



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 07:05 PM UTC

PDB ID : 3HVR / pdb_00003hvr
Title : Crystal structure of T. thermophilus Argonaute complexed with DNA guide strand and 19-nt RNA target strand with two Mg²⁺ at the cleavage site
Authors : Wang, Y.; Li, H.; Sheng, G.; Patel, D.J.
Deposited on : 2009-06-16
Resolution : 3.21 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

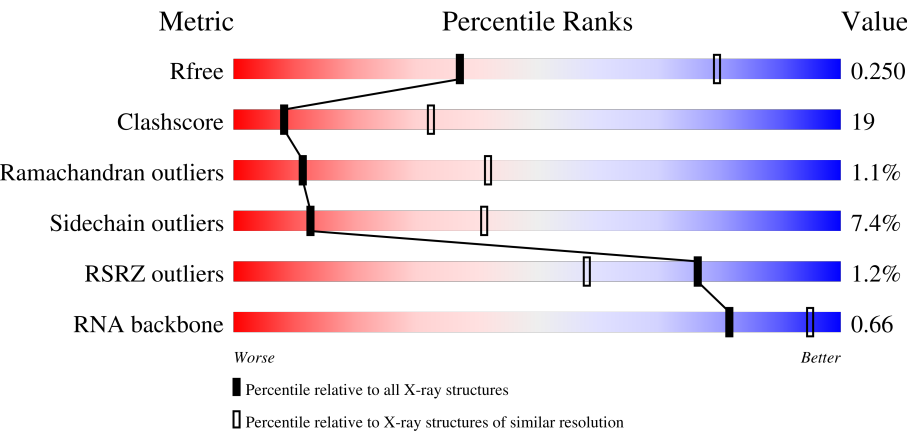
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.21 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.



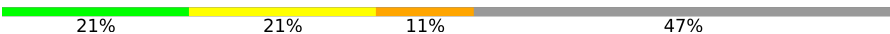
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1768 (3.24-3.20)
Clashscore	190562	1879 (3.24-3.20)
Ramachandran outliers	187476	1844 (3.24-3.20)
Sidechain outliers	187428	1843 (3.24-3.20)
RSRZ outliers	180081	1768 (3.24-3.20)
RNA backbone	3983	1055 (3.50-2.94)

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	685	% 60%31%5%.
1	B	685	% 61%32%. .
2	C	21	33%43%24%
2	M	21	33%38%5%24%

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Mol	Chain	Length	Quality of chain
3	D	19	
3	N	19	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	PO4	A	689	-	X	-	-
5	PO4	B	689	-	X	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10637 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	660	Total	C	N	O	S	0	0	0
			4754	3030	889	830	5			
1	B	657	Total	C	N	O	S	0	0	0
			4759	3038	886	830	5			

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

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	16	Total	C	N	O	P	0	0	0
			338	160	62	100	16			
2	M	16	Total	C	N	O	P	0	0	0
			338	160	62	100	16			

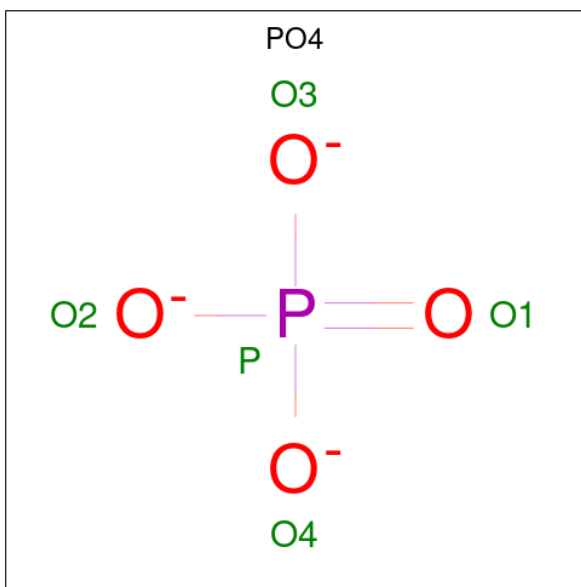
- Molecule 3 is a RNA chain called 5'-R(*UP*AP*UP*AP*CP*AP*AP*CP*CP*UP*AP*CP*UP*AP*CP*CP*UP*CP*G)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	12	Total	C	N	O	P	0	0	0
			228	101	34	81	12			
3	N	10	Total	C	N	O	P	0	0	0
			204	92	31	71	10			

- Molecule 4 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Mg	0	0
			3	3		
4	B	3	Total	Mg	0	0
			3	3		

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).

