:LEU:HD13	1.90	0.54
1:D:51:ILE:HG21	1:E:74:THR:HG23	1.89	0.54
1:F:52:ASP:HB3	1:F:69:LEU:HD11	1.90	0.54
1:F:59:SER:HA	1:F:64:MET:HE2	1.90	0.53
1:D:98:LEU:O	1:D:102:ILE:HD12	2.08	0.53
1:F:11:PHE:O	1:F:16:ARG:HB2	2.08	0.53
1:A:41:PHE:HE1	1:E:5:THR:HG23	1.74	0.53
1:B:22:PRO:HG2	1:B:25:ARG:H	1.72	0.53
1:B:73:LEU:O	1:B:77:LYS:HG3	2.09	0.52
1:B:61:GLY:HA3	1:B:91:PHE:HD2	1.74	0.52
1:B:30:SER:O	1:B:33:ILE:HG22	2.10	0.52
1:B:94:VAL:O	1:B:98:LEU:HB2	2.09	0.52
1:D:9:LYS:HB2	1:D:9:LYS:NZ	2.25	0.52
1:D:58:VAL:HG12	1:D:62:LEU:HD12	1.92	0.52
1:F:52:ASP:N	1:F:52:ASP:OD1	2.43	0.52
1:A:46:GLU:C	1:A:47:LYS:CG	2.84	0.51
1:E:16:ARG:HD2	1:E:105:ASN:HD22	1.75	0.51
1:E:16:ARG:NH2	1:E:102:ILE:HG13	2.23	0.51
1:C:8:PHE:O	1:C:12:PHE:HB2	2.11	0.50
1:D:10:ARG:HH22	1:D:15:VAL:HB	1.75	0.50
1:D:26:GLY:O	1:D:29:LEU:HB2	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:61:GLY:HA3	1:F:91:PHE:HD2	1.76	0.50
1:C:90:LEU:O	1:C:94:VAL:HG23	2.11	0.50
1:A:62:LEU:HD23	1:A:63:PRO:HD2	1.93	0.50
1:D:51:ILE:HG12	1:E:74:THR:HG23	1.93	0.50
1:B:51:ILE:HD11	1:C:50:VAL:HG23	1.92	0.50
1:E:25:ARG:NE	1:E:97:SER:HB3	2.27	0.50
1:A:10:ARG:CZ	1:A:15:VAL:HG23	2.41	0.49
1:C:62:LEU:C	1:C:64:MET:H	2.20	0.49
1:F:61:GLY:HA3	1:F:91:PHE:CD2	2.47	0.49
1:B:25:ARG:HE	1:B:97:SER:HB3	1.77	0.49
1:D:9:LYS:HE3	1:E:34:ILE:HG23	1.94	0.49
1:D:17:THR:HA	1:D:20:ARG:HH11	1.75	0.49
1:B:51:ILE:O	1:B:55:PHE:N	2.32	0.49
1:D:58:VAL:HG21	1:E:81:LEU:HD22	1.93	0.49
1:C:20:ARG:O	1:C:25:ARG:NE	2.44	0.49
1:F:56:TYR:O	1:F:60:VAL:HG13	2.12	0.49
1:A:16:ARG:NH1	1:A:17:THR:HG23	2.27	0.49
1:A:94:VAL:O	1:A:98:LEU:HD12	2.12	0.49
1:B:46:GLU:OE2	1:B:72:THR:N	2.41	0.49
1:D:54:LEU:O	1:D:58:VAL:HG23	2.13	0.49
1:A:60:VAL:HA	1:E:64:MET:HA	1.95	0.48
1:D:13:GLY:O	1:E:86:LEU:HD21	2.13	0.48
1:B:34:ILE:HG22	1:B:83:TYR:HE1	1.75	0.48
1:D:10:ARG:NH1	1:D:15:VAL:HB	2.27	0.48
1:C:11:PHE:O	1:C:16:ARG:HB2	2.14	0.48
1:D:88:VAL:O	1:D:92:VAL:HG23	2.13	0.48
1:A:79:PHE:HD2	1:E:6:LEU:HB2	1.78	0.48
1:A:6:LEU:HD12	1:A:6:LEU:O	2.14	0.48
1:D:99:ALA:HA	1:D:102:ILE:CD1	2.43	0.48
1:C:10:ARG:NH1	1:C:14:ALA:CB	2.76	0.48
1:B:99:ALA:CB	1:D:89:GLY:HA3	2.44	0.48
1:C:88:VAL:O	1:C:92:VAL:HG23	2.14	0.48
