



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 29, 2017 – 06:46 PM EDT

PDB ID : 1I1G
Title : CRYSTAL STRUCTURE OF THE LRP-LIKE TRANSCRIPTIONAL REGULATOR FROM THE ARCHAEON PYROCOCCUS FURIOSUS
Authors : Leonard, P.M.; Smits, S.H.J.; Sedelnikova, S.E.; Brinkman, A.B.; de Vos, W.M.; van der Oost, J.; Rice, D.W.; Rafferty, J.B.
Deposited on : 2001-02-01
Resolution : 2.90 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20029077
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20029077

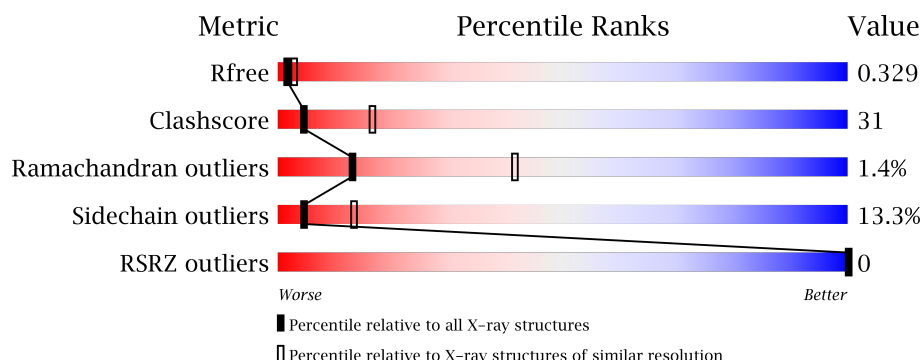
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

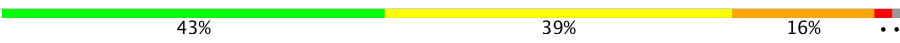
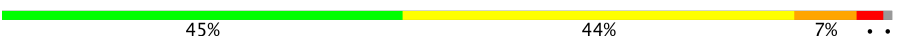
The reported resolution of this entry is 2.90 Å.

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}	100719	1586 (2.90-2.90)
Clashscore	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	141	
1	B	141	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 2138 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.

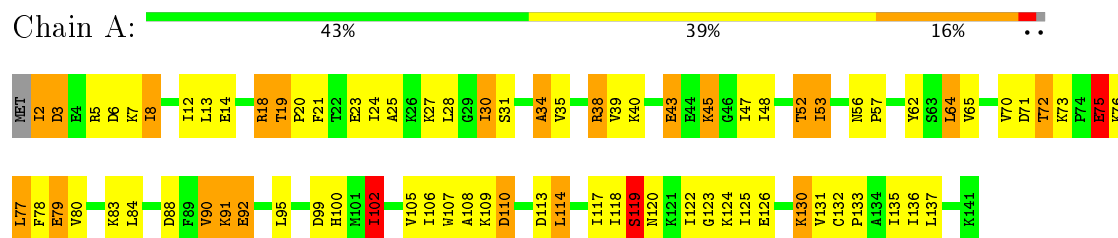
- Molecule 1 is a protein called TRANSCRIPTIONAL REGULATOR LRPA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	140	Total	C	N	O	S	0	0	0
			1069	693	170	203	3			
1	B	140	Total	C	N	O	S	0	0	0
			1069	693	170	203	3			

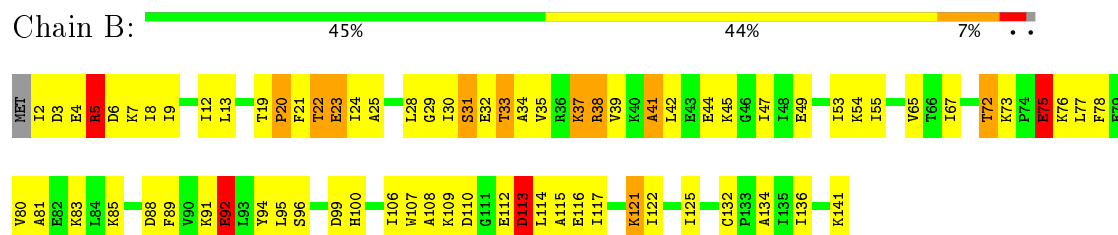
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: TRANSCRIPTIONAL REGULATOR LRPA