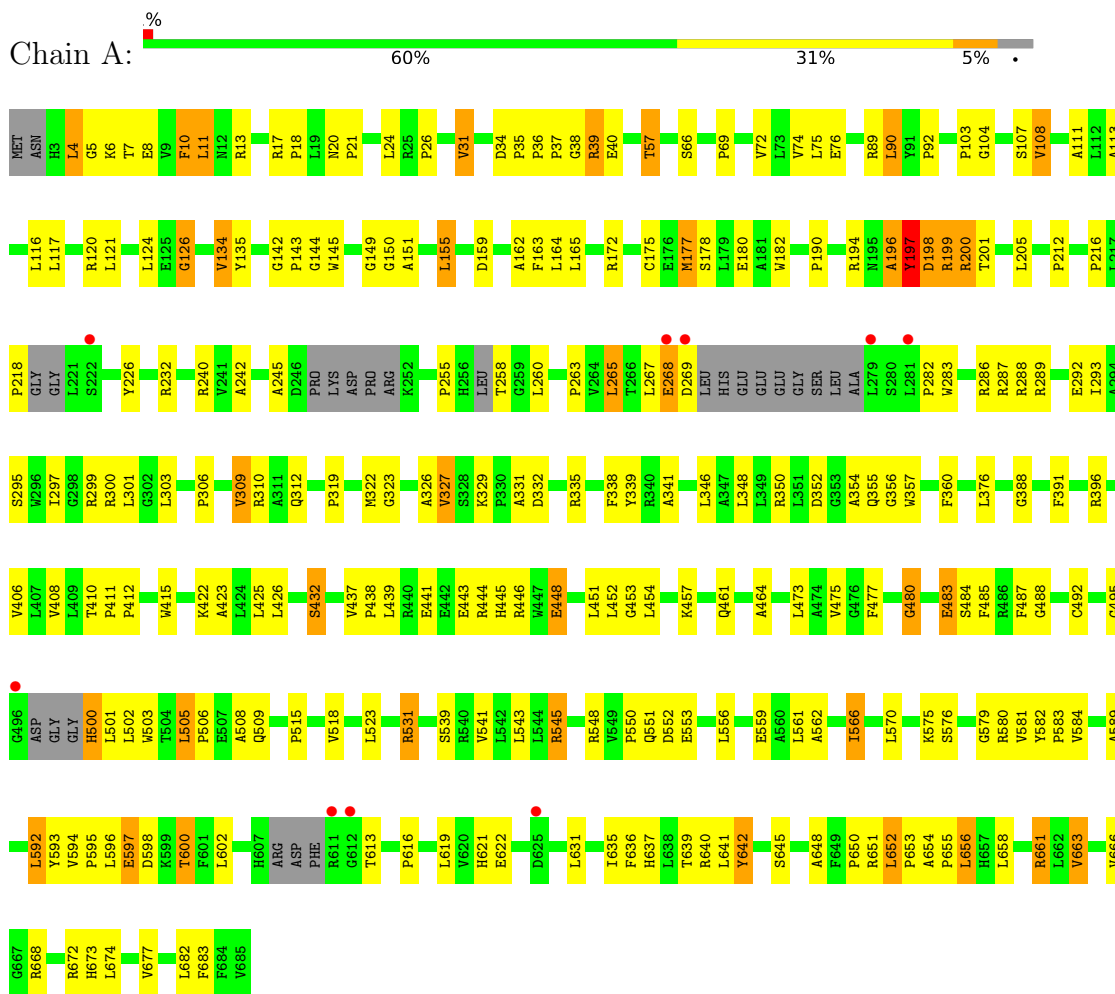


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	P	0	0
			5	4	1		
5	B	1	Total	O	P	0	0
			5	4	1		

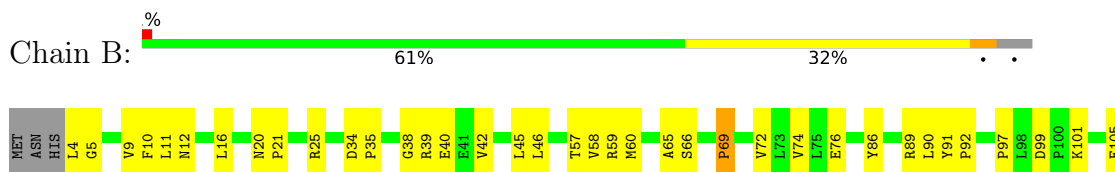
3 Residue-property plots

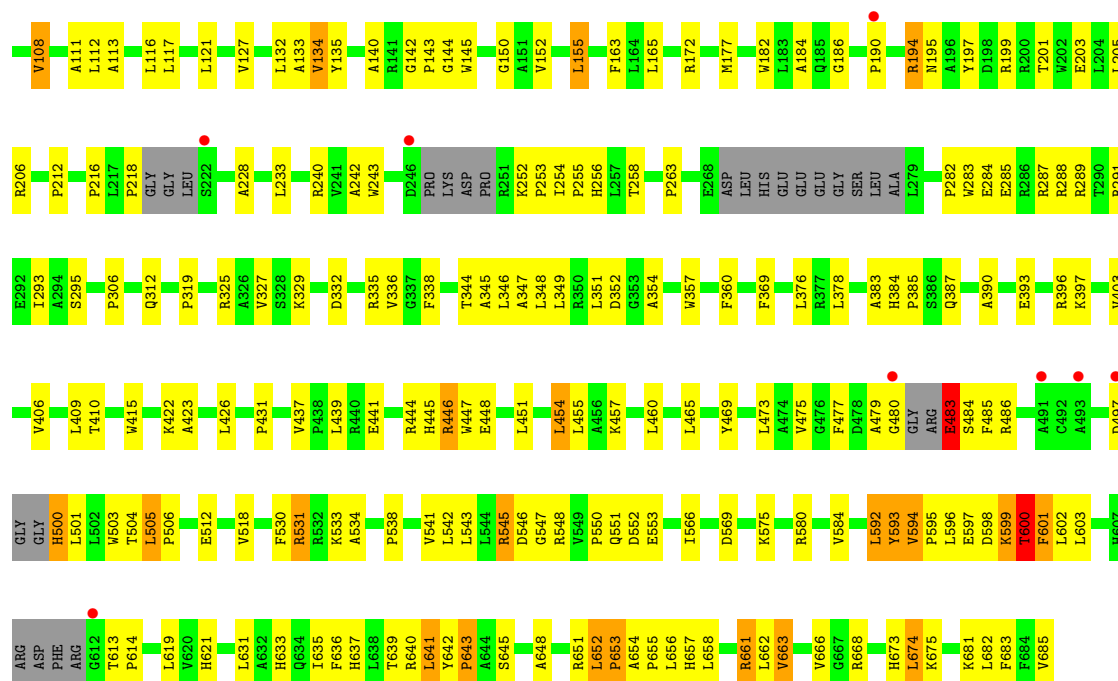
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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



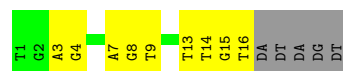
• Molecule 1: Argonaute





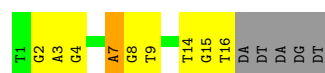
• Molecule 2: 5'-D(P*TP*GP*AP*GP*GP*TP*AP*GP*TP*AP*GP*GP*TP*TP*TP*GP*AP*TP*AP*GP*T)-3'

Chain C: 33% 43% 24%



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

Chain M: 33% 38% 5% 24%



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

Chain D: 21% 42% 37%



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

Chain N: 21% 21% 11% 47%

U	A	U	A	C	A	A	C	C9	U10	A11	A14	C15	C18	G
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	111.25Å 116.90Å 170.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.68 – 3.21 37.68 – 3.21	Depositor EDS
% Data completeness (in resolution range)	94.4 (37.68-3.21) 94.3 (37.68-3.21)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.59 (at 3.18Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.226 , 0.273 (Not available) , 0.250	Depositor DCC
R_{free} test set	1744 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	100.3	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 99.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.046 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10637	wwPDB-VP
Average B, all atoms (Å ²)	97.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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, PO4

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	0.47	4/4859 (0.1%)	0.91	22/6624 (0.3%)
1	B	0.56	7/4866 (0.1%)	0.92	25/6635 (0.4%)
2	C	0.31	0/379	0.71	0/584
2	M	0.32	0/379	0.69	1/584 (0.2%)
3	D	0.44	0/251	0.71	0/387
3	N	0.54	0/225	0.93	2/346 (0.6%)
All	All	0.51	11/10959 (0.1%)	0.90	50/15160 (0.3%)

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
1	B	0	1
All	All	0	5

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	199	ARG	C-N	11.23	1.48	1.33
1	B	593	TYR	C-N	10.91	1.44	1.33
1	B	600	THR	C-N	-9.80	1.20	1.33
1	B	599	LYS	C-N	9.00	1.46	1.33
1	B	594	VAL	CA-C	-7.37	1.46	1.52
1	B	601	PHE	C-N	-7.28	1.22	1.33
1	A	196	ALA	C-N	-7.09	1.22	1.33
1	B	594	VAL	CA-CB	-6.98	1.48	1.54
1	A	197	TYR	C-N	-5.78	1.25	1.33
1	A	199	ARG	CA-C	-5.53	1.46	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	194	ARG	C-N	5.17	1.40	1.33