1:F:10:ARG:HH12	1:F:15:VAL:HG23	1.79	0.48
1:A:89:GLY:HA3	1:E:99:ALA:HB2	1.95	0.47
1:C:58:VAL:HG23	1:C:64:MET:HE1	1.95	0.47
1:A:9:LYS:O	1:A:13:GLY:HA3	2.13	0.47
1:D:64:MET:HA	1:E:60:VAL:CA	2.44	0.47
1:B:79:PHE:O	1:B:83:TYR:HB3	2.14	0.47
1:F:79:PHE:CD1	1:F:79:PHE:C	2.92	0.47
1:C:40:ILE:O	1:C:44:LEU:HD11	2.05	0.47
1:F:78:ILE:O	1:F:82:VAL:HG23	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:20:ARG:HB3	1:C:25:ARG:NH1	2.30	0.47
1:E:85:ILE:O	1:E:88:VAL:HG12	2.15	0.47
1:B:12:PHE:O	1:B:16:ARG:NE	2.48	0.46
1:C:66:ASN:HB2	1:C:69:LEU:O	2.15	0.46
1:A:12:PHE:O	1:A:16:ARG:NH1	2.49	0.46
1:B:46:GLU:O	1:B:47:LYS:CG	2.61	0.46
1:F:10:ARG:NH1	1:F:15:VAL:HG23	2.30	0.46
1:A:26:GLY:HA2	1:A:29:LEU:HD23	1.97	0.46
1:C:98:LEU:O	1:C:102:ILE:HG12	2.16	0.46
1:A:82:VAL:HG12	1:E:14:ALA:HB3	1.96	0.46
1:B:14:ALA:HB2	1:D:82:VAL:HG13	1.97	0.46
1:A:19:TRP:C	1:A:20:ARG:HH21	2.23	0.46
1:B:99:ALA:HB2	1:D:89:GLY:HA3	1.97	0.46
1:C:47:LYS:HB2	1:D:73:LEU:HA	1.97	0.46
1:B:20:ARG:HB3	1:B:25:ARG:NH2	2.29	0.46
1:A:16:ARG:HD2	1:A:17:THR:CG2	2.45	0.45
1:A:20:ARG:CB	1:A:25:ARG:NH2	2.55	0.45
1:C:5:THR:OG1	1:C:8:PHE:CD2	2.62	0.45
1:D:16:ARG:HD3	1:D:17:THR:N	2.31	0.45
1:E:10:ARG:NH1	1:E:15:VAL:HG21	2.32	0.45
1:D:102:ILE:HD12	1:D:102:ILE:H	1.82	0.45
1:E:41:PHE:CZ	1:E:76:SER:HA	2.52	0.45
1:F:98:LEU:O	1:F:102:ILE:HG13	2.17	0.45
1:D:16:ARG:HD3	1:D:17:THR:HG23	1.98	0.45
1:E:34:ILE:HG22	1:E:83:TYR:CD1	2.51	0.45
1:D:11:PHE:O	1:D:16:ARG:CB	2.66	0.44
1:F:66:ASN:ND2	1:F:69:LEU:O	2.36	0.44
1:B:17:THR:HG21	1:B:98:LEU:O	2.18	0.44
1:F:55:PHE:O	1:F:59:SER:HB3	2.18	0.44
1:B:8:PHE:O	1:B:12:PHE:HB2	2.16	0.44
1:C:58:VAL:HG12	1:C:91:PHE:HZ	1.82	0.44
1:E:26:GLY:HA2	1:E:29:LEU:HD23	2.00	0.44
1:A:46:GLU:O	1:A:47:LYS:CG	2.66	0.43
1:C:20:ARG:HD2	1:C:104:GLN:NE2	2.33	0.43
1:E:9:LYS:O	1:E:13:GLY:HA3	2.18	0.43
1:B:49:SER:CB	1:C:51:ILE:HD13	2.47	0.43
1:D:74:THR:O	1:D:78:ILE:HG13	2.19	0.43
1:B:33:ILE:HD12	1:B:33:ILE:HA	1.95	0.43
1:D:51:ILE:H	1:D:51:ILE:CD1	2.27	0.43
1:D:16:ARG:HH21	1:D:105:ASN:HD22	1.65	0.43
1:B:61:GLY:HA3	1:B:91:PHE:CD2	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:ARG:NE	1:B:97:SER:HB3	2.33	0.43
1:B:38:ALA:HB2	1:B:83:TYR:CE1	2.54	0.43
