



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 08:54 am BST

PDB ID : 5XOW
Title : Crystal structure of T. thermophilus Argonaute protein complexed with a bulge 6'A7' on the target strand
Authors : Sheng, G.; Wang, J.; Zhao, H.; Wang, Y.
Deposited on : 2017-05-31
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

<https://www.wwpdb.org/validation/2017/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)	:	1.13
EDS	:	2.11
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)	:	2.11

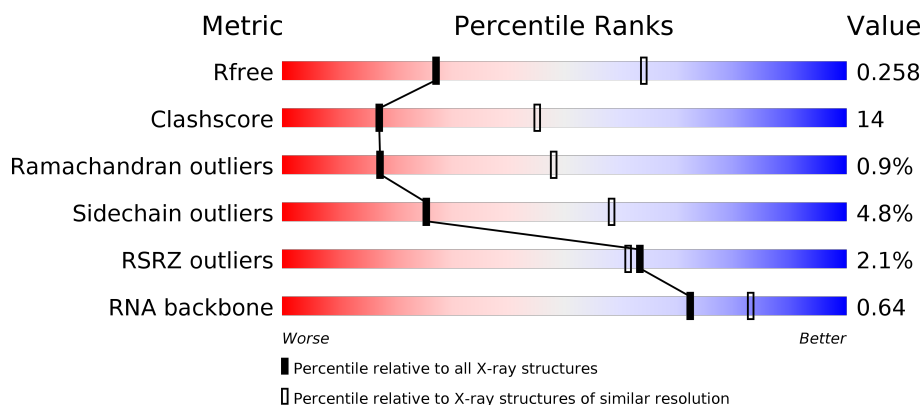
1 Overall quality at a glance

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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)
RNA backbone	3102	1007 (3.16-2.64)

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 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	685	<div> <div>2%</div> <div>67%</div> <div>29%</div> <div>..</div> </div>
2	C	21	<div> <div>52%</div> <div>43%</div> <div>5%</div> </div>
3	G	20	<div> <div>5%</div> <div>75%</div> <div>15%</div> <div>10%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6040 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 TtAgo (D546N).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	684	Total	C	N	O	S	0	0	0
			5241	3350	979	906	6			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	546	ASN	ASP	engineered mutation	UNP Q746M7

- Molecule 2 is a DNA chain called DNA (5'-D(P*(TD)P*GP*AP*GP*GP*TP*AP*GP*TP*AP*GP*GP*TP*TP*GP*TP*AP*TP*AP*GP*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	21	Total	C	N	O	P	0	0	0
			442	210	81	130	21			

- Molecule 3 is a RNA chain called RNA (5'-R(P*UP*AP*CP*AP*AP*CP*CP*UP*AP*CP*UP*AP*AP*CP*CP*UP*CP*G)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	18	Total	C	N	O	P	0	0	0
			356	159	59	120	18			