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	483	GLU	O-C-N	-10.52	106.17	123.00
1	B	201	THR	CA-C-N	-10.39	105.73	122.36
1	B	201	THR	C-N-CA	-10.39	105.73	122.36
1	B	201	THR	O-C-N	9.56	135.92	123.11
1	A	92	PRO	N-CA-CB	8.94	110.45	103.30
1	A	200	ARG	O-C-N	-8.88	112.39	122.86
1	A	199	ARG	N-CA-C	-8.68	101.31	114.16
1	B	92	PRO	N-CA-CB	8.57	110.99	103.27
1	B	216	PRO	N-CA-CB	8.27	110.22	103.36
1	A	197	TYR	O-C-N	-8.23	111.19	122.46
1	B	263	PRO	N-CA-CB	8.13	110.42	103.35
1	A	198	ASP	O-C-N	-8.09	111.84	122.59
1	A	255	PRO	N-CA-CB	7.99	110.31	103.35
1	B	282	PRO	N-CA-CB	7.84	110.29	103.31
1	A	282	PRO	N-CA-CB	7.32	109.69	103.25
1	A	642	TYR	CA-C-N	7.06	128.66	119.84
1	A	642	TYR	C-N-CA	7.06	128.66	119.84
1	A	263	PRO	N-CA-CB	7.05	110.20	102.72
1	B	594	VAL	CB-CA-C	-7.02	102.72	110.52
3	N	9	C	P-O3'-C3'	7.01	130.72	120.20
1	B	143	PRO	N-CA-CB	6.98	110.58	103.25
1	A	218	PRO	N-CA-CB	6.95	110.64	103.00
1	B	69	PRO	CA-C-N	6.89	127.46	119.47
1	B	69	PRO	C-N-CA	6.89	127.46	119.47
1	A	143	PRO	N-CA-CB	6.87	110.46	103.25
1	B	218	PRO	N-CA-CB	6.78	110.46	103.00
1	A	216	PRO	N-CA-CB	6.65	110.23	103.25
1	A	212	PRO	N-CA-CB	6.62	110.20	103.25
1	A	103	PRO	N-CA-CB	6.60	110.29	103.23
3	N	9	C	C2'-C3'-O3'	6.49	123.43	113.70
1	B	212	PRO	N-CA-CB	6.45	110.36	103.52
1	B	469	TYR	CA-C-N	6.42	125.65	118.97
1	B	469	TYR	C-N-CA	6.42	125.65	118.97
1	A	151	ALA	N-CA-C	-6.26	99.70	109.72
1	A	265	LEU	O-C-N	-6.21	115.32	122.89
1	B	599	LYS	O-C-N	-5.99	115.87	121.56
1	A	196	ALA	O-C-N	-5.78	114.20	122.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	199	ARG	O-C-N	5.72	128.95	122.20
1	B	652	LEU	CA-C-N	5.69	126.95	119.84
1	B	652	LEU	C-N-CA	5.69	126.95	119.84
2	M	7	DA	P-O3'-C3'	-5.41	112.09	120.20
1	B	199	ARG	CA-C-N	-5.34	112.56	120.94
1	B	199	ARG	C-N-CA	-5.34	112.56	120.94
1	A	267	LEU	N-CA-C	5.26	117.92	111.82
1	A	268	GLU	N-CA-C	5.20	117.53	110.35
1	B	593	TYR	O-C-N	-5.18	117.21	123.27
1	B	592	LEU	CA-C-N	-5.12	115.55	122.77
1	B	592	LEU	C-N-CA	-5.12	115.55	122.77
1	A	39	ARG	N-CA-C	-5.11	106.15	112.38
1	A	483	GLU	N-CA-C	5.05	116.84	110.33

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	149	GLY	Peptide
1	A	197	TYR	Mainchain
1	A	200	ARG	Mainchain
1	A	265	LEU	Mainchain
1	B	483	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	4754	0	4408	184	0
1	B	4759	0	4449	183	0
2	C	338	0	183	6	0
2	M	338	0	183	8	0
3	D	228	0	118	8	0
3	N	204	0	108	9	0
4	A	3	0	0	0	0
4	B	3	0	0	0	0
5	A	5	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	5	0	0	1	0
All	All	10637	0	9449	386	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 19.