• Molecule 1: TRANSCRIPTIONAL REGULATOR LRPA



4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	101.32Å 101.32Å 245.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.90 19.87 – 2.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.90) 91.9 (19.87-2.90)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.99 (at 2.88Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.313 , 0.382 0.276 , 0.329	Depositor DCC
R_{free} test set	698 reflections (5.50%)	DCC
Wilson B-factor (Å ²)	54.2	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	2138	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.77	3/1082 (0.3%)	1.79	15/1460 (1.0%)
1	B	0.63	0/1082	1.78	17/1460 (1.2%)
All	All	0.70	3/2164 (0.1%)	1.79	32/2920 (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	5
1	B	0	7
All	All	0	12

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	3	ASP	N-CA	-7.05	1.32	1.46
1	A	3	ASP	C-O	6.82	1.36	1.23
1	A	2	ILE	C-N	-5.78	1.20	1.34

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	3	ASP	O-C-N	12.00	141.90	122.70
1	B	92	GLU	CB-CA-C	9.94	130.29	110.40
1	B	99	ASP	CB-CG-OD1	9.88	127.19	118.30
1	B	113	ASP	CB-CG-OD1	9.13	126.52	118.30
1	B	75	GLU	CA-CB-CG	8.60	132.31	113.40
1	B	5	ARG	CD-NE-CZ	8.37	135.31	123.60
1	B	5	ARG	NE-CZ-NH1	-7.40	116.60	120.30
1	B	38	ARG	NE-CZ-NH1	7.27	123.93	120.30
1	A	13	LEU	CA-CB-CG	6.90	131.17	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	137	LEU	CA-C-O	-6.86	105.70	120.10
1	B	44	GLU	CA-CB-CG	6.80	128.37	113.40
1	A	34	ALA	N-CA-CB	-6.79	100.60	110.10
1	A	110	ASP	CB-CG-OD1	6.59	124.24	118.30
1	A	62	TYR	CB-CG-CD2	-6.50	117.10	121.00
1	B	134	ALA	N-CA-CB	6.40	119.06	110.10
1	B	41	ALA	N-CA-CB	6.30	118.92	110.10
1	B	99	ASP	CB-CG-OD2	-6.29	112.64	118.30
1	A	64	LEU	CA-CB-CG	6.07	129.25	115.30
1	A	3	ASP	CA-C-N	-6.02	103.95	117.20
1	B	141	LYS	CA-C-O	-5.98	107.54	120.10
1	A	43	GLU	OE1-CD-OE2	5.71	130.15	123.30
1	A	62	TYR	CB-CG-CD1	5.66	124.40	121.00
1	A	3	ASP	N-CA-CB	-5.64	100.45	110.60
1	B	121	LYS	CA-C-O	-5.48	108.59	120.10
1	B	37	LYS	CB-CA-C	5.48	121.36	110.40
1	B	5	ARG	NE-CZ-NH2	5.27	122.93	120.30
1	B	113	ASP	CB-CG-OD2	-5.25	113.58	118.30
1	B	121	LYS	CB-CA-C	5.17	120.74	110.40
1	A	99	ASP	CB-CG-OD1	-5.16	113.65	118.30
1	A	19	THR	N-CA-CB	5.08	119.94	110.30
1	A	38	ARG	NE-CZ-NH1	-5.07	117.76	120.30
1	A	102	ILE	CB-CA-C	5.03	121.66	111.60

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	102	ILE	Mainchain
1	A	119	SER	Mainchain
1	A	75	GLU	Mainchain
1	A	90	VAL	Mainchain
1	A	92	GLU	Mainchain
1	B	113	ASP	Mainchain
1	B	23	GLU	Mainchain
1	B	33	THR	Mainchain
1	B	37	LYS	Mainchain
1	B	47	ILE	Mainchain
1	B	88	ASP	Mainchain
1	B	92	GLU	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	1069	0	1108	64	0
1	B	1069	0	1109	72	0
All	All	2138	0	2217	136	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 31.