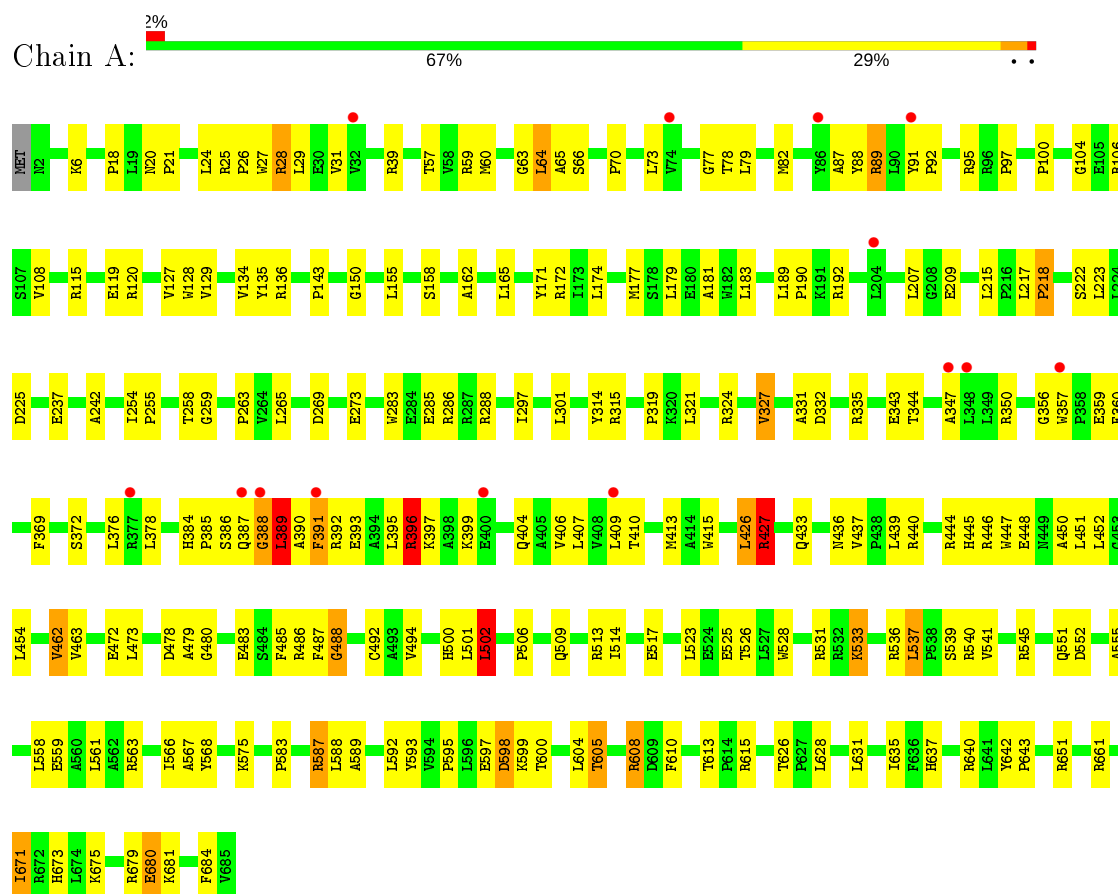
- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	Mg	0	0
			1	1		

3 Residue-property plots

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.

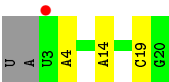
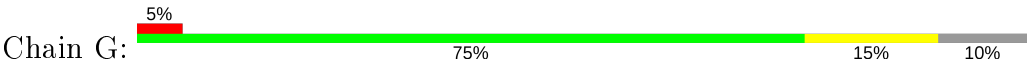
• Molecule 1: TtAgo (D546N)



• Molecule 2: DNA (5'-D(P*(TD)P*GP*AP*GP*GP*TP*AP*GP*TP*AP*GP*GP*TP*TP*GP*TP*AP*TP*AP*GP*T)-3')



• Molecule 3: RNA (5'-R(P*UP*AP*CP*AP*AP*CP*CP*UP*AP*CP*UP*AP*AP*CP*CP*UP*CP*G)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 3 2	Depositor
Cell constants a, b, c, α , β , γ	200.82Å 200.82Å 200.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.71 – 2.90 48.71 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.3 (48.71-2.90) 99.3 (48.71-2.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.214 , 0.257 0.217 , 0.258	Depositor DCC
R_{free} test set	1563 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	65.9	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 49.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6040	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.92% 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

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

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	0.42	1/5369 (0.0%)	0.81	18/7308 (0.2%)
2	C	0.75	1/496 (0.2%)	0.95	0/765
3	G	0.25	0/395	0.78	0/611
All	All	0.44	2/6260 (0.0%)	0.82	18/8684 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	DT	OP3-P	-10.86	1.48	1.61
1	A	528	TRP	CB-CG	-6.55	1.38	1.50

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	389	LEU	CB-CG-CD2	-22.00	73.60	111.00
1	A	389	LEU	CA-CB-CG	-10.16	91.93	115.30
1	A	502	LEU	CB-CG-CD1	-8.05	97.32	111.00
1	A	537	LEU	CA-CB-CG	7.25	131.97	115.30
1	A	396	ARG	NE-CZ-NH1	6.39	123.50	120.30
1	A	533	LYS	CA-CB-CG	6.22	127.08	113.40
1	A	533	LYS	N-CA-CB	6.11	121.59	110.60
1	A	427	ARG	NE-CZ-NH1	6.07	123.33	120.30
1	A	64	LEU	CA-CB-CG	6.00	129.10	115.30
1	A	39	ARG	NE-CZ-NH1	-5.93	117.33	120.30
1	A	396	ARG	CG-CD-NE	-5.89	99.44	111.80
1	A	396	ARG	NE-CZ-NH2	-5.83	117.38	120.30
1	A	218	PRO	N-CA-CB	5.55	109.97	103.30
1	A	396	ARG	CB-CG-CD	5.55	126.02	111.60
1	A	89	ARG	NE-CZ-NH1	-5.43	117.58	120.30
1	A	598	ASP	N-CA-C	-5.38	96.47	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	388	GLY	N-CA-C	-5.28	99.89	113.10
1	A	396	ARG	CD-NE-CZ	5.17	130.84	123.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5241	0	5198	156	0
2	C	442	0	240	9	0
3	G	356	0	183	3	0
4	C	1	0	0	0	0
All	All	6040	0	5621	161	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 14.

