



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:38 PM UTC

PDB ID : 6D92 / pdb_00006d92
Title : Ternary RsAgo Complex with Guide RNA and Target DNA Containing A-A non-canonical pair at position 3
Authors : Liu, Y.; Esyunina, D.; Olovnikov, I.; Teplova, M.; Patel, D.J.
Deposited on : 2018-04-27
Resolution : 1.81 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

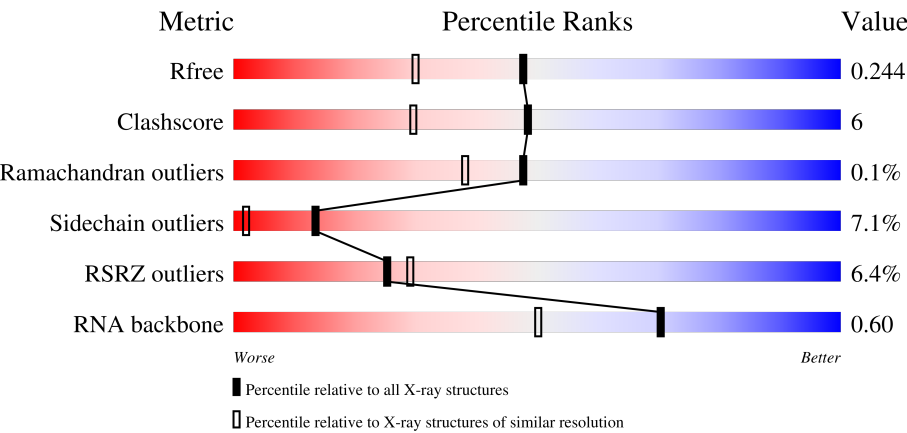
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.81 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	1112 (1.82-1.82)
Clashscore	190562	1148 (1.82-1.82)
Ramachandran outliers	187476	1140 (1.82-1.82)
Sidechain outliers	187428	1140 (1.82-1.82)
RSRZ outliers	180081	1112 (1.82-1.82)
RNA backbone	3983	1006 (2.22-1.42)

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	791	
1	F	791	
2	C	18	
2	H	18	

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Mol	Chain	Length	Quality of chain
3	G	19	 84%16%
4	J	21	 5%76%24%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 14158 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 Uncharacterized protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	758	Total	C	N	O	S	0	1	0
			5895	3742	1057	1080	16			
1	F	758	Total	C	N	O	S	0	0	0
			5800	3693	1032	1059	16			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	initiating methionine	UNP A4WYU7
A	-12	HIS	-	expression tag	UNP A4WYU7
A	-11	HIS	-	expression tag	UNP A4WYU7
A	-10	HIS	-	expression tag	UNP A4WYU7
A	-9	HIS	-	expression tag	UNP A4WYU7
A	-8	HIS	-	expression tag	UNP A4WYU7
A	-7	HIS	-	expression tag	UNP A4WYU7
A	-6	ASP	-	expression tag	UNP A4WYU7
A	-5	TYR	-	expression tag	UNP A4WYU7
A	-4	LYS	-	expression tag	UNP A4WYU7
A	-3	ASP	-	expression tag	UNP A4WYU7
A	-2	ASP	-	expression tag	UNP A4WYU7
A	-1	ASP	-	expression tag	UNP A4WYU7
A	0	ASP	-	expression tag	UNP A4WYU7
A	1	LYS	-	expression tag	UNP A4WYU7
F	-13	MET	-	initiating methionine	UNP A4WYU7
F	-12	HIS	-	expression tag	UNP A4WYU7
F	-11	HIS	-	expression tag	UNP A4WYU7
F	-10	HIS	-	expression tag	UNP A4WYU7
F	-9	HIS	-	expression tag	UNP A4WYU7
F	-8	HIS	-	expression tag	UNP A4WYU7
F	-7	HIS	-	expression tag	UNP A4WYU7
F	-6	ASP	-	expression tag	UNP A4WYU7
F	-5	TYR	-	expression tag	UNP A4WYU7
F	-4	LYS	-	expression tag	UNP A4WYU7

