



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:36 PM UTC

PDB ID : 7R8J / pdb_00007r8j
Title : Crystal structure of Pseudooceanicola lipolyticus Argonaute bound to 5' p
guide DNA in the presence of Mg²⁺
Authors : Shin, Y.; Murakami, K.S.
Deposited on : 2021-06-26
Resolution : 2.70 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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

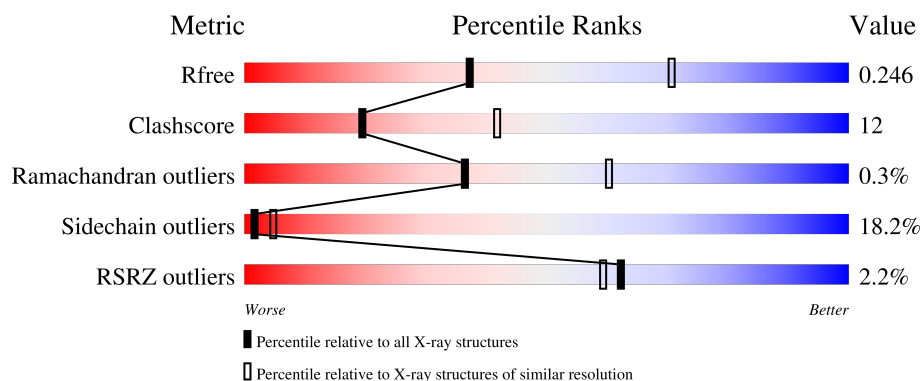
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

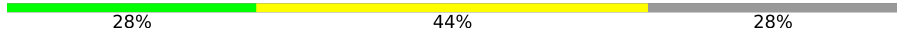
The reported resolution of this entry is 2.70 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	789	
2	T	18	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6264 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 Argonaute.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	764	Total	C	N	O	S	0	0	0
			5996	3806	1074	1087	29			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	13	THR	LEU	conflict	UNP A0A2M8J4C7
A	14	GLY	GLU	conflict	UNP A0A2M8J4C7
A	15	ALA	GLY	conflict	UNP A0A2M8J4C7
A	16	CYS	LEU	conflict	UNP A0A2M8J4C7
A	17	GLY	THR	conflict	UNP A0A2M8J4C7