All (161) 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:407:LEU:HD13	1:A:454:LEU:HD22	1.60	0.82
1:A:537:LEU:HD13	1:A:566:ILE:HD11	1.62	0.82
1:A:393:GLU:HB3	1:A:396:ARG:CZ	2.10	0.80
1:A:598:ASP:O	1:A:599:LYS:HG2	1.82	0.80
1:A:540:ARG:NH1	1:A:626:THR:OG1	2.18	0.76
1:A:480:GLY:HA2	1:A:487:PHE:H	1.54	0.71
1:A:473:LEU:HB3	1:A:541:VAL:HG12	1.73	0.71
1:A:158:SER:HB2	1:A:162:ALA:H	1.57	0.69
1:A:28:ARG:NH1	1:A:29:LEU:O	2.27	0.68
1:A:436:ASN:HD21	3:G:14:A:H2	1.42	0.67
1:A:285:GLU:OE2	1:A:288:ARG:NH2	2.28	0.67
1:A:427:ARG:HG2	1:A:673:HIS:CD2	2.31	0.66
1:A:393:GLU:HA	1:A:396:ARG:HB3	1.77	0.65
1:A:321:LEU:HB3	1:A:327:VAL:HG12	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:3:DA:H2'	2:C:4:DG:C8	2.33	0.63
1:A:350:ARG:NH2	1:A:356:GLY:O	2.32	0.62
1:A:386:SER:C	1:A:388:GLY:H	2.03	0.62
1:A:344:THR:HG23	1:A:404:GLN:OE1	2.00	0.62
1:A:559:GLU:O	1:A:563:ARG:HG3	1.99	0.61
1:A:396:ARG:HG2	1:A:397:LYS:N	2.15	0.61
1:A:388:GLY:O	1:A:391:PHE:HD1	1.84	0.60
1:A:390:ALA:HA	1:A:393:GLU:OE1	2.02	0.59
1:A:155:LEU:HD13	1:A:165:LEU:HD13	1.84	0.59
1:A:513:ARG:NH2	1:A:551:GLN:O	2.36	0.58
1:A:27:TRP:HA	1:A:95:ARG:HG2	1.84	0.58
1:A:315:ARG:NH2	1:A:589:ALA:O	2.36	0.58
1:A:478:ASP:O	1:A:488:GLY:HA2	2.03	0.58
1:A:57:THR:HG21	1:A:73:LEU:HD21	1.85	0.57
1:A:18:PRO:HA	1:A:162:ALA:HA	1.87	0.57
1:A:436:ASN:O	1:A:446:ARG:NH2	2.38	0.57
1:A:494:VAL:HB	1:A:500:HIS:HB2	1.86	0.57
1:A:319:PRO:HG3	1:A:637:HIS:CD2	2.40	0.56
1:A:179:LEU:HD12	1:A:263:PRO:HG3	1.87	0.56
1:A:593:TYR:CZ	1:A:595:PRO:HG3	2.40	0.56
1:A:134:VAL:O	1:A:150:GLY:HA3	2.06	0.56
1:A:393:GLU:O	1:A:396:ARG:HG2	2.06	0.56
2:C:3:DA:H2'	2:C:4:DG:H8	1.71	0.55
1:A:331:ALA:HA	1:A:452:LEU:HD11	1.89	0.55
1:A:396:ARG:CG	1:A:397:LYS:N	2.70	0.54
1:A:588:LEU:HD11	1:A:604:LEU:HD23	1.90	0.54
1:A:599:LYS:HG3	1:A:599:LYS:O	2.06	0.54
1:A:57:THR:HG22	1:A:66:SER:OG	2.08	0.54
1:A:135:TYR:CZ	1:A:172:ARG:HG3	2.43	0.53
1:A:446:ARG:HG3	2:C:2:DG:C8	2.42	0.53
1:A:100:PRO:O	1:A:106:ARG:HD3	2.09	0.53
1:A:389:LEU:O	1:A:389:LEU:HG	2.08	0.53
1:A:605:THR:HG21	1:A:651:ARG:O	2.09	0.53
1:A:462:VAL:HG23	1:A:463:VAL:HG13	1.91	0.52
1:A:192:ARG:NH2	2:C:10:DA:O4'	2.42	0.52
1:A:177:MET:HG2	1:A:181:ALA:HB3	1.90	0.52
1:A:631:LEU:O	1:A:635:ILE:HG12	2.09	0.52
1:A:115:ARG:NH2	1:A:119:GLU:OE2	2.42	0.52
1:A:433:GLN:HG2	1:A:450:ALA:O	2.10	0.52
1:A:479:ALA:HA	1:A:488:GLY:HA2	1.91	0.52