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Chain	Residue	Modelled	Actual	Comment	Reference
F	-3	ASP	-	expression tag	UNP A4WYU7
F	-2	ASP	-	expression tag	UNP A4WYU7
F	-1	ASP	-	expression tag	UNP A4WYU7
F	0	ASP	-	expression tag	UNP A4WYU7
F	1	LYS	-	expression tag	UNP A4WYU7

- Molecule 2 is a RNA chain called RNA (5'-R(P*UP*UP*AP*CP*UP*GP*CP*AP*CP*AP*GP*GP*UP*GP*AP*CP*GP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	18	Total	C	N	O	P	0	0	0
			386	172	70	126	18			
2	H	18	Total	C	N	O	P	0	0	0
			386	172	70	126	18			

- Molecule 3 is a DNA chain called DNA (5'-D(P*TP*CP*GP*TP*CP*AP*CP*CP*TP*GP*TP*GP*CP*AP*GP*AP*AP*AP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	G	19	Total	C	N	O	P	0	0	0
			387	184	71	113	19			

- Molecule 4 is a DNA chain called DNA (5'-D(P*TP*GP*TP*CP*GP*TP*CP*AP*CP*CP*TP*GP*TP*GP*CP*AP*GP*AP*AP*AP*C)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	J	21	Total	C	N	O	P	0	0	0
			430	204	78	127	21			

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	F	1	Total C O 4 2 2	0	0
5	F	1	Total C O 4 2 2	0	0

- Molecule 6 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	C	1	Total Mg 1 1	0	0
6	H	1	Total Mg 1 1	0	0

- Molecule 7 is (4S)-2-METHYL-2,4-PENTANEDIOL (CCD ID: MPD) (formula: C₆H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	H	1	Total	C	O	0	0
			8	6	2		

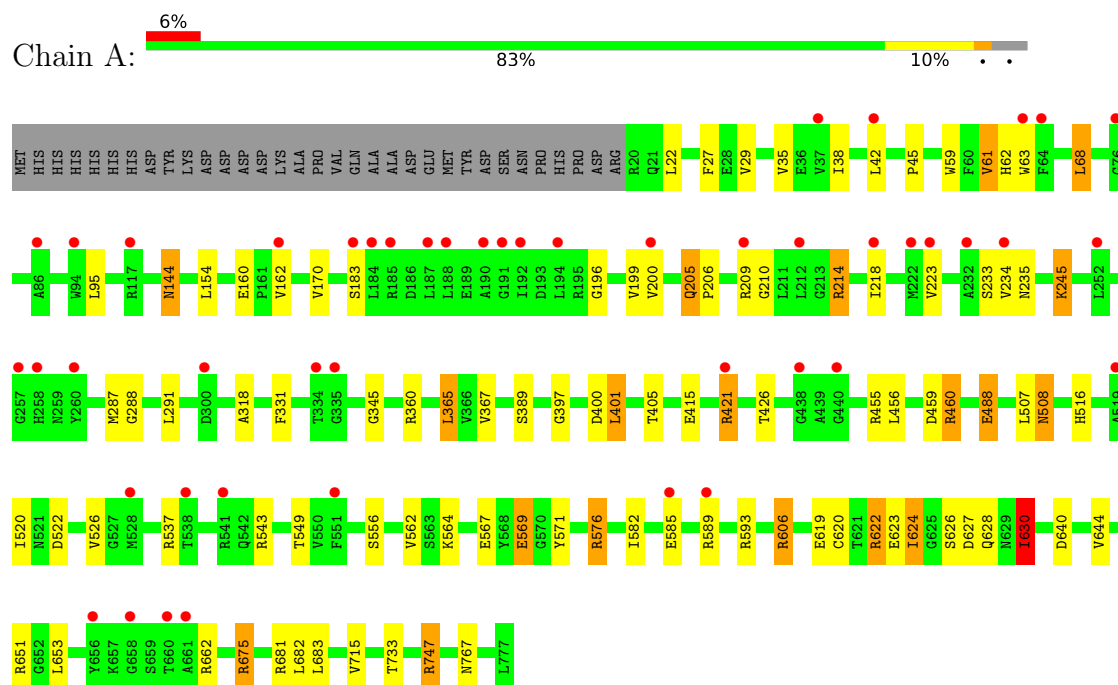
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	343	Total	O	0	0
			343	343		
8	C	50	Total	O	0	0
			50	50		
8	G	41	Total	O	0	0
			41	41		
8	F	318	Total	O	0	0
			318	318		
8	H	56	Total	O	0	0
			56	56		
8	J	40	Total	O	0	0
			40	40		