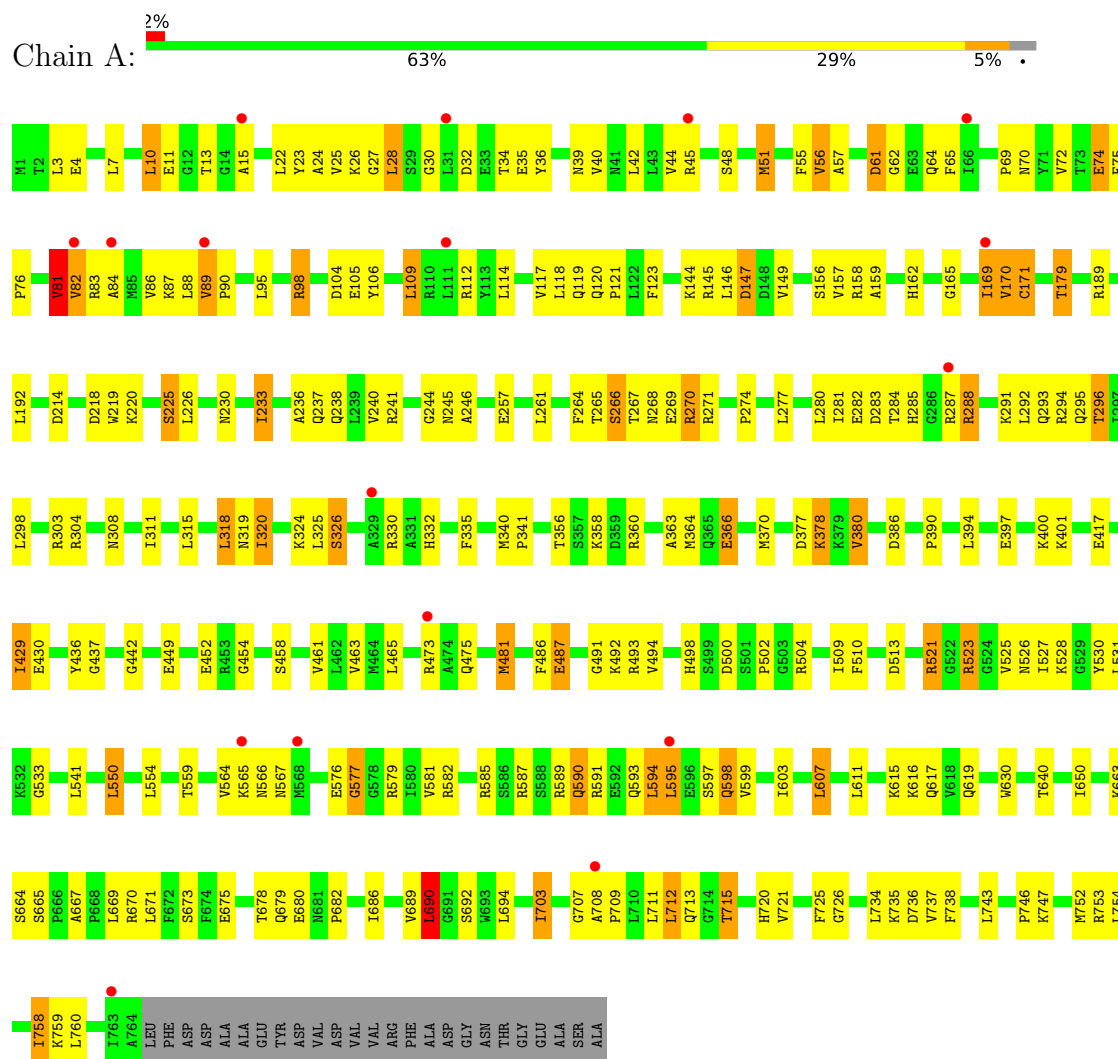
- Molecule 2 is DNA/RNA hybrid called DNA (5'-D(*TP*TP*AP*CP*TP*GP*CP*AP*CP*AP*GP*GP*TP*GP*AP*CP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	T	13	Total	C	N	O	P	0	0	0
			268	127	50	78	13			

3 Residue-property plots

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: Argonaute



• Molecule 2: DNA (5'-D(*TP*TP*AP*CP*TP*GP*CP*AP*CP*AP*GP*GP*TP*GP*AP*CP*GP*A)-3')



T1	T2	T3	T4	T5	G6	G7	A8	DC	DA	DG	DG	DT	G14	A15	C16	G17	A18
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4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, α , β , γ	199.19Å 199.19Å 199.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.54 – 2.70 44.54 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.5 (44.54-2.70) 99.4 (44.54-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.40 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.206 , 0.244 0.209 , 0.246	Depositor DCC
R_{free} test set	1985 reflections (5.49%)	wwPDB-VP
Wilson B-factor (Å ²)	80.8	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 59.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.027 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6264	wwPDB-VP
Average B, all atoms (Å ²)	104.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.93% 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	0.40	0/6130	0.66	7/8289 (0.1%)
2	T	0.48	0/299	0.69	0/455
All	All	0.41	0/6429	0.66	7/8744 (0.1%)

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	4

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	735	LYS	CA-C-N	-7.88	107.41	122.06
1	A	735	LYS	C-N-CA	-7.88	107.41	122.06
1	A	576	GLU	CA-C-N	-5.48	110.67	121.41
1	A	576	GLU	C-N-CA	-5.48	110.67	121.41
1	A	147	ASP	CB-CA-C	-5.29	110.04	117.23
1	A	690	LEU	CA-C-N	-5.19	109.23	122.78
1	A	690	LEU	C-N-CA	-5.19	109.23	122.78

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	436	TYR	Peptide
1	A	491	GLY	Peptide
1	A	577	GLY	Peptide
1	A	707	GLY	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5996	0	6039	144	1
2	T	268	0	148	13	0
All	All	6264	0	6187	144	1

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 12.