All (386) 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:480:GLY:HA2	1:B:663:VAL:HG13	1.37	1.01
1:A:480:GLY:HA2	1:A:663:VAL:HG13	1.48	0.96
1:B:465:LEU:O	1:B:497:ASP:CB	2.14	0.95
1:A:473:LEU:HB3	1:A:541:VAL:HG12	1.50	0.94
1:B:661:ARG:CG	1:B:661:ARG:HH21	1.81	0.93
1:A:661:ARG:HG2	1:A:661:ARG:HH21	1.35	0.91
1:A:596:LEU:O	1:A:597:GLU:O	1.89	0.90
1:B:465:LEU:HB2	1:B:497:ASP:CB	2.03	0.89
1:B:642:TYR:HD2	1:B:645:SER:HB3	1.38	0.88
1:B:661:ARG:HH21	1:B:661:ARG:HG2	1.40	0.85
1:A:197:TYR:CE1	1:A:232:ARG:NH1	2.46	0.84
1:A:483:GLU:O	1:A:485:PHE:N	2.12	0.82
1:B:473:LEU:HB3	1:B:541:VAL:HG12	1.62	0.81
1:B:598:ASP:O	1:B:599:LYS:HB2	1.80	0.80
1:A:197:TYR:CE1	1:A:232:ARG:CZ	2.65	0.80
1:A:120:ARG:HB3	1:A:301:LEU:HD11	1.62	0.79
1:A:575:LYS:HB3	1:A:651:ARG:HH22	1.47	0.79
1:A:7:THR:HG22	1:A:8:GLU:H	1.45	0.78
1:A:661:ARG:HH21	1:A:661:ARG:CG	1.97	0.78
1:A:453:GLY:O	1:A:457:LYS:HG3	1.86	0.76
1:A:593:TYR:CZ	1:A:595:PRO:HG3	2.20	0.75
1:A:197:TYR:CG	1:A:232:ARG:HD3	2.22	0.75
1:B:465:LEU:H	1:B:497:ASP:CB	2.00	0.75
1:B:117:LEU:HD22	1:B:155:LEU:HB2	1.68	0.74
1:B:242:ALA:HB2	1:B:258:THR:HG22	1.69	0.73
1:B:352:ASP:HB3	1:B:437:VAL:HG21	1.69	0.73
1:A:121:LEU:HD13	1:A:134:VAL:HG21	1.69	0.73
1:A:319:PRO:HA	1:A:637:HIS:CE1	2.24	0.73
1:A:319:PRO:HG2	1:A:640:ARG:HD3	1.70	0.72
1:B:597:GLU:O	1:B:600:THR:HG23	1.89	0.72
3:D:9:C:H2'	3:D:10:U:H6	1.54	0.71
1:B:319:PRO:HG2	1:B:640:ARG:HD3	1.71	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:182:TRP:HZ3	1:B:190:PRO:HD3	1.56	0.70
1:B:287:ARG:O	1:B:291:ARG:HG3	1.92	0.70
1:B:666:VAL:HG22	1:B:674:LEU:HD11	1.72	0.70
1:B:12:ASN:HD22	1:B:580:ARG:HB2	1.57	0.69
1:B:483:GLU:O	1:B:484:SER:C	2.33	0.69
1:A:197:TYR:CZ	1:A:232:ARG:NH1	2.60	0.69
1:B:58:VAL:HG11	1:B:111:ALA:HB1	1.75	0.69
1:B:444:ARG:O	1:B:448:GLU:HB2	1.92	0.69
1:B:58:VAL:HG23	1:B:112:LEU:HD23	1.75	0.69
1:B:465:LEU:CB	1:B:497:ASP:CB	2.71	0.69
1:A:113:ALA:HB1	1:A:155:LEU:HD13	1.75	0.68
1:A:654:ALA:HB3	1:A:655:PRO:HD3	1.74	0.68
1:A:312:GLN:CD	1:A:312:GLN:H	2.02	0.67
1:B:635:ILE:HG23	1:B:653:PRO:HG3	1.77	0.67
1:B:423:ALA:HB1	1:B:673:HIS:CE1	2.30	0.67
1:A:289:ARG:O	1:A:293:ILE:HG12	1.95	0.67
1:B:346:LEU:HD23	1:B:454:LEU:HD13	1.78	0.66
1:A:642:TYR:HD2	1:A:645:SER:HB3	1.61	0.66
1:B:182:TRP:CZ3	1:B:190:PRO:HG3	2.32	0.65
2:M:8:DG:H3'	2:M:9:DT:H71	1.78	0.65
1:A:240:ARG:O	1:A:258:THR:HG23	1.97	0.65
1:B:480:GLY:HA2	1:B:663:VAL:CG1	2.20	0.65
1:B:422:LYS:O	1:B:426:LEU:HD13	1.97	0.64
1:B:465:LEU:CA	1:B:497:ASP:CB	2.75	0.64
2:M:15:DG:H2''	2:M:16:DT:H72	1.79	0.64
1:A:415:TRP:HZ3	1:A:668:ARG:HD2	1.61	0.64
1:A:531:ARG:HD3	1:A:531:ARG:C	2.23	0.64
1:B:661:ARG:CG	1:B:661:ARG:NH2	2.51	0.64
1:B:38:GLY:C	1:B:40:GLU:H	2.05	0.64
1:A:575:LYS:HB3	1:A:651:ARG:NH2	2.12	0.63
1:A:445:HIS:NE2	3:D:18:C:H2'	2.13	0.63
1:A:635:ILE:HG23	1:A:653:PRO:HG3	1.80	0.63
1:B:465:LEU:N	1:B:497:ASP:CB	2.62	0.62
1:B:652:LEU:HB2	1:B:653:PRO:HD2	1.80	0.62
1:B:542:LEU:C	1:B:543:LEU:HD12	2.24	0.62
1:B:633:HIS:O	1:B:637:HIS:HD2	1.82	0.62
1:A:197:TYR:CD2	1:A:232:ARG:HD3	2.35	0.62
1:B:360:PHE:CG	1:B:441:GLU:HG2	2.35	0.62
1:A:197:TYR:CD1	1:A:232:ARG:CZ	2.83	0.61
1:B:596:LEU:O	1:B:597:GLU:C	2.41	0.61
1:A:672:ARG:HA	1:B:505:LEU:HD12	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:621:HIS:ND1	1:A:631:LEU:HD11	2.15	0.61
1:A:39:ARG:NH1	5:B:689:PO4:O3	2.33	0.61
1:B:329:LYS:HG2	1:B:332:ASP:OD2	2.02	0.60
1:A:155:LEU:HD21	1:A:163:PHE:HB3	1.83	0.60
1:B:57:THR:HG22	1:B:66:SER:CB	2.32	0.60
1:A:350:ARG:HD3	1:A:354:ALA:O	2.02	0.59
1:B:601:PHE:HD2	1:B:602:LEU:O	1.86	0.59
1:A:543:LEU:N	1:A:543:LEU:HD12	2.17	0.59
1:B:445:HIS:NE2	3:N:18:C:H2'	2.17	0.59
1:B:465:LEU:C	1:B:497:ASP:CB	2.76	0.58
1:A:75:LEU:HA	1:A:90:LEU:HD23	1.85	0.58
3:N:9:C:O2'	3:N:10:U:H5'	2.04	0.58
1:A:242:ALA:HB2	1:A:258:THR:HG22	1.86	0.58
1:A:322:MET:HG3	1:A:326:ALA:HA	1.85	0.58
1:B:45:LEU:HD11	1:B:86:TYR:CD2	2.37	0.58
1:B:34:ASP:HA	1:B:35:PRO:C	2.29	0.58
1:A:619:LEU:N	1:A:619:LEU:HD12	2.19	0.57
1:B:135:TYR:HA	1:B:150:GLY:HA3	1.87	0.57
1:B:500:HIS:NE2	1:B:533:LYS:HG3	2.19	0.57
1:B:12:ASN:ND2	1:B:580:ARG:HB2	2.19	0.57
3:D:9:C:O2'	3:D:10:U:H5'	2.05	0.57
1:A:396:ARG:NH1	1:B:531:ARG:HH12	2.03	0.57
1:A:10:PHE:HB3	1:A:310:ARG:HA	1.87	0.56