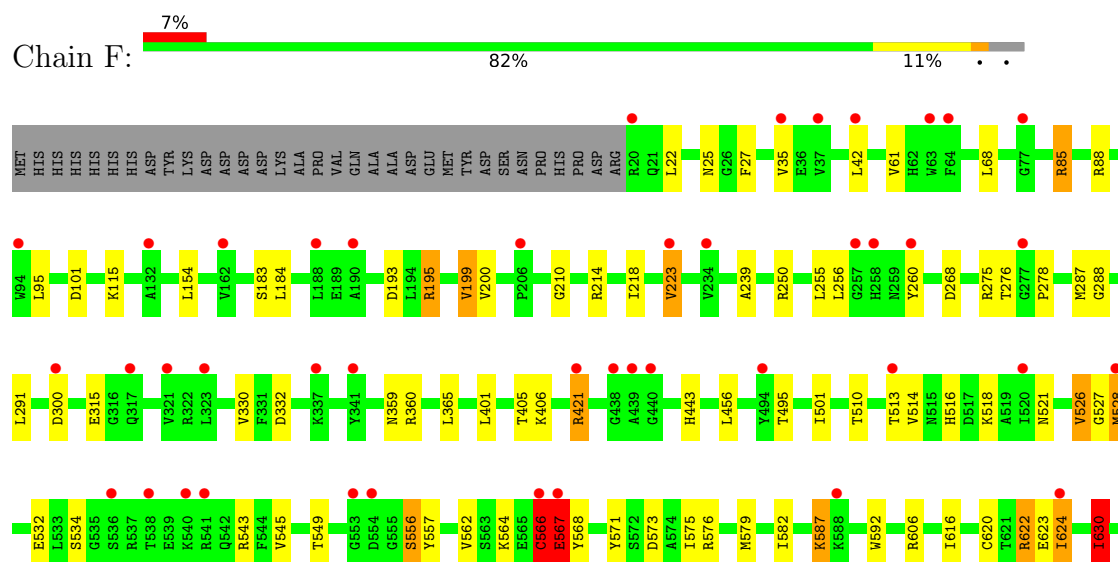
3 Residue-property plots [i](#)

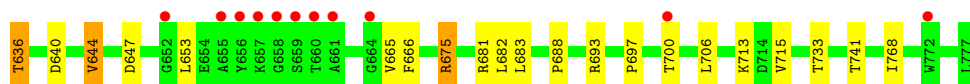
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

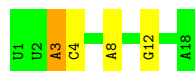
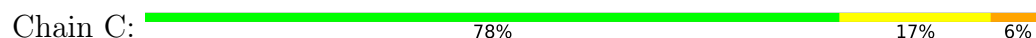


• Molecule 1: Uncharacterized protein

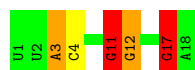




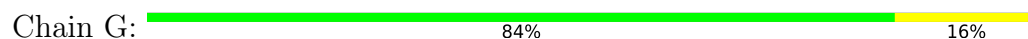
- Molecule 2: RNA (5'-R(P*UP*UP*AP*CP*UP*GP*CP*AP*CP*AP*GP*GP*UP*GP*AP*C P*GP*A)-3')