1:E:105:ASN:O	1:E:106:ASN:ND2	2.52	0.43
1:C:103:VAL:C	1:C:105:ASN:H	2.26	0.43
1:D:10:ARG:NH2	1:D:15:VAL:HB	2.33	0.43
1:D:22:PRO:HG2	1:D:25:ARG:H	1.83	0.43
1:D:16:ARG:NE	1:D:17:THR:HG23	2.34	0.42
1:E:20:ARG:O	1:E:25:ARG:NH2	2.51	0.42
1:B:58:VAL:CG1	1:B:62:LEU:HD12	2.49	0.42
1:F:20:ARG:CD	1:F:101:ALA:HB1	2.49	0.42
1:C:12:PHE:HA	1:C:16:ARG:HD2	2.01	0.42
1:C:52:ASP:HB3	1:C:69:LEU:HD21	2.02	0.42
1:A:62:LEU:HD11	1:B:85:ILE:HD12	2.01	0.42
1:E:16:ARG:HE	1:E:16:ARG:HB3	1.65	0.42
1:F:28:VAL:HG22	1:F:94:VAL:HG13	2.01	0.42
1:D:100:SER:O	1:D:103:VAL:HG23	2.20	0.42
1:B:58:VAL:HB	1:D:81:LEU:HD22	2.02	0.42
1:D:98:LEU:O	1:D:101:ALA:HB3	2.20	0.42
1:B:31:LEU:HD13	1:B:90:LEU:HB2	2.02	0.42
1:A:41:PHE:CE1	1:E:5:THR:HG23	2.54	0.41
1:F:54:LEU:O	1:F:58:VAL:HG13	2.20	0.41
1:A:25:ARG:NE	1:A:97:SER:HB3	2.34	0.41
1:A:79:PHE:CD1	1:A:79:PHE:C	2.98	0.41
1:B:26:GLY:HA2	1:B:29:LEU:HD23	2.02	0.41
1:E:56:TYR:OH	1:E:71:PRO:HD3	2.20	0.41
1:D:57:ALA:O	1:D:60:VAL:HG22	2.20	0.41
1:C:78:ILE:O	1:C:82:VAL:HG23	2.20	0.41
1:F:13:GLY:H	1:F:16:ARG:HB2	1.85	0.41
1:D:58:VAL:CG2	1:E:81:LEU:HD22	2.51	0.41
1:E:75:LEU:O	1:E:78:ILE:HG13	2.20	0.41
1:B:66:ASN:ND2	1:B:69:LEU:O	2.37	0.41
1:E:80:THR:O	1:E:84:ALA:HB2	2.20	0.41
1:C:46:GLU:O	1:C:47:LYS:CG	2.66	0.41
1:D:6:LEU:HD12	1:D:10:ARG:HB2	2.02	0.41
1:A:20:ARG:CA	1:A:20:ARG:HE	2.32	0.40
1:A:22:PRO:HG2	1:A:25:ARG:H	1.86	0.40
1:C:46:GLU:C	1:C:47:LYS:CG	2.82	0.40
1:A:62:LEU:HA	1:A:63:PRO:HD3	1.94	0.40
1:D:13:GLY:HA3	1:E:86:LEU:HD11	2.04	0.40
1:A:83:TYR:HB2	1:E:9:LYS:HD3	2.03	0.40
1:C:5:THR:HG1	1:C:8:PHE:HD2	1.59	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:33:ILE:HD12	1:E:33:ILE:HA	1.90	0.40
1:B:63:PRO:O	1:B:64:MET:C	2.64	0.40
1:A:73:LEU:HD12	1:F:47:LYS:HB3	2.03	0.40
1:D:98:LEU:O	1:D:102:ILE:CD1	2.70	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	100/129 (78%)	89 (89%)	11 (11%)	0	100	100
1	B	100/129 (78%)	89 (89%)	9 (9%)	2 (2%)	6	29
1	C	100/129 (78%)	92 (92%)	8 (8%)	0	100	100
1	D	100/129 (78%)	88 (88%)	10 (10%)	2 (2%)	6	29
1	E	100/129 (78%)	90 (90%)	10 (10%)	0	100	100
1	F	100/129 (78%)	90 (90%)	9 (9%)	1 (1%)	12	42
All	All	600/774 (78%)	538 (90%)	57 (10%)	5 (1%)	16	46