All (144) 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:746:PRO:HD2	1:A:752:MET:HE1	1.56	0.88
1:A:590:GLN:HG3	1:A:593:GLN:HB3	1.65	0.79
1:A:554:LEU:HD22	1:A:736:ASP:HB2	1.67	0.77
1:A:144:LYS:NZ	1:A:295:GLN:OE1	2.18	0.75
1:A:270:ARG:H	1:A:270:ARG:HD3	1.54	0.71
1:A:567:ASN:ND2	1:A:589:ARG:O	2.26	0.68
1:A:593:GLN:HG2	1:A:594:LEU:H	1.57	0.67
1:A:754:LEU:HD22	1:A:758:ILE:HD11	1.77	0.66
1:A:35:GLU:O	1:A:39:ASN:ND2	2.29	0.65
1:A:311:ILE:HD13	1:A:315:LEU:HD12	1.78	0.64
1:A:303:ARG:NH1	2:T:6:DG:OP1	2.30	0.64
1:A:145:ARG:NH2	1:A:147:ASP:O	2.31	0.64
1:A:400:LYS:HE3	1:A:430:GLU:HB3	1.79	0.63
1:A:298:LEU:O	1:A:715:THR:HG22	1.99	0.62
1:A:594:LEU:HD22	1:A:599:VAL:HG23	1.80	0.62
1:A:15:ALA:HA	1:A:98:ARG:HH22	1.66	0.61
1:A:28:LEU:HD13	1:A:30:GLY:H	1.66	0.60
1:A:189:ARG:NH2	1:A:218:ASP:HA	2.16	0.60
1:A:233:ILE:HD12	1:A:237:GLN:HG2	1.81	0.60
1:A:663:LYS:O	1:A:753:ARG:NH2	2.34	0.60
1:A:189:ARG:HH22	1:A:218:ASP:HA	1.65	0.60
1:A:24:ALA:HB3	1:A:89:VAL:HG13	1.83	0.59
1:A:24:ALA:O	1:A:88:LEU:HA	2.02	0.59
1:A:117:VAL:HB	1:A:320:ILE:HG21	1.84	0.59
1:A:106:TYR:O	1:A:109:LEU:HB2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:437:GLY:HA2	1:A:442:GLY:HA3	1.85	0.59
1:A:303:ARG:HD2	1:A:715:THR:HG21	1.85	0.58
1:A:74:GLU:OE1	1:A:75:PHE:N	2.37	0.58
1:A:394:LEU:HD13	1:A:429:ILE:HG22	1.85	0.57
1:A:162:HIS:HE1	1:A:332:HIS:HD2	1.50	0.57
1:A:703:ILE:H	1:A:703:ILE:HD12	1.69	0.57
1:A:607:LEU:HD22	1:A:611:LEU:HG	1.87	0.56
1:A:593:GLN:HG2	1:A:594:LEU:N	2.20	0.56
1:A:283:ASP:H	1:A:287:ARG:NH2	2.04	0.56
1:A:51:MET:HE2	1:A:55:PHE:HZ	1.71	0.55
1:A:40:VAL:O	1:A:44:VAL:HG23	2.06	0.55
1:A:112:ARG:HH22	2:T:14:DG:H3'	1.72	0.55
1:A:554:LEU:HD22	1:A:736:ASP:CB	2.36	0.55
1:A:157:VAL:O	1:A:158:ARG:HD2	2.06	0.54
1:A:23:TYR:O	1:A:65:PHE:HA	2.07	0.54
1:A:675:GLU:HB3	1:A:686:ILE:HD13	1.89	0.54
1:A:179:THR:HG23	1:A:280:LEU:O	2.08	0.54
1:A:11:GLU:HB2	1:A:326:SER:HB3	1.91	0.53
1:A:747:LYS:HE3	2:T:2:DT:O2	2.08	0.53
1:A:120:GLN:HG2	2:T:15:DA:H62	1.75	0.52
1:A:121:PRO:HG2	1:A:318:LEU:HG	1.91	0.52
1:A:303:ARG:HH11	1:A:715:THR:HG23	1.74	0.52
1:A:274:PRO:HG2	1:A:277:LEU:HD12	1.91	0.52