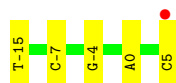
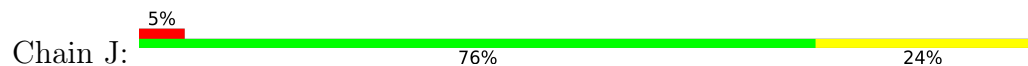
- Molecule 2: RNA (5'-R(P*UP*UP*AP*CP*UP*GP*CP*AP*CP*AP*GP*GP*UP*GP*AP*C P*GP*A)-3')



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



- Molecule 4: DNA (5'-D(P*TP*GP*TP*CP*GP*TP*CP*AP*CP*CP*TP*GP*TP*GP*CP*A P*GP*AP*AP*AP*C)-3')



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.24Å 118.94Å 118.56Å 90.00° 95.72° 90.00°	Depositor
Resolution (Å)	42.81 – 1.81 42.81 – 1.81	Depositor EDS
% Data completeness (in resolution range)	93.4 (42.81-1.81) 93.6 (42.81-1.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 1.81Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.198 , 0.240 0.202 , 0.244	Depositor DCC
R_{free} test set	8084 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å ²)	34.5	Xtriage
Anisotropy	0.285	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 32.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14158	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.30% 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, ACT, MPD

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	0.72	0/6024	0.89	1/8179 (0.0%)
1	F	0.69	0/5929	0.89	6/8063 (0.1%)
2	C	0.52	0/431	0.95	1/668 (0.1%)
2	H	0.50	0/431	1.06	3/668 (0.4%)
3	G	0.43	0/433	0.78	0/665
4	J	0.39	0/481	0.79	0/738
All	All	0.68	0/13729	0.89	11/18981 (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	F	0	1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	11	G	C2'-C3'-O3'	-10.56	97.86	113.70
2	H	3	A	C4'-C3'-O3'	-5.85	104.23	113.00
1	A	630	ILE	CB-CA-C	-5.84	101.94	110.81
2	C	3	A	C4'-C3'-O3'	-5.60	104.60	113.00
1	F	568	TYR	N-CA-CB	5.58	118.05	109.91
2	H	17	G	C4'-C3'-O3'	-5.53	104.70	113.00
1	F	510	THR	CA-C-N	5.43	124.62	118.97
1	F	510	THR	C-N-CA	5.43	124.62	118.97
1	F	567	GLU	N-CA-C	5.42	124.38	113.31
1	F	630	ILE	CB-CA-C	-5.20	102.90	110.81

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	F	700	THR	N-CA-C	-5.02	103.38	110.31

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	566	CYS	Peptide

5.2 Too-close contacts [i](#)

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	5895	0	5810	56	0
1	F	5800	0	5663	70	0
2	C	386	0	195	3	0
2	H	386	0	195	7	0
3	G	387	0	214	2	0
4	J	430	0	237	7	0
5	A	8	0	6	0	0
5	F	8	0	6	0	0
6	C	1	0	0	0	0
6	H	1	0	0	0	0
7	H	8	0	14	1	0
8	A	343	0	0	7	1
8	C	50	0	0	1	0
8	F	318	0	0	10	1
8	G	41	0	0	0	0
8	H	56	0	0	3	0
8	J	40	0	0	2	0
All	All	14158	0	12340	141	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 6.