1:A:13:ARG:HB2	1:A:309:VAL:CG1	2.35	0.56
1:A:24:LEU:C	1:A:26:PRO:HD3	2.30	0.56
1:A:17:ARG:HE	1:A:303:LEU:HD23	1.70	0.56
1:B:661:ARG:HH21	1:B:661:ARG:HG3	1.65	0.56
1:B:661:ARG:N	1:B:661:ARG:HD2	2.19	0.56
1:B:360:PHE:HD1	1:B:360:PHE:H	1.53	0.56
1:A:322:MET:HG3	1:A:326:ALA:N	2.20	0.56
1:B:348:LEU:HD11	1:B:409:LEU:HD12	1.88	0.55
1:B:501:LEU:HD21	1:B:641:LEU:HD13	1.87	0.55
1:A:437:VAL:O	1:A:439:LEU:N	2.39	0.55
1:A:483:GLU:O	1:A:484:SER:C	2.48	0.55
1:A:144:GLY:HA2	1:A:177:MET:HE2	1.88	0.55
1:A:661:ARG:CG	1:A:661:ARG:NH2	2.65	0.55
1:A:666:VAL:HG22	1:A:674:LEU:HD11	1.89	0.55
1:A:339:TYR:CE1	1:A:464:ALA:HB3	2.42	0.55
1:B:338:PHE:CZ	1:B:455:LEU:HD13	2.41	0.55
1:A:589:ALA:HB3	1:A:592:LEU:HD11	1.89	0.55
1:B:12:ASN:C	1:B:12:ASN:OD1	2.49	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:LEU:HD22	1:A:4:LEU:O	2.06	0.55
1:A:598:ASP:O	1:A:600:THR:HG22	2.07	0.55
1:B:415:TRP:HZ3	1:B:668:ARG:HD2	1.71	0.55
1:B:530:PHE:CD2	1:B:538:PRO:HG3	2.42	0.54
2:M:3:DA:H2'	2:M:4:DG:C8	2.42	0.54
2:M:14:DT:C4	2:M:15:DG:C2	2.95	0.54
1:A:18:PRO:HA	1:A:162:ALA:HA	1.89	0.54
1:B:89:ARG:HG3	1:B:91:TYR:CE1	2.42	0.54
1:A:645:SER:HB2	1:A:648:ALA:O	2.07	0.54
1:A:422:LYS:O	1:A:426:LEU:HD13	2.07	0.54
1:A:487:PHE:CE1	1:A:508:ALA:HB2	2.43	0.54
1:B:530:PHE:CG	1:B:538:PRO:HG3	2.43	0.54
1:B:551:GLN:O	1:B:552:ASP:HB2	2.08	0.54
1:A:501:LEU:HD13	1:A:658:LEU:HD11	1.90	0.54
1:A:652:LEU:HB2	1:A:653:PRO:HD2	1.90	0.53
1:A:332:ASP:HA	1:A:335:ARG:HD3	1.89	0.53
1:B:319:PRO:HA	1:B:637:HIS:CE1	2.43	0.53
2:C:13:DT:H73	2:C:14:DT:C4	2.43	0.53
1:A:268:GLU:O	1:A:269:ASP:CB	2.57	0.53
1:A:503:TRP:HZ2	1:A:677:VAL:HG12	1.73	0.53
1:A:505:LEU:HG	1:B:675:LYS:HB2	1.91	0.53
1:B:295:SER:HA	1:B:306:PRO:HG2	1.90	0.53
1:A:539:SER:O	1:A:566:ILE:HD12	2.09	0.53
1:A:621:HIS:CE1	1:A:631:LEU:HD11	2.44	0.53
1:B:640:ARG:HD2	1:B:640:ARG:N	2.24	0.53
1:B:501:LEU:HD13	1:B:658:LEU:HD11	1.91	0.53
1:A:322:MET:HG3	1:A:326:ALA:CA	2.38	0.53
1:A:20:ASN:HB2	1:A:21:PRO:HD2	1.89	0.52
1:A:327:VAL:CG2	1:A:332:ASP:HB2	2.39	0.52
1:B:446:ARG:HD2	2:M:2:DG:C8	2.44	0.52
1:B:38:GLY:C	1:B:40:GLU:N	2.68	0.52
1:A:10:PHE:HA	1:A:310:ARG:HA	1.91	0.52
1:B:483:GLU:O	1:B:485:PHE:N	2.42	0.52
1:B:437:VAL:O	1:B:439:LEU:N	2.41	0.52
1:B:594:VAL:O	1:B:601:PHE:HB2	2.10	0.52
1:B:357:TRP:CZ2	1:B:378:LEU:HD12	2.44	0.52
1:A:7:THR:HG22	1:A:8:GLU:N	2.20	0.52
1:B:657:HIS:O	1:B:661:ARG:HD3	2.09	0.52
1:B:57:THR:HG22	1:B:66:SER:OG	2.10	0.52
1:A:545:ARG:NH1	1:A:553:GLU:OE1	2.42	0.52
1:B:480:GLY:CA	1:B:663:VAL:HG13	2.26	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:PHE:CG	1:A:441:GLU:HG2	2.45	0.51
2:C:15:DG:H2'	2:C:16:DT:OP2	2.10	0.51
1:A:483:GLU:C	1:A:485:PHE:H	2.17	0.51
1:B:58:VAL:HG12	1:B:59:ARG:H	1.73	0.51
3:D:14:A:H2'	3:D:15:C:C6	2.46	0.51
1:B:390:ALA:O	1:B:393:GLU:HB3	2.10	0.51
1:A:8:GLU:O	1:A:584:VAL:HG23	2.11	0.51
1:A:355:GLN:HG2	1:A:356:GLY:N	2.26	0.51
1:A:360:PHE:CD1	1:A:441:GLU:HG2	2.46	0.51
1:B:140:ALA:HB1	1:B:145:TRP:HZ3	1.75	0.51
1:B:543:LEU:HD12	1:B:543:LEU:N	2.25	0.51
1:A:596:LEU:O	1:A:597:GLU:C	2.54	0.51
1:A:443:GLU:O	1:A:446:ARG:HB2	2.11	0.51
1:A:596:LEU:C	1:A:597:GLU:O	2.53	0.51
1:B:69:PRO:O	1:B:72:VAL:HG22	2.11	0.51
2:C:7:DA:H2'	2:C:8:DG:H8	1.75	0.51
1:A:523:LEU:HD21	1:A:561:LEU:HD11	1.93	0.50
1:A:121:LEU:HD13	1:A:134:VAL:CG2	2.39	0.50
1:B:12:ASN:OD1	1:B:12:ASN:O	2.29	0.50
1:B:16:LEU:HB2	1:B:163:PHE:HB2	1.94	0.50
1:B:383:ALA:HA	1:B:387:GLN:OE1	2.12	0.50
1:B:545:ARG:HD3	1:B:545:ARG:C	2.37	0.50
3:N:9:C:H2'	3:N:10:U:H6	1.77	0.50
1:A:144:GLY:O	1:A:175:CYS:HB2	2.12	0.50
1:A:286:ARG:HD2	1:A:613:THR:HG21	1.93	0.50
1:A:652:LEU:HB2	1:A:653:PRO:CD	2.41	0.49
1:B:58:VAL:HG12	1:B:59:ARG:N	2.27	0.49
1:A:425:LEU:HD12	1:A:432:SER:HB3	1.94	0.49
1:B:105:GLU:O	1:B:108:VAL:HG13	2.11	0.49
1:B:182:TRP:CZ3	1:B:190:PRO:CG	2.95	0.49
1:B:682:LEU:O	1:B:685:VAL:HG22	2.13	0.49
1:B:197:TYR:OH	1:B:256:HIS:CE1	2.66	0.49
1:B:593:TYR:CZ	1:B:595:PRO:HG3	2.48	0.49
1:A:475:VAL:HG23	1:A:477:PHE:CE1	2.48	0.49
1:B:116:LEU:C	1:B:116:LEU:HD23	2.37	0.49
1:B:142:GLY:HA3	1:B:145:TRP:CZ2	2.48	0.48
1:A:113:ALA:HB1	1:A:155:LEU:CD1	2.42	0.48
1:A:640:ARG:HD2	1:A:640:ARG:N	2.28	0.48
1:B:285:GLU:HG2	1:B:288:ARG:NH2	2.29	0.48
1:B:345:ALA:C	1:B:346:LEU:HD12	2.37	0.48