All (136) 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:28:LEU:HB3	1:B:30:ILE:HD12	1.24	1.11
1:B:5:ARG:HH11	1:B:30:ILE:HD11	1.24	1.02
1:B:5:ARG:HD2	1:B:30:ILE:CD1	2.03	0.86
1:A:72:THR:HG21	1:A:77:LEU:HA	1.63	0.81
1:B:28:LEU:CB	1:B:30:ILE:HD12	2.11	0.80
1:B:30:ILE:HG22	1:B:34:ALA:HB3	1.64	0.79
1:B:28:LEU:HB3	1:B:30:ILE:CD1	2.11	0.79
1:B:29:GLY:C	1:B:30:ILE:HG13	2.03	0.78
1:A:21:PHE:HE2	1:A:39:VAL:HG21	1.49	0.76
1:A:30:ILE:HG22	1:A:31:SER:H	1.48	0.76
1:B:2:ILE:HD11	1:B:7:LYS:HG2	1.69	0.75
1:B:2:ILE:HD12	1:B:6:ASP:HB2	1.71	0.71
1:A:21:PHE:HA	1:A:24:ILE:HD12	1.71	0.71
1:B:5:ARG:HH11	1:B:30:ILE:CD1	2.03	0.70
1:B:5:ARG:NH1	1:B:30:ILE:HD11	2.01	0.70
1:A:106:ILE:CD1	1:A:122:ILE:HD11	2.21	0.70
1:A:21:PHE:CE2	1:A:39:VAL:HG21	2.27	0.69
1:A:24:ILE:HB	1:A:35:VAL:HG11	1.76	0.68
1:B:5:ARG:HD2	1:B:30:ILE:HD13	1.74	0.68
1:A:12:ILE:HD13	1:A:27:LYS:HD3	1.76	0.68
1:B:8:ILE:O	1:B:12:ILE:HG13	1.94	0.68
1:B:5:ARG:CD	1:B:30:ILE:HD11	2.25	0.66
1:B:75:GLU:HB2	1:B:76:LYS:HG3	1.77	0.66
1:B:5:ARG:HD2	1:B:30:ILE:HD11	1.79	0.64
1:A:120:ASN:O	1:A:124:LYS:HG3	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:38:ARG:HG2	1:B:38:ARG:HH11	1.63	0.63
1:A:5:ARG:HD3	1:A:30:ILE:HD11	1.80	0.62
1:B:41:ALA:O	1:B:45:LYS:HG3	1.99	0.62
1:B:92:GLU:HG3	1:B:94:TYR:OH	2.00	0.61
1:B:106:ILE:CD1	1:B:122:ILE:HD11	2.31	0.61
1:A:118:ILE:O	1:A:123:GLY:N	2.34	0.61
1:B:67:ILE:HD11	1:B:136:ILE:HD11	1.82	0.60
1:A:106:ILE:HD11	1:A:122:ILE:HD11	1.84	0.60
1:A:92:GLU:HG2	1:A:105:VAL:HB	1.84	0.60
1:B:65:VAL:HG12	1:B:136:ILE:HD12	1.83	0.60
1:A:118:ILE:HD11	1:A:133:PRO:HG3	1.85	0.59
1:A:84:LEU:HD21	1:A:122:ILE:HG12	1.84	0.59
1:B:21:PHE:CE2	1:B:39:VAL:HG21	2.38	0.59
1:B:21:PHE:HD2	1:B:24:ILE:HD12	1.68	0.59
1:A:76:LYS:O	1:A:80:VAL:HG23	2.03	0.59
1:A:40:LYS:O	1:A:43:GLU:HB2	2.02	0.58
1:A:118:ILE:HG12	1:A:131:VAL:HG11	1.85	0.57
1:B:100:HIS:CE1	1:B:132:CYS:HG	2.23	0.57
1:B:100:HIS:CE1	1:B:132:CYS:SG	2.98	0.57
1:B:30:ILE:CG2	1:B:34:ALA:HB3	2.33	0.57
1:A:6:ASP:OD1	1:A:38:ARG:HD2	2.05	0.56
1:A:19:THR:HG23	1:A:23:GLU:HB3	1.87	0.56
1:A:108:ALA:HB2	1:A:117:ILE:HD12	1.86	0.56
1:B:23:GLU:O	1:B:24:ILE:C	2.43	0.56
1:A:64:LEU:HD13	1:A:135:ILE:HD13	1.86	0.56
1:A:75:GLU:HG3	1:A:76:LYS:HG3	1.88	0.55
1:B:100:HIS:NE2	1:B:132:CYS:SG	2.78	0.55
1:B:39:VAL:O	1:B:42:LEU:HB2	2.06	0.55
1:A:31:SER:O	1:A:34:ALA:HB3	2.06	0.55
1:B:20:PRO:HG2	1:B:23:GLU:OE2	2.07	0.55
1:A:100:HIS:CE1	1:A:132:CYS:SG	2.99	0.55
1:A:100:HIS:CE1	1:A:132:CYS:HG	2.25	0.55
1:A:52:THR:HG23	1:A:53:ILE:N	2.22	0.54
1:B:72:THR:HG21	1:B:77:LEU:HA	1.91	0.53
1:B:20:PRO:HG2	1:B:23:GLU:CD	2.29	0.53
1:B:19:THR:HG22	1:B:20:PRO:O	2.10	0.52
1:B:108:ALA:HB2	1:B:117:ILE:HD12	1.91	0.52
1:B:2:ILE:HG13	1:B:3:ASP:N	2.25	0.52
1:B:4:GLU:HG3	1:B:8:ILE:HD11	1.92	0.51
1:B:8:ILE:HG22	1:B:12:ILE:HD11	1.93	0.51
1:B:77:LEU:HD11	1:B:95:LEU:CD2	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:2:ILE:C	1:A:3:ASP:O	2.45	0.50