All (141) 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)
4:J:-15:DT:O4	8:J:101:HOH:O	1.69	1.08
1:F:513:THR:HG23	1:F:557:TYR:O	1.65	0.97
1:A:61:VAL:HG13	1:A:68:LEU:HD21	1.59	0.85
1:A:585:GLU:OE2	1:A:589:ARG:NH2	2.13	0.81
1:F:193:ASP:OD1	1:F:195:ARG:HD3	1.80	0.80
4:J:-4:DG:OP2	8:J:102:HOH:O	2.06	0.73
1:A:576:ARG:HD3	1:A:619:GLU:HG3	1.69	0.73
1:F:573:ASP:OD1	1:F:576:ARG:NH1	2.25	0.70
1:A:360:ARG:HB3	1:A:405:THR:HG23	1.73	0.69
1:F:513:THR:HG22	1:F:514:VAL:N	2.07	0.69
1:F:200:VAL:HG21	8:F:1006:HOH:O	1.93	0.67
1:A:624:ILE:HD11	1:A:630:ILE:HD12	1.79	0.65
1:A:421:ARG:HG2	1:A:456:LEU:HD21	1.79	0.64
1:A:624:ILE:CD1	1:A:630:ILE:HD12	2.29	0.63
1:F:260:TYR:OH	4:J:0:DA:OP1	2.15	0.62
1:A:154:LEU:HD11	1:A:287:MET:HE2	1.80	0.62
1:A:537:ARG:HD3	8:C:201:HOH:O	1.98	0.62
1:F:88:ARG:NH1	8:F:902:HOH:O	2.26	0.61
2:H:11:G:H2'	8:H:241:HOH:O	2.00	0.60
1:F:526:VAL:HG13	1:F:579:MET:HE2	1.83	0.60
1:F:85:ARG:HG2	1:F:85:ARG:HH21	1.65	0.60
1:F:624:ILE:HG12	1:F:630:ILE:HD12	1.83	0.59
1:A:488:GLU:OE1	1:A:488:GLU:HA	2.03	0.59
1:F:675:ARG:HG3	1:F:681:ARG:CZ	2.32	0.59
1:A:681:ARG:CZ	1:A:715:VAL:HG13	2.33	0.59
1:F:513:THR:HG21	1:F:556:SER:HB3	1.85	0.59
1:A:205:GLN:HB3	1:A:206:PRO:CD	2.33	0.59
1:A:624:ILE:CG1	1:A:630:ILE:HD12	2.32	0.59
2:H:17:G:H5''	2:H:17:G:H8	1.67	0.58
1:A:589:ARG:N	1:A:589:ARG:HD3	2.18	0.58
1:F:513:THR:HG21	1:F:556:SER:CB	2.33	0.58
1:F:360:ARG:HB3	1:F:405:THR:HG23	1.83	0.58
1:F:268:ASP:OD2	1:F:693:ARG:NH2	2.34	0.58
1:A:144:ASN:HB2	8:A:1118:HOH:O	2.02	0.57
1:F:276:THR:HG22	1:F:278:PRO:HD2	1.86	0.57
1:A:624:ILE:HG12	1:A:630:ILE:HD12	1.85	0.57
1:F:545:VAL:CG2	1:F:567:GLU:HB3	2.34	0.57
1:F:566:CYS:HB2	8:F:914:HOH:O	2.04	0.57
1:F:526:VAL:CG1	1:F:579:MET:HE2	2.35	0.56
1:A:233:SER:O	1:A:234:VAL:HG23	2.05	0.56
1:F:549:THR:HG21	1:F:582:ILE:CD1	2.35	0.56
1:A:630:ILE:N	1:A:630:ILE:HD13	2.21	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:LEU:HD13	1:A:367:VAL:CG2	2.36	0.56
1:A:624:ILE:HD11	1:A:630:ILE:CD1	2.36	0.55
1:F:200:VAL:HG23	1:F:210:GLY:C	2.32	0.54
1:A:624:ILE:HD12	1:A:628:GLN:HB2	1.88	0.54
1:F:528:MET:HE1	1:F:571:TYR:HB2	1.89	0.54
1:F:359:ASN:HD21	1:F:443:HIS:CG	2.26	0.54
1:F:300:ASP:OD1	8:F:901:HOH:O	2.18	0.53
1:F:513:THR:CG2	1:F:514:VAL:N	2.71	0.53