1:A:713:GLN:NE2	2:T:5:DT:O2	2.42	0.52
1:A:593:GLN:HG3	1:A:630:TRP:CD2	2.45	0.51
1:A:364:MET:HB2	1:A:709:PRO:HG2	1.93	0.50
1:A:703:ILE:HG13	1:A:734:LEU:HD22	1.91	0.50
1:A:708:ALA:HB3	1:A:709:PRO:HD3	1.93	0.50
1:A:56:VAL:HG22	1:A:112:ARG:HD2	1.94	0.50
1:A:42:LEU:HD23	1:A:82:VAL:HG13	1.94	0.49
1:A:360:ARG:O	1:A:364:MET:HG2	2.13	0.49
1:A:32:ASP:C	1:A:34:THR:H	2.19	0.49
1:A:162:HIS:CE1	1:A:332:HIS:HD2	2.29	0.49
1:A:680:GLU:O	1:A:682:PRO:HD3	2.13	0.49
1:A:669:LEU:HD21	1:A:671:LEU:HD21	1.94	0.48
1:A:703:ILE:HG21	1:A:738:PHE:HB2	1.95	0.48
1:A:593:GLN:HG3	1:A:630:TRP:CG	2.49	0.48
1:A:754:LEU:HD13	1:A:758:ILE:HD11	1.95	0.48
1:A:340:MET:HE2	1:A:360:ARG:NH1	2.28	0.48
1:A:170:VAL:HG21	1:A:667:ALA:HB1	1.94	0.48
1:A:244:GLY:O	1:A:246:ALA:N	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:689:VAL:HG22	1:A:690:LEU:H	1.79	0.48
1:A:219:TRP:HB2	1:A:225:SER:HB3	1.96	0.48
1:A:288:ARG:HE	1:A:288:ARG:HB2	1.49	0.48
1:A:670:ARG:NH1	1:A:715:THR:O	2.47	0.47
1:A:156:SER:O	1:A:171:CYS:HA	2.15	0.47
1:A:34:THR:HG21	1:A:595:LEU:HD22	1.96	0.47
1:A:266:SER:OG	2:T:16:DC:OP1	2.32	0.47
1:A:303:ARG:HD2	1:A:715:THR:CG2	2.44	0.47
1:A:3:LEU:HD11	1:A:692:SER:HB3	1.96	0.47
1:A:270:ARG:HD3	1:A:270:ARG:N	2.25	0.47
1:A:587:ARG:O	1:A:598:GLN:NE2	2.47	0.47
1:A:665:SER:OG	1:A:720:HIS:ND1	2.32	0.47
1:A:162:HIS:HB3	1:A:165:GLY:O	2.14	0.47
1:A:27:GLY:HA2	1:A:86:VAL:HG12	1.98	0.46
1:A:61:ASP:HB2	1:A:62:GLY:H	1.66	0.46
1:A:725:PHE:CG	1:A:726:GLY:N	2.84	0.46
1:A:81:VAL:HG12	1:A:82:VAL:H	1.81	0.45
1:A:500:ASP:O	1:A:504:ARG:HG3	2.16	0.45
1:A:120:GLN:HG2	2:T:15:DA:N6	2.31	0.45
1:A:296:THR:HG23	2:T:6:DG:H4'	1.99	0.45
1:A:589:ARG:HB2	1:A:590:GLN:NE2	2.31	0.45
1:A:481:MET:HE3	1:A:481:MET:HB2	1.84	0.45
1:A:577:GLY:H	1:A:579:ARG:HH11	1.63	0.45
1:A:550:LEU:HD22	1:A:743:LEU:HD11	1.99	0.45
1:A:39:ASN:HA	1:A:42:LEU:HB3	1.99	0.45
1:A:358:LYS:HB2	1:A:363:ALA:HB2	1.99	0.44
1:A:10:LEU:HD12	1:A:10:LEU:HA	1.56	0.44
1:A:502:PRO:HD3	1:A:530:TYR:CZ	2.52	0.44
1:A:56:VAL:HG12	1:A:57:ALA:H	1.82	0.44
1:A:429:ILE:HD13	1:A:429:ILE:HA	1.61	0.44
1:A:390:PRO:HB2	1:A:458:SER:HB2	2.00	0.44