1:B:91:LYS:HG3	1:B:107:TRP:CE2	2.47	0.50
1:A:91:LYS:HD2	1:A:107:TRP:CE3	2.47	0.50
1:B:29:GLY:O	1:B:30:ILE:HG13	2.11	0.50
1:B:25:ALA:HB2	1:B:35:VAL:HG21	1.94	0.49
1:A:25:ALA:CA	1:A:35:VAL:HG21	2.42	0.49
1:A:70:VAL:HB	1:A:102:ILE:HB	1.94	0.49
1:B:77:LEU:HD11	1:B:95:LEU:HD22	1.94	0.49
1:B:72:THR:OG1	1:B:80:VAL:HG21	2.13	0.49
1:A:110:ASP:OD2	1:A:113:ASP:N	2.39	0.48
1:A:64:LEU:HD13	1:A:135:ILE:CD1	2.43	0.48
1:A:65:VAL:HG12	1:A:136:ILE:HD12	1.96	0.48
1:A:24:ILE:HG21	1:A:35:VAL:HG13	1.95	0.48
1:B:21:PHE:HB3	1:B:35:VAL:HG11	1.94	0.48
1:B:5:ARG:NH1	1:B:30:ILE:CD1	2.69	0.47
1:B:4:GLU:HG3	1:B:8:ILE:CD1	2.44	0.47
1:B:83:LYS:HD2	1:B:125:ILE:HG23	1.96	0.47
1:B:29:GLY:C	1:B:30:ILE:CG1	2.81	0.47
1:A:84:LEU:CD2	1:A:122:ILE:HG12	2.44	0.47
1:B:22:THR:O	1:B:25:ALA:HB3	2.15	0.47
1:A:20:PRO:O	1:A:23:GLU:N	2.41	0.47
1:A:70:VAL:HG21	1:A:84:LEU:HD11	1.97	0.46
1:B:122:ILE:O	1:B:125:ILE:HG13	2.15	0.46
1:A:114:LEU:HD22	1:A:118:ILE:CD1	2.45	0.46
1:A:2:ILE:HD12	1:A:6:ASP:CB	2.46	0.46
1:A:119:SER:OG	1:A:120:ASN:N	2.46	0.46
1:B:106:ILE:HD11	1:B:122:ILE:HD11	1.96	0.46
1:A:25:ALA:N	1:A:35:VAL:HG21	2.30	0.46
1:A:28:LEU:HD23	1:A:28:LEU:HA	1.66	0.46
1:B:32:GLU:O	1:B:35:VAL:HB	2.16	0.45
1:A:125:ILE:O	1:A:126:GLU:C	2.52	0.45
1:B:21:PHE:HD2	1:B:21:PHE:HA	1.70	0.45
1:B:38:ARG:NH1	1:B:38:ARG:HG2	2.31	0.45
1:A:100:HIS:NE2	1:A:132:CYS:SG	2.90	0.45
1:A:72:THR:HG23	1:A:73:LYS:O	2.17	0.45
1:A:84:LEU:O	1:A:90:VAL:HG21	2.16	0.45
1:A:2:ILE:O	1:A:3:ASP:C	2.55	0.45
1:B:89:PHE:HE2	1:B:121:LYS:HE2	1.82	0.44
1:B:30:ILE:CG2	1:B:34:ALA:CB	2.95	0.44
1:B:9:ILE:HG22	1:B:13:LEU:CD1	2.47	0.44
1:A:5:ARG:HD3	1:A:30:ILE:CD1	2.46	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:LYS:HB3	1:A:47:ILE:HD12	2.00	0.44
1:A:19:THR:HG23	1:A:20:PRO:HD2	2.00	0.43
1:B:21:PHE:HZ	1:B:39:VAL:HG11	1.83	0.43
1:A:18:ARG:O	1:A:19:THR:C	2.55	0.43
1:B:9:ILE:HG22	1:B:13:LEU:HD12	2.00	0.43
1:A:91:LYS:HD2	1:A:107:TRP:CD2	2.53	0.43
1:B:67:ILE:CD1	1:B:136:ILE:HD11	2.48	0.43
1:A:18:ARG:H	1:A:18:ARG:HG3	1.62	0.43
1:B:80:VAL:O	1:B:81:ALA:C	2.57	0.43
1:A:30:ILE:HG22	1:A:31:SER:N	2.24	0.42
1:A:56:ASN:HA	1:A:57:PRO:HD2	1.65	0.42
1:A:2:ILE:HD12	1:A:6:ASP:C	2.39	0.42
1:B:112:GLU:O	1:B:115:ALA:HB3	2.19	0.42
1:B:78:PHE:CE2	1:B:95:LEU:HD11	2.55	0.42
1:A:118:ILE:HG23	1:A:131:VAL:HG21	2.01	0.42
1:B:25:ALA:HB1	1:B:30:ILE:O	2.19	0.42
1:A:7:LYS:O	1:A:8:ILE:C	2.58	0.42
1:B:8:ILE:HG21	1:B:28:LEU:HD21	2.01	0.42
1:B:31:SER:O	1:B:35:VAL:HG23	2.20	0.42
1:B:9:ILE:HD11	1:B:38:ARG:HD3	2.02	0.42
1:A:78:PHE:CE2	1:A:95:LEU:HD11	2.54	0.41
1:B:53:ILE:CD1	1:B:55:ILE:HD11	2.49	0.41
1:A:106:ILE:CG2	1:A:107:TRP:N	2.83	0.41
1:A:39:VAL:O	1:A:40:LYS:C	2.56	0.41
1:A:72:THR:HG21	1:A:77:LEU:CA	2.44	0.41
1:B:72:THR:HG23	1:B:73:LYS:O	2.21	0.41
1:A:79:GLU:O	1:A:83:LYS:HG3	2.20	0.41
1:B:2:ILE:CD1	1:B:7:LYS:HG2	2.45	0.41
1:B:109:LYS:HG3	1:B:113:ASP:OD2	2.21	0.41