1:A:324:ARG:HD3	1:A:372:SER:HA	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:392:ARG:O	1:A:395:LEU:HB2	2.10	0.52
1:A:347:ALA:HB3	1:A:406:VAL:HG22	1.93	0.51
1:A:513:ARG:HH22	1:A:551:GLN:CG	2.22	0.51
1:A:513:ARG:HH22	1:A:551:GLN:HG3	1.76	0.51
1:A:427:ARG:HH11	1:A:427:ARG:HG3	1.75	0.51
1:A:283:TRP:CZ3	1:A:583:PRO:HD2	2.46	0.51
1:A:315:ARG:NH2	1:A:592:LEU:HD13	2.26	0.51
1:A:324:ARG:CD	1:A:372:SER:HA	2.41	0.50
1:A:558:LEU:HD22	1:A:568:TYR:CE1	2.47	0.50
1:A:384:HIS:H	1:A:387:GLN:NE2	2.10	0.50
1:A:384:HIS:H	1:A:387:GLN:HE21	1.59	0.50
1:A:385:PRO:HA	1:A:391:PHE:CZ	2.46	0.50
1:A:615:ARG:O	1:A:651:ARG:NH1	2.43	0.50
1:A:327:VAL:HG22	1:A:332:ASP:HB2	1.94	0.49
1:A:485:PHE:HB2	1:A:509:GLN:O	2.12	0.49
1:A:514:ILE:HG13	1:A:545:ARG:HH21	1.78	0.49
1:A:445:HIS:NE2	3:G:19:C:H2'	2.27	0.49
1:A:661:ARG:HB3	1:A:684:PHE:HD1	1.78	0.49
1:A:551:GLN:O	1:A:551:GLN:HG3	2.12	0.49
1:A:24:LEU:O	1:A:26:PRO:HD3	2.11	0.48
1:A:396:ARG:HG2	1:A:397:LYS:H	1.78	0.48
1:A:385:PRO:HA	1:A:391:PHE:CE1	2.49	0.48
1:A:506:PRO:O	1:A:671:ILE:HD11	2.13	0.48
1:A:78:THR:HA	1:A:87:ALA:HA	1.95	0.48
1:A:6:LYS:NZ	1:A:314:TYR:HE2	2.11	0.48
1:A:513:ARG:NH2	1:A:551:GLN:HG3	2.29	0.48
1:A:597:GLU:H	1:A:600:THR:HG23	1.79	0.48
1:A:427:ARG:HH11	1:A:427:ARG:CG	2.27	0.47
1:A:242:ALA:HB2	1:A:258:THR:HG22	1.96	0.47
1:A:396:ARG:O	1:A:399:LYS:HG3	2.14	0.47
1:A:143:PRO:HD2	1:A:273:GLU:OE2	2.14	0.47
1:A:357:TRP:CD2	1:A:378:LEU:HD23	2.50	0.47
1:A:357:TRP:CE2	1:A:378:LEU:HD23	2.49	0.47
1:A:335:ARG:NH2	1:A:448:GLU:OE1	2.46	0.47
1:A:502:LEU:HG	1:A:679:ARG:O	2.14	0.47
1:A:502:LEU:HD11	1:A:680:GLU:HB3	1.96	0.47
1:A:237:GLU:O	1:A:259:GLY:HA3	2.15	0.47
1:A:70:PRO:HA	1:A:73:LEU:HD12	1.97	0.47
1:A:393:GLU:HB3	1:A:396:ARG:NE	2.29	0.46
1:A:531:ARG:HD3	1:A:537:LEU:HA	1.98	0.46
1:A:540:ARG:HD3	1:A:567:ALA:CB	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:410:THR:O	1:A:436:ASN:HA	2.16	0.46
1:A:369:PHE:CD1	1:A:376:LEU:HD22	2.51	0.46
1:A:480:GLY:CA	1:A:487:PHE:H	2.27	0.46
1:A:28:ARG:HH12	1:A:63:GLY:HA3	1.81	0.46
1:A:60:MET:HE2	1:A:65:ALA:HB2	1.97	0.46
1:A:390:ALA:O	1:A:393:GLU:OE1	2.34	0.45
1:A:472:GLU:OE1	1:A:536:ARG:NH2	2.47	0.45
1:A:127:VAL:HG23	1:A:129:VAL:HG23	1.99	0.45
1:A:222:SER:HB2	1:A:225:ASP:HB2	1.99	0.45
1:A:444:ARG:HA	1:A:447:TRP:NE1	2.31	0.45
1:A:592:LEU:HD12	1:A:592:LEU:HA	1.86	0.44
1:A:451:LEU:HA	1:A:451:LEU:HD23	1.83	0.44