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	7	MET
1	B	22	PRO
1	D	22	PRO
1	D	13	GLY
1	F	22	PRO

### 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	84/107 (78%)	66 (79%)	18 (21%)	1	7
1	B	84/107 (78%)	71 (84%)	13 (16%)	2	14
1	C	84/107 (78%)	73 (87%)	11 (13%)	4	19
1	D	84/107 (78%)	69 (82%)	15 (18%)	2	11
1	E	84/107 (78%)	65 (77%)	19 (23%)	1	6
1	F	84/107 (78%)	68 (81%)	16 (19%)	1	9
All	All	504/642 (78%)	412 (82%)	92 (18%)	2	10

All (92) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	6	LEU
1	A	9	LYS
1	A	10	ARG
1	A	12	PHE
1	A	16	ARG
1	A	20	ARG
1	A	44	LEU
1	A	47	LYS
1	A	49	SER
1	A	51	ILE
1	A	60	VAL
1	A	62	LEU
1	A	73	LEU
1	A	82	VAL
1	A	100	SER
1	A	103	VAL
1	A	106	ASN
1	B	6	LEU
1	B	8	PHE
1	B	9	LYS
1	B	17	THR

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	B	20	ARG
1	B	30	SER
1	B	34	ILE
1	B	44	LEU
1	B	49	SER
1	B	54	LEU
1	B	70	SER
1	B	88	VAL
1	B	91	PHE
1	C	9	LYS
1	C	16	ARG
1	C	44	LEU
1	C	47	LYS
1	C	51	ILE
1	C	60	VAL
1	C	73	LEU
1	C	80	THR
1	C	93	THR
1	C	98	LEU
1	C	105	ASN
1	D	6	LEU
1	D	9	LYS
1	D	10	ARG
1	D	15	VAL
1	D	16	ARG
1	D	17	THR
1	D	20	ARG
1	D	29	LEU
1	D	44	LEU
1	D	70	SER
1	D	73	LEU
1	D	75	LEU
1	D	86	LEU
1	D	102	ILE
1	D	106	ASN
1	E	5	THR
1	E	6	LEU
1	E	9	LYS
1	E	10	ARG
1	E	20	ARG
1	E	28	VAL
1	E	29	LEU

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Mol	Chain	Res	Type
1	E	30	SER
1	E	33	ILE
1	E	35	VAL
1	E	69	LEU
1	E	73	LEU
1	E	78	ILE
1	E	85	ILE
1	E	87	VAL
1	E	94	VAL
1	E	97	SER
1	E	98	LEU
1	E	102	ILE
1	F	6	LEU
1	F	9	LYS
1	F	12	PHE
1	F	16	ARG
1	F	18	SER
1	F	33	ILE
1	F	40	ILE
1	F	47	LYS
1	F	52	ASP
1	F	53	SER
1	F	58	VAL
1	F	62	LEU
1	F	75	LEU
1	F	82	VAL
1	F	94	VAL
1	F	103	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	106	ASN
1	D	105	ASN
1	D	106	ASN
1	E	104	GLN
1	E	105	ASN
1	E	106	ASN
1	F	104	GLN
1	F	105	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](#)

Of 3 ligands modelled in this entry, 3 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 [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, 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	102/129 (79%)	0.04	3 (2%) 53 32	66, 99, 136, 163	0
1	B	102/129 (79%)	0.06	3 (2%) 53 32	70, 104, 157, 201	0
1	C	102/129 (79%)	0.68	12 (11%) 9 9	91, 126, 178, 208	0
1	D	102/129 (79%)	0.13	7 (6%) 23 16	66, 102, 145, 178	0
1	E	102/129 (79%)	0.29	5 (4%) 35 21	74, 102, 157, 170	0
1	F	102/129 (79%)	0.63	14 (13%) 6 7	63, 119, 166, 197	0
All	All	612/774 (79%)	0.31	44 (7%) 21 15	63, 109, 164, 208	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	92	VAL	11.7
1	F	92	VAL	9.0
1	C	90	LEU	6.9
1	C	94	VAL	5.5
1	D	92	VAL	5.0
1	C	88	VAL	4.8
1	F	94	VAL	4.7
1	E	44	LEU	4.5
1	C	91	PHE	4.2
1	F	88	VAL	4.1
1	F	95	GLY	4.0
1	D	29	LEU	3.7
1	C	63	PRO	3.7
1	C	105	ASN	3.6
1	E	18	SER	3.5
1	F	96	GLY	3.3
1	E	14	ALA	3.2
1	D	44	LEU	3.1
1	F	91	PHE	3.0

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Mol	Chain	Res	Type	RSRZ
1	F	90	LEU	3.0
1	C	61	GLY	3.0
1	B	18	SER	2.8
1	E	50	VAL	2.8
1	E	39	THR	2.7
1	A	44	LEU	2.7
1	B	10	ARG	2.6
1	D	14	ALA	2.5
1	F	45	ALA	2.5
1	F	93	THR	2.5
1	F	75	LEU	2.4
1	F	61	GLY	2.3
1	C	44	LEU	2.3
1	D	27	ALA	2.3
1	F	65	GLY	2.3
1	C	93	THR	2.3
1	D	83	TYR	2.2
1	F	44	LEU	2.2
1	A	29	LEU	2.2
1	A	75	LEU	2.2
1	D	40	ILE	2.1
1	C	89	GLY	2.1
1	B	14	ALA	2.1
1	C	82	VAL	2.1
1	F	85	ILE	2.1

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

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

## 6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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( $\text{\AA}^2$ )	Q<0.9
2	CA	B	201	1/1	0.83	0.15	148,148,148,148	0
2	CA	E	201	1/1	0.91	0.08	76,76,76,76	0
2	CA	A	201	1/1	0.94	0.07	97,97,97,97	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.