1:A:556:LEU:O	1:A:556:LEU:HD23	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:636:PHE:O	1:A:639:THR:HB	2.13	0.48
1:A:74:VAL:HG12	1:A:76:GLU:O	2.13	0.48
1:B:12:ASN:ND2	1:B:580:ARG:H	2.11	0.48
1:B:172:ARG:HD2	2:M:9:DT:OP1	2.13	0.48
1:B:312:GLN:H	1:B:312:GLN:CD	2.21	0.48
1:B:545:ARG:NH1	1:B:553:GLU:OE1	2.46	0.48
1:B:642:TYR:CD2	1:B:645:SER:HB3	2.30	0.48
1:A:135:TYR:CD2	1:A:150:GLY:HA3	2.48	0.48
1:A:327:VAL:HG23	1:A:332:ASP:HB2	1.94	0.48
1:A:339:TYR:CD1	1:A:464:ALA:HB3	2.48	0.48
1:B:25:ARG:HH11	1:B:97:PRO:HG3	1.78	0.48
1:A:4:LEU:O	1:A:6:LYS:N	2.46	0.48
1:B:289:ARG:O	1:B:293:ILE:HG12	2.13	0.48
1:B:99:ASP:OD2	1:B:101:LYS:HG2	2.13	0.48
3:N:10:U:C2	3:N:11:A:C8	3.02	0.48
1:B:661:ARG:HG2	1:B:661:ARG:NH2	2.18	0.47
1:A:594:VAL:HB	1:A:602:LEU:HB2	1.96	0.47
1:B:57:THR:HG22	1:B:66:SER:HB2	1.97	0.47
1:A:226:TYR:C	1:A:226:TYR:CD2	2.92	0.47
1:A:480:GLY:CA	1:A:663:VAL:HG13	2.31	0.47
1:B:113:ALA:HB1	1:B:155:LEU:HD13	1.95	0.47
1:B:203:GLU:O	1:B:205:LEU:HD13	2.14	0.47
1:B:548:ARG:O	1:B:550:PRO:HD3	2.13	0.47
1:B:134:VAL:O	1:B:150:GLY:HA3	2.14	0.47
1:B:531:ARG:O	1:B:534:ALA:O	2.31	0.47
1:B:661:ARG:NH2	1:B:661:ARG:HG3	2.25	0.47
3:D:12:C:H2'	3:D:13:U:C6	2.49	0.47
1:B:531:ARG:C	1:B:531:ARG:HD3	2.40	0.47
1:A:108:VAL:O	1:A:111:ALA:HB3	2.14	0.47
1:A:116:LEU:C	1:A:116:LEU:HD23	2.40	0.47
1:A:338:PHE:HB2	1:A:341:ALA:HB2	1.96	0.47
1:A:76:GLU:HG2	1:A:89:ARG:HD2	1.96	0.47
1:A:299:ARG:HH12	1:A:300:ARG:HH11	1.62	0.47
1:B:444:ARG:HA	1:B:447:TRP:NE1	2.30	0.47
1:A:182:TRP:HZ3	1:A:190:PRO:HD3	1.80	0.47
1:A:331:ALA:HA	1:A:452:LEU:HD11	1.97	0.47
1:B:121:LEU:HD13	1:B:134:VAL:HG21	1.96	0.47
3:N:9:C:C2	3:N:10:U:C6	3.03	0.47
1:A:475:VAL:HG23	1:A:477:PHE:HE1	1.79	0.47
1:B:60:MET:SD	1:B:108:VAL:HG21	2.55	0.47
2:M:2:DG:H2'	2:M:3:DA:C8	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:548:ARG:O	1:A:550:PRO:HD3	2.15	0.46
1:B:133:ALA:HA	1:B:152:VAL:HA	1.96	0.46
1:B:325:ARG:HB2	1:B:336:VAL:CG1	2.45	0.46
1:B:410:THR:N	1:B:437:VAL:HG23	2.30	0.46
1:A:475:VAL:HG12	1:A:492:CYS:HA	1.96	0.46
1:A:35:PRO:HA	1:A:36:PRO:HD3	1.85	0.46
1:B:144:GLY:HA2	1:B:177:MET:HE2	1.98	0.46
1:B:327:VAL:CG2	1:B:332:ASP:HB2	2.46	0.46
1:A:295:SER:HA	1:A:306:PRO:HG2	1.98	0.46
1:B:503:TRP:CZ2	1:B:683:PHE:HE1	2.33	0.46
1:B:45:LEU:HD11	1:B:86:TYR:CE2	2.50	0.46
1:A:329:LYS:HE2	1:A:329:LYS:HB3	1.79	0.46
1:A:350:ARG:HB3	1:A:352:ASP:OD1	2.16	0.46
1:A:36:PRO:HA	1:A:37:PRO:HD3	1.82	0.45
1:B:654:ALA:HB3	1:B:655:PRO:HD3	1.98	0.45
1:A:388:GLY:O	1:A:391:PHE:HB3	2.17	0.45
1:B:206:ARG:HG3	1:B:243:TRP:HB2	1.98	0.45
1:A:551:GLN:O	1:A:552:ASP:HB2	2.16	0.45
1:A:76:GLU:HG2	1:A:89:ARG:CD	2.47	0.45
1:A:172:ARG:HD2	2:C:9:DT:OP1	2.15	0.45
1:B:9:VAL:HA	1:B:584:VAL:HG23	1.99	0.45
1:B:415:TRP:CZ3	1:B:668:ARG:HD2	2.49	0.45
1:B:592:LEU:O	1:B:603:LEU:HD12	2.16	0.45
1:B:479:ALA:O	1:B:480:GLY:O	2.34	0.45
1:A:142:GLY:HA3	1:A:145:TRP:CZ2	2.51	0.45
1:B:38:GLY:O	1:B:40:GLU:N	2.50	0.45
1:A:69:PRO:O	1:A:72:VAL:HG22	2.17	0.45
1:A:198:ASP:CG	1:A:199:ARG:H	2.25	0.45
1:B:410:THR:CA	1:B:437:VAL:HG23	2.47	0.45
1:A:348:LEU:HB2	1:A:357:TRP:CE2	2.51	0.45
1:A:31:VAL:HG13	1:A:90:LEU:CD1	2.47	0.44
1:B:132:LEU:HD12	1:B:132:LEU:N	2.32	0.44
1:A:444:ARG:O	1:A:448:GLU:HB2	2.16	0.44
1:A:38:GLY:O	1:A:39:ARG:C	2.61	0.44
1:A:178:SER:C	1:A:180:GLU:H	2.25	0.44
1:B:140:ALA:HB1	1:B:145:TRP:CZ3	2.52	0.44
1:B:360:PHE:CD1	1:B:360:PHE:N	2.83	0.44
3:N:14:A:H2'	3:N:15:C:C6	2.52	0.44
1:A:515:PRO:HB2	1:A:518:VAL:CG2	2.47	0.44
1:B:503:TRP:CE2	1:B:683:PHE:HE1	2.35	0.44
1:A:531:ARG:C	1:A:531:ARG:CD	2.90	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:656:LEU:HD23	1:A:656:LEU:HA	1.54	0.44
1:B:42:VAL:O	1:B:46:LEU:HB3	2.17	0.44
1:A:197:TYR:CD1	1:A:232:ARG:NE	2.85	0.44
1:A:483:GLU:C	1:A:485:PHE:N	2.64	0.44
1:A:593:TYR:O	1:A:595:PRO:HD3	2.17	0.44
1:B:240:ARG:O	1:B:258:THR:HG23	2.18	0.44
1:A:34:ASP:HA	1:A:35:PRO:C	2.41	0.44
1:A:124:LEU:C	1:A:126:GLY:H	2.25	0.44
1:A:301:LEU:HD12	1:A:301:LEU:HA	1.74	0.44
1:B:503:TRP:HB2	1:B:662:LEU:HD22	2.00	0.44
1:B:20:ASN:HB2	1:B:21:PRO:HD2	1.99	0.43
1:B:384:HIS:ND1	1:B:385:PRO:HD2	2.33	0.43
1:B:545:ARG:HD3	1:B:546:ASP:N	2.33	0.43
1:A:287:ARG:HB3	1:A:582:TYR:CD1	2.53	0.43