1:A:31:VAL:HG21	1:A:64:LEU:HD13	1.97	0.44
1:A:413:MET:O	3:G:14:A:N6	2.47	0.44
1:A:426:LEU:HA	1:A:426:LEU:HD12	1.42	0.44
1:A:480:GLY:H	1:A:488:GLY:H	1.66	0.44
1:A:500:HIS:CE1	1:A:533:LYS:HD3	2.53	0.44
1:A:254:ILE:HA	1:A:255:PRO:HD3	1.84	0.44
1:A:409:LEU:HA	1:A:409:LEU:HD23	1.75	0.44
2:C:6:DT:H2"	2:C:7:DA:H5"	1.99	0.44
1:A:20:ASN:HB2	1:A:21:PRO:HD2	1.99	0.43
1:A:359:GLU:N	1:A:359:GLU:OE1	2.45	0.43
1:A:479:ALA:HB2	1:A:545:ARG:HH22	1.82	0.43
1:A:319:PRO:HG2	1:A:640:ARG:HG3	1.98	0.43
1:A:480:GLY:HA3	1:A:486:ARG:HD2	2.00	0.43
1:A:427:ARG:HG2	1:A:673:HIS:HD2	1.82	0.43
1:A:472:GLU:OE2	1:A:539:SER:OG	2.34	0.43
1:A:25:ARG:HA	1:A:97:PRO:HA	2.01	0.43
1:A:128:TRP:CZ3	1:A:174:LEU:HD21	2.53	0.43
1:A:189:LEU:HA	1:A:190:PRO:HD3	1.84	0.43
1:A:286:ARG:HD2	1:A:613:THR:HG21	2.00	0.43
1:A:642:TYR:HA	1:A:643:PRO:HD3	1.82	0.43
1:A:77:GLY:HA3	1:A:88:TYR:CZ	2.53	0.43
1:A:583:PRO:CG	1:A:588:LEU:HB2	2.49	0.43
1:A:533:LYS:HE2	1:A:533:LYS:HB3	1.44	0.42
1:A:531:ARG:CD	1:A:537:LEU:HD23	2.48	0.42
1:A:480:GLY:HA3	1:A:486:ARG:HB3	2.01	0.42
1:A:285:GLU:CD	1:A:288:ARG:HH21	2.21	0.42
1:A:359:GLU:HG2	1:A:360:PHE:N	2.34	0.42
1:A:523:LEU:HD21	1:A:561:LEU:HD11	2.01	0.42
1:A:265:LEU:HD22	1:A:269:ASP:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:492:CYS:SG	1:A:526:THR:HB	2.60	0.42
1:A:552:ASP:HB2	1:A:555:ALA:HB2	2.02	0.42
1:A:120:ARG:HA	1:A:120:ARG:HD3	1.76	0.42
1:A:183:LEU:HD13	1:A:207:LEU:HD11	2.02	0.42
1:A:215:LEU:HD23	1:A:223:LEU:HD22	2.01	0.42
1:A:404:GLN:O	1:A:681:LYS:HD3	2.20	0.41
1:A:480:GLY:HA2	1:A:487:PHE:N	2.28	0.41
1:A:593:TYR:OH	1:A:628:LEU:HB3	2.20	0.41
1:A:60:MET:HG3	1:A:108:VAL:HG21	2.02	0.41
1:A:135:TYR:CE2	1:A:172:ARG:HG3	2.55	0.41
1:A:297:ILE:O	1:A:301:LEU:HB2	2.20	0.41
2:C:2:DG:H2''	2:C:3:DA:H8	1.85	0.41
1:A:89:ARG:HE	1:A:91:TYR:HE1	1.60	0.41
2:C:16:DT:H2'	2:C:17:DA:C8	2.56	0.41
1:A:437:VAL:O	1:A:439:LEU:N	2.53	0.41
1:A:575:LYS:HB3	1:A:651:ARG:NH2	2.36	0.41
1:A:587:ARG:HD2	1:A:587:ARG:HH11	1.70	0.41
1:A:217:LEU:HA	1:A:217:LEU:HD12	1.87	0.40
1:A:540:ARG:HD3	1:A:567:ALA:HB3	2.03	0.40
1:A:171:TYR:O	2:C:8:DG:H5'	2.21	0.40
1:A:492:CYS:O	1:A:501:LEU:HA	2.20	0.40
1:A:104:GLY:O	1:A:108:VAL:HG23	2.20	0.40
1:A:389:LEU:O	1:A:389:LEU:CG	2.67	0.40
1:A:415:TRP:HE3	2:C:1:DT:O4	2.05	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	682/685 (100%)	637 (93%)	39 (6%)	6 (1%)	17 48