1:A:690:LEU:HD22	1:A:709:PRO:HD2	2.00	0.44
1:A:465:LEU:O	1:A:498:HIS:HA	2.18	0.44
1:A:293:GLN:O	1:A:296:THR:HG22	2.18	0.43
1:A:690:LEU:HD13	1:A:690:LEU:HA	1.84	0.43
1:A:270:ARG:H	1:A:270:ARG:HH11	1.66	0.43
1:A:521:ARG:HE	1:A:521:ARG:HB2	1.70	0.43
1:A:56:VAL:HG13	1:A:112:ARG:HG3	2.01	0.43
1:A:171:CYS:SG	1:A:311:ILE:HD11	2.58	0.43
1:A:236:ALA:O	1:A:240:VAL:HG23	2.19	0.43
1:A:378:LYS:H	1:A:378:LYS:HG3	1.59	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:615:LYS:HD3	1:A:615:LYS:HA	1.51	0.43
1:A:341:PRO:HD3	1:A:364:MET:HE1	2.00	0.42
1:A:366:GLU:O	1:A:370:MET:HB2	2.19	0.42
1:A:23:TYR:CE2	1:A:90:PRO:HG3	2.54	0.42
1:A:69:PRO:O	1:A:72:VAL:HG22	2.19	0.42
1:A:123:PHE:CZ	2:T:17:DG:N2	2.87	0.42
1:A:673:SER:HB3	1:A:686:ILE:HG13	2.01	0.42
1:A:564:VAL:CG2	1:A:594:LEU:HD12	2.50	0.42
1:A:112:ARG:NH2	2:T:14:DG:H3'	2.33	0.42
1:A:233:ILE:HG23	1:A:237:GLN:HB3	2.02	0.41
1:A:487:GLU:OE2	1:A:493:ARG:HA	2.19	0.41
1:A:712:LEU:HD22	1:A:712:LEU:HA	1.85	0.41
1:A:159:ALA:HA	1:A:169:ILE:HD13	2.02	0.41
1:A:114:LEU:HD23	1:A:159:ALA:HB1	2.02	0.41
1:A:377:ASP:HB3	1:A:380:VAL:HG13	2.02	0.41
1:A:752:MET:HE3	1:A:752:MET:HB2	1.88	0.41
1:A:28:LEU:HB2	1:A:84:ALA:HB3	2.01	0.41
1:A:220:LYS:HD3	1:A:257:GLU:HA	2.01	0.41
1:A:452:GLU:C	1:A:454:GLY:H	2.28	0.41
1:A:712:LEU:HB3	1:A:713:GLN:OE1	2.21	0.41
1:A:118:LEU:CD1	1:A:169:ILE:HD11	2.51	0.41
1:A:264:PHE:HE1	1:A:270:ARG:HB3	1.85	0.41
1:A:119:GLN:O	1:A:123:PHE:HB2	2.21	0.41
1:A:486:PHE:HB3	1:A:492:LYS:HB2	2.03	0.41
1:A:4:GLU:OE2	1:A:304:ARG:HD2	2.20	0.41
1:A:458:SER:O	1:A:492:LYS:HE3	2.21	0.41
1:A:523:ARG:O	1:A:523:ARG:HD2	2.20	0.41
1:A:533:GLY:HA3	2:T:1:DT:O2	2.21	0.41
1:A:581:VAL:O	1:A:582:ARG:HD3	2.21	0.41
1:A:120:GLN:HB2	1:A:121:PRO:HD3	2.02	0.41
1:A:45:ARG:CZ	2:T:14:DG:N7	2.83	0.40
1:A:75:PHE:HA	1:A:76:PRO:HD3	1.92	0.40
1:A:296:THR:HG21	2:T:6:DG:H2''	2.02	0.40
1:A:527:ILE:O	1:A:531:LEU:HG	2.21	0.40
1:A:603:ILE:HG22	1:A:640:THR:HG21	2.02	0.40
1:A:565:LYS:O	1:A:566:ASN:HB2	2.21	0.40
1:A:36:TYR:O	1:A:40:VAL:HG23	2.22	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:241:ARG:NH1	1:A:397:GLU:OE2[6_555]	2.15	0.05