1:A:589:ALA:O	1:A:592:LEU:HD13	2.18	0.43
1:A:593:TYR:CE2	1:A:595:PRO:HG3	2.51	0.43
1:A:652:LEU:CB	1:A:653:PRO:CD	2.96	0.43
1:B:506:PRO:HG2	1:B:666:VAL:HG21	2.00	0.43
1:A:515:PRO:HB2	1:A:518:VAL:HG23	1.99	0.43
1:B:384:HIS:CG	1:B:385:PRO:HD2	2.53	0.43
1:A:194:ARG:CB	1:A:201:THR:HG22	2.49	0.43
1:A:196:ALA:HB3	1:A:260:LEU:O	2.19	0.43
1:B:652:LEU:CB	1:B:653:PRO:HD2	2.46	0.43
1:A:411:PRO:O	1:A:412:PRO:C	2.62	0.43
1:B:349:LEU:HD21	1:B:351:LEU:HD21	2.01	0.43
1:B:636:PHE:O	1:B:639:THR:HB	2.19	0.43
1:A:117:LEU:HD22	1:A:155:LEU:HB2	2.00	0.43
1:B:329:LYS:HE2	1:B:329:LYS:HB3	1.82	0.43
1:A:652:LEU:H	1:A:652:LEU:HD12	1.84	0.43
1:B:475:VAL:HG23	1:B:477:PHE:HE1	1.84	0.43
1:B:645:SER:HB2	1:B:648:ALA:O	2.19	0.43
1:B:194:ARG:O	1:B:195:ASN:C	2.61	0.43
1:B:431:PRO:HB2	1:B:457:LYS:HB3	2.00	0.43
1:A:457:LYS:HG2	1:A:682:LEU:O	2.18	0.43
1:A:543:LEU:N	1:A:543:LEU:CD1	2.82	0.43
1:A:545:ARG:HD3	1:A:545:ARG:C	2.43	0.43
1:A:288:ARG:HD2	1:A:289:ARG:HH12	1.84	0.42
1:B:74:VAL:HG12	1:B:76:GLU:O	2.19	0.42
1:B:347:ALA:HB2	1:B:403:VAL:HG11	2.01	0.42
1:B:621:HIS:CE1	1:B:631:LEU:HD11	2.55	0.42
1:A:11:LEU:HD13	1:A:13:ARG:HG3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:LEU:HD11	1:A:245:ALA:HB2	2.02	0.42
1:B:332:ASP:HA	1:B:335:ARG:HD3	2.01	0.42
1:B:327:VAL:HG23	1:B:332:ASP:HB2	2.01	0.42
2:C:7:DA:H2'	2:C:8:DG:C8	2.53	0.42
1:A:4:LEU:HD12	1:A:4:LEU:H	1.84	0.42
1:A:408:VAL:HG12	1:A:410:THR:HG22	2.02	0.42
3:D:8:C:H2'	3:D:8:C:O2	2.20	0.42
1:A:576:SER:O	3:D:8:C:H5'	2.20	0.42
1:B:531:ARG:C	1:B:531:ARG:CD	2.93	0.42
1:B:344:THR:HG22	1:B:369:PHE:HE2	1.84	0.42
1:B:455:LEU:HD22	1:B:460:LEU:HD22	2.00	0.42
1:A:346:LEU:HD23	1:A:454:LEU:HD13	2.01	0.42
1:B:184:ALA:C	1:B:186:GLY:H	2.28	0.42
1:B:344:THR:HG22	1:B:369:PHE:CE2	2.54	0.42
1:B:575:LYS:HB3	1:B:651:ARG:HH22	1.83	0.42
1:B:656:LEU:HD23	1:B:656:LEU:HA	1.69	0.42
1:B:546:ASP:OD1	1:B:575:LYS:HE3	2.20	0.42
1:B:619:LEU:HD12	1:B:619:LEU:N	2.35	0.42
1:A:10:PHE:CB	1:A:310:ARG:HA	2.49	0.41
1:A:38:GLY:C	1:A:40:GLU:N	2.77	0.41
1:A:580:ARG:O	1:A:581:VAL:HG23	2.19	0.41
1:B:182:TRP:CZ3	1:B:190:PRO:HD3	2.46	0.41
1:A:297:ILE:O	1:A:301:LEU:HB2	2.21	0.41
1:A:621:HIS:ND1	1:A:631:LEU:CD1	2.83	0.41
1:A:672:ARG:CZ	1:B:518:VAL:HG22	2.50	0.41
1:B:228:ALA:HA	1:B:233:LEU:CB	2.50	0.41
1:A:10:PHE:HB3	1:A:310:ARG:CA	2.50	0.41
1:A:488:GLY:HA3	1:A:509:GLN:NE2	2.35	0.41
1:B:60:MET:SD	1:B:65:ALA:HB2	2.60	0.41
1:B:205:LEU:HB3	1:B:206:ARG:HG2	2.02	0.41
1:B:252:LYS:HA	1:B:253:PRO:HD3	1.87	0.41
1:A:506:PRO:HG2	1:A:666:VAL:HG21	2.03	0.41
1:A:639:THR:HG23	1:A:650:PRO:O	2.20	0.41
1:B:20:ASN:OD1	1:B:20:ASN:C	2.64	0.41
1:B:369:PHE:HE1	1:B:455:LEU:HD21	1.84	0.41
1:B:197:TYR:OH	1:B:256:HIS:HE1	2.02	0.41
3:D:14:A:H2'	3:D:15:C:H6	1.85	0.41
1:A:4:LEU:H	1:A:4:LEU:CD1	2.34	0.41
1:B:486:ARG:NH1	1:B:512:GLU:HB2	2.35	0.41
3:N:9:C:C2	3:N:10:U:C5	3.09	0.41
1:A:13:ARG:HB2	1:A:309:VAL:HG11	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:396:ARG:HH12	1:B:397:LYS:HE3	1.85	0.41
1:B:431:PRO:HG2	1:B:681:LYS:HD3	2.02	0.41
1:B:642:TYR:HA	1:B:643:PRO:HD3	1.72	0.41
1:A:288:ARG:O	1:A:292:GLU:HG3	2.20	0.41
1:A:319:PRO:HA	1:A:637:HIS:HE1	1.80	0.41
1:A:423:ALA:HB1	1:A:673:HIS:CE1	2.56	0.41
1:B:254:ILE:HA	1:B:255:PRO:HD3	1.77	0.41
2:M:7:DA:C2	3:N:14:A:C2	3.08	0.41
1:A:197:TYR:OH	5:A:689:PO4:P	2.79	0.41
1:A:352:ASP:HB3	1:A:437:VAL:HG21	2.03	0.41
1:A:144:GLY:CA	1:A:177:MET:HE2	2.50	0.40
1:A:582:TYR:HA	1:A:583:PRO:HD3	1.87	0.40
1:B:613:THR:HA	1:B:614:PRO:HD3	1.87	0.40
1:A:57:THR:HG22	1:A:66:SER:OG	2.21	0.40
1:A:155:LEU:HA	1:A:164:LEU:O	2.21	0.40
1:A:579:GLY:O	1:A:616:PRO:HD2	2.21	0.40
1:A:503:TRP:CZ2	1:A:683:PHE:HE1	2.39	0.40
1:B:597:GLU:O	1:B:598:ASP:C	2.65	0.40
1:B:661:ARG:N	1:B:661:ARG:CD	2.84	0.40
2:C:3:DA:H2'	2:C:4:DG:C8	2.57	0.40
1:A:319:PRO:HA	1:A:637:HIS:ND1	2.36	0.40
1:A:559:GLU:O	1:A:562:ALA:HB3	2.21	0.40
1:B:284:GLU:HA	1:B:287:ARG:HD2	2.03	0.40
1:B:547:GLY:HA2	3:N:9:C:OP1	2.20	0.40
1:A:104:GLY:O	1:A:107:SER:OG	2.36	0.40
1:A:461:GLN:CD	1:A:500:HIS:HB3	2.47	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	646/685 (94%)	567 (88%)	72 (11%)	7 (1%)	11	42
1	B	643/685 (94%)	577 (90%)	59 (9%)	7 (1%)	11	42
All	All	1289/1370 (94%)	1144 (89%)	131 (10%)	14 (1%)	11	42