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	138/141 (98%)	112 (81%)	23 (17%)	3 (2%)	8	29
1	B	138/141 (98%)	122 (88%)	15 (11%)	1 (1%)	25	60
All	All	276/282 (98%)	234 (85%)	38 (14%)	4 (1%)	13	41

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	75	GLU
1	A	130	LYS
1	A	30	ILE
1	B	20	PRO

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	113/123 (92%)	96 (85%)	17 (15%)	3	10
1	B	113/123 (92%)	100 (88%)	13 (12%)	6	20
All	All	226/246 (92%)	196 (87%)	30 (13%)	4	13

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ILE
1	A	14	GLU
1	A	18	ARG
1	A	45	LYS
1	A	48	ILE
1	A	52	THR
1	A	53	ILE
1	A	71	ASP
1	A	72	THR
1	A	77	LEU
1	A	79	GLU
1	A	88	ASP

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Mol	Chain	Res	Type
1	A	91	LYS
1	A	109	LYS
1	A	114	LEU
1	A	119	SER
1	A	130	LYS
1	B	5	ARG
1	B	22	THR
1	B	31	SER
1	B	33	THR
1	B	49	GLU
1	B	54	LYS
1	B	72	THR
1	B	75	GLU
1	B	85	LYS
1	B	96	SER
1	B	110	ASP
1	B	114	LEU
1	B	116	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	140/141 (99%)	-0.33	0 100 100	62, 62, 62, 62	0
1	B	140/141 (99%)	-0.38	0 100 100	62, 62, 62, 62	0
All	All	280/282 (99%)	-0.35	0 100 100	62, 62, 62, 62	0

There are no RSRZ outliers to report.

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 carbohydrates 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.