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	762/789 (97%)	710 (93%)	50 (7%)	2 (0%)	36 60

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	81	VAL
1	A	245	ASN

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	639/657 (97%)	523 (82%)	116 (18%)	2 5

All (116) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	10	LEU
1	A	13	THR
1	A	22	LEU

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Mol	Chain	Res	Type
1	A	25	VAL
1	A	26	LYS
1	A	28	LEU
1	A	48	SER
1	A	51	MET
1	A	56	VAL
1	A	61	ASP
1	A	64	GLN
1	A	70	ASN
1	A	74	GLU
1	A	81	VAL
1	A	82	VAL
1	A	83	ARG
1	A	87	LYS
1	A	89	VAL
1	A	95	LEU
1	A	98	ARG
1	A	104	ASP
1	A	105	GLU
1	A	109	LEU
1	A	146	LEU
1	A	149	VAL
1	A	169	ILE
1	A	170	VAL
1	A	171	CYS
1	A	179	THR
1	A	192	LEU
1	A	214	ASP
1	A	225	SER
1	A	226	LEU
1	A	230	ASN
1	A	233	ILE
1	A	238	GLN
1	A	261	LEU
1	A	265	THR
1	A	266	SER
1	A	267	THR
1	A	268	ASN
1	A	269	GLU
1	A	270	ARG
1	A	271	ARG
1	A	281	ILE

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Mol	Chain	Res	Type
1	A	282	GLU
1	A	284	THR
1	A	285	HIS
1	A	288	ARG
1	A	291	LYS
1	A	292	LEU
1	A	294	ARG
1	A	296	THR
1	A	308	ASN
1	A	318	LEU
1	A	319	ASN
1	A	320	ILE
1	A	324	LYS
1	A	325	LEU
1	A	326	SER
1	A	330	ARG
1	A	335	PHE
1	A	356	THR
1	A	366	GLU
1	A	378	LYS
1	A	380	VAL
1	A	386	ASP
1	A	401	LYS
1	A	417	GLU
1	A	429	ILE
1	A	449	GLU
1	A	461	VAL
1	A	463	VAL
1	A	473	ARG
1	A	475	GLN
1	A	481	MET
1	A	487	GLU
1	A	494	VAL
1	A	509	ILE
1	A	510	PHE
1	A	513	ASP
1	A	521	ARG
1	A	523	ARG
1	A	525	VAL
1	A	526	ASN
1	A	528	LYS
1	A	541	LEU

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Mol	Chain	Res	Type
1	A	550	LEU
1	A	559	THR
1	A	585	ARG
1	A	590	GLN
1	A	591	ARG
1	A	594	LEU
1	A	595	LEU
1	A	597	SER
1	A	598	GLN
1	A	607	LEU
1	A	616	LYS
1	A	617	GLN
1	A	619	GLN
1	A	650	ILE
1	A	664	SER
1	A	678	THR
1	A	679	GLN
1	A	690	LEU
1	A	694	LEU
1	A	703	ILE
1	A	711	LEU
1	A	712	LEU
1	A	715	THR
1	A	721	VAL
1	A	737	VAL
1	A	758	ILE
1	A	759	LYS
1	A	760	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	126	HIS
1	A	162	HIS
1	A	238	GLN
1	A	319	ASN
1	A	332	HIS
1	A	365	GLN
1	A	416	GLN
1	A	443	GLN
1	A	555	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	764/789 (96%)	0.06	17 (2%) 62 59	54, 93, 181, 252	0
2	T	13/18 (72%)	-0.40	0 100 100	65, 96, 125, 162	0
All	All	777/807 (96%)	0.05	17 (2%) 62 59	54, 93, 181, 252	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	329	ALA	4.2
1	A	763	ILE	2.9
1	A	708	ALA	2.7
1	A	15	ALA	2.7
1	A	568	MET	2.5
1	A	66	ILE	2.4
1	A	287	ARG	2.3
1	A	169	ILE	2.3
1	A	89	VAL	2.2
1	A	31	LEU	2.2
1	A	82	VAL	2.2
1	A	565	LYS	2.2
1	A	473	ARG	2.1
1	A	45	ARG	2.1
1	A	595	LEU	2.1
1	A	84	ALA	2.1
1	A	111	LEU	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 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.