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	597	GLU
1	A	5	GLY
1	B	39	ARG
1	B	354	ALA
1	A	480	GLY
1	B	446	ARG
1	A	126	GLY
1	B	5	GLY
1	A	495	GLY
1	B	653	PRO
1	A	323	GLY
1	B	127	VAL
1	A	438	PRO
1	B	643	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	400/549 (73%)	365 (91%)	35 (9%)	9	34
1	B	407/549 (74%)	382 (94%)	25 (6%)	17	47
All	All	807/1098 (74%)	747 (93%)	60 (7%)	13	41

All (60) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	LEU
1	A	10	PHE

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Mol	Chain	Res	Type
1	A	11	LEU
1	A	31	VAL
1	A	57	THR
1	A	90	LEU
1	A	108	VAL
1	A	134	VAL
1	A	155	LEU
1	A	159	ASP
1	A	165	LEU
1	A	177	MET
1	A	283	TRP
1	A	309	VAL
1	A	327	VAL
1	A	376	LEU
1	A	406	VAL
1	A	432	SER
1	A	448	GLU
1	A	451	LEU
1	A	500	HIS
1	A	502	LEU
1	A	505	LEU
1	A	531	ARG
1	A	545	ARG
1	A	566	ILE
1	A	570	LEU
1	A	592	LEU
1	A	600	THR
1	A	622	GLU
1	A	641	LEU
1	A	652	LEU
1	A	656	LEU
1	A	661	ARG
1	A	663	VAL
1	B	4	LEU
1	B	10	PHE
1	B	11	LEU
1	B	90	LEU
1	B	108	VAL
1	B	134	VAL
1	B	155	LEU
1	B	165	LEU
1	B	283	TRP

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Mol	Chain	Res	Type
1	B	376	LEU
1	B	406	VAL
1	B	451	LEU
1	B	454	LEU
1	B	500	HIS
1	B	504	THR
1	B	505	LEU
1	B	531	ARG
1	B	545	ARG
1	B	566	ILE
1	B	569	ASP
1	B	600	THR
1	B	641	LEU
1	B	661	ARG
1	B	663	VAL
1	B	674	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	382	HIS
1	A	387	GLN
1	A	404	GLN
1	A	461	GLN
1	A	509	GLN
1	B	12	ASN
1	B	139	HIS
1	B	256	HIS
1	B	355	GLN
1	B	404	GLN
1	B	633	HIS
1	B	634	GLN
1	B	637	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
3	D	10/19 (52%)	0	0
3	N	9/19 (47%)	1 (11%)	0
All	All	19/38 (50%)	1 (5%)	0

All (1) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
3	N	10	U

There are no RNA pucker outliers to report.

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 8 ligands modelled in this entry, 6 are monoatomic - leaving 2 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	PO4	B	689	-	4,4,4	3.07	4 (100%)	6,6,6	0.46	0
5	PO4	A	689	-	4,4,4	3.07	4 (100%)	6,6,6	0.46	0

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	689	PO4	P-O1	4.69	1.61	1.50
5	B	689	PO4	P-O1	4.69	1.61	1.50
5	A	689	PO4	P-O3	2.38	1.61	1.54
5	B	689	PO4	P-O3	2.37	1.61	1.54
5	B	689	PO4	P-O2	2.35	1.61	1.54
5	A	689	PO4	P-O2	2.34	1.61	1.54
5	A	689	PO4	P-O4	-2.14	1.48	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	689	PO4	P-O4	-2.13	1.48	1.54

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	689	PO4	1	0
5	A	689	PO4	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	660/685 (96%)	-0.19	9 (1%) 73 54	54, 93, 143, 187	0
1	B	657/685 (95%)	-0.19	8 (1%) 76 58	55, 89, 151, 292	0
2	C	16/21 (76%)	-0.31	0 100 100	76, 101, 134, 140	0
2	M	16/21 (76%)	-0.30	0 100 100	81, 98, 140, 154	0
3	D	12/19 (63%)	-0.25	0 100 100	98, 116, 124, 159	0
3	N	10/19 (52%)	-0.50	0 100 100	105, 113, 117, 118	0
All	All	1371/1450 (94%)	-0.19	17 (1%) 76 58	54, 92, 148, 292	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	480	GLY	5.7
1	B	497	ASP	3.6
1	A	496	GLY	3.5
1	A	612	GLY	3.2
1	A	611	ARG	3.1
1	A	268	GLU	3.0
1	B	246	ASP	2.9
1	A	222	SER	2.8
1	B	491	ALA	2.6
1	B	222	SER	2.6
1	A	625	ASP	2.4
1	B	612	GLY	2.3
1	A	269	ASP	2.3
1	A	279	LEU	2.2
1	B	190	PRO	2.2
1	B	493	ALA	2.1
1	A	281	LEU	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, 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($\text{\AA}^2$)	Q<0.9
5	PO4	B	689	5/5	0.63	0.14	140,140,140,140	0
4	MG	B	687	1/1	0.68	0.12	95,95,95,95	0
5	PO4	A	689	5/5	0.70	0.12	129,129,129,129	0
4	MG	B	688	1/1	0.85	0.15	101,101,101,101	0
4	MG	A	686	1/1	0.93	0.08	66,66,66,66	0
4	MG	A	687	1/1	0.94	0.06	61,61,61,61	0
4	MG	A	688	1/1	0.95	0.06	68,68,68,68	0
4	MG	B	686	1/1	0.98	0.11	55,55,55,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.