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	218	PRO
1	A	610	PHE
1	A	82	MET
1	A	92	PRO
1	A	488	GLY
1	A	608	ARG

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	502/549 (91%)	478 (95%)	24 (5%)	25 58

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

Mol	Chain	Res	Type
1	A	28	ARG
1	A	59	ARG
1	A	79	LEU
1	A	136	ARG
1	A	209	GLU
1	A	327	VAL
1	A	343	GLU
1	A	389	LEU
1	A	391	PHE
1	A	396	ARG
1	A	426	LEU
1	A	427	ARG
1	A	440	ARG
1	A	462	VAL
1	A	483	GLU
1	A	502	LEU
1	A	517	GLU
1	A	525	GLU
1	A	587	ARG

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Mol	Chain	Res	Type
1	A	605	THR
1	A	608	ARG
1	A	671	ILE
1	A	675	LYS
1	A	680	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	387	GLN
1	A	673	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	G	16/20 (80%)	1 (6%)	0

All (1) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	G	4	A

There are no RNA pucker outliers to report.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 1 ligands modelled in this entry, 1 is 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

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 ⓘ

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, 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	684/685 (99%)	0.13	14 (2%) 65 63	40, 75, 121, 198	0
2	C	21/21 (100%)	-0.35	0 100 100	55, 64, 98, 118	0
3	G	18/20 (90%)	-0.07	1 (5%) 24 20	55, 73, 108, 153	1 (5%)
All	All	723/726 (99%)	0.11	15 (2%) 63 61	40, 75, 121, 198	1 (0%)

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	388	GLY	4.9
1	A	391	PHE	3.9
1	A	348	LEU	2.8
3	G	3	U	2.8
1	A	91	TYR	2.7
1	A	357	TRP	2.6
1	A	377	ARG	2.5
1	A	409	LEU	2.4
1	A	387	GLN	2.3
1	A	74	VAL	2.3
1	A	347	ALA	2.3
1	A	400	GLU	2.2
1	A	32	VAL	2.2
1	A	86	TYR	2.1
1	A	204	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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](#)

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, 95th 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(\AA^2)	Q<0.9
4	MG	C	101	1/1	0.96	0.25	61,61,61,61	0

6.5 Other polymers [i](#)

There are no such residues in this entry.