1:A:640:ASP:O	3:G:-7:DC:H3'	2.09	0.53
4:J:5:DC:C2'	4:J:5:DC:O2	2.57	0.53
1:F:516:HIS:HB3	1:F:557:TYR:CE2	2.44	0.52
1:F:218:ILE:HD13	1:F:223:VAL:HA	1.90	0.52
1:F:606:ARG:NH1	8:F:904:HOH:O	2.34	0.52
1:A:767:ASN:HB2	8:A:1017:HOH:O	2.10	0.51
1:A:287:MET:HE3	1:A:291:LEU:HG	1.92	0.51
1:A:162:VAL:HG12	1:A:318:ALA:HA	1.93	0.51
1:F:624:ILE:CG1	1:F:630:ILE:HD12	2.41	0.51
1:A:200:VAL:HG23	1:A:210:GLY:C	2.36	0.51
1:A:218:ILE:CD1	1:A:223:VAL:HG12	2.41	0.50
2:H:11:G:H2'	2:H:12:G:O4'	2.10	0.50
1:F:513:THR:CG2	1:F:557:TYR:O	2.50	0.50
1:F:624:ILE:CD1	1:F:630:ILE:HD12	2.41	0.50
1:F:521:ASN:CB	8:F:1071:HOH:O	2.60	0.50
1:F:545:VAL:HG22	1:F:567:GLU:HB3	1.92	0.50
4:J:5:DC:O2	4:J:5:DC:H2'	2.10	0.50
1:F:27:PHE:CZ	1:F:288:GLY:HA3	2.46	0.50
1:A:160:GLU:OE1	1:A:675:ARG:NH2	2.44	0.50
1:A:662:ARG:CD	8:A:1097:HOH:O	2.59	0.50
2:H:11:G:H3'	8:H:241:HOH:O	2.12	0.49
1:F:624:ILE:HD11	1:F:630:ILE:CD1	2.41	0.49
1:F:526:VAL:HG11	1:F:579:MET:HG2	1.95	0.49
1:A:218:ILE:HD13	1:A:223:VAL:HG12	1.94	0.48
1:A:549:THR:HG21	1:A:582:ILE:CD1	2.43	0.48
1:F:255:LEU:HB2	1:F:256:LEU:HD22	1.96	0.48
1:A:622:ARG:HG3	1:A:623:GLU:N	2.28	0.47
2:H:17:G:H5''	2:H:17:G:C8	2.48	0.47
1:F:154:LEU:HD11	1:F:287:MET:HE2	1.95	0.47
1:F:532:GLU:OE2	1:F:543:ARG:HD3	2.14	0.47
1:F:218:ILE:HD12	1:F:223:VAL:HG12	1.97	0.47
1:F:647:ASP:O	1:F:666:PHE:HA	2.15	0.46
1:F:575:ILE:CG2	1:F:579:MET:HE3	2.45	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:250:ARG:HD2	8:F:1000:HOH:O	2.16	0.46
1:F:287:MET:HE3	1:F:291:LEU:HG	1.98	0.46
1:A:95:LEU:C	1:A:95:LEU:HD13	2.40	0.46
1:F:516:HIS:CD2	1:F:518:LYS:H	2.34	0.45
1:F:154:LEU:HD21	1:F:287:MET:HE2	1.97	0.45
1:A:569:GLU:H	1:A:569:GLU:CD	2.23	0.45
1:A:38:ILE:HD12	1:A:59:TRP:CH2	2.52	0.45
1:A:606:ARG:HH11	1:A:606:ARG:HG3	1.80	0.45
1:A:45:PRO:HG3	1:A:63:TRP:CZ2	2.52	0.45
1:F:636:THR:HG22	8:F:1143:HOH:O	2.16	0.45
1:F:624:ILE:HD11	1:F:630:ILE:HD12	1.99	0.45
1:A:397:GLY:O	1:A:401:LEU:HB2	2.17	0.45
1:A:245:LYS:HE3	2:C:8:A:C1'	2.46	0.45
1:F:25:ASN:CG	1:F:644:VAL:HG13	2.42	0.45
1:F:95:LEU:C	1:F:95:LEU:HD13	2.41	0.45
1:A:627:ASP:HB2	8:A:1044:HOH:O	2.16	0.44
1:A:459:ASP:OD1	1:A:460:ARG:HG3	2.17	0.44
1:F:85:ARG:HH21	1:F:85:ARG:CG	2.30	0.44
1:F:184:LEU:CB	8:F:1173:HOH:O	2.65	0.44
1:F:681:ARG:CZ	1:F:715:VAL:HG13	2.48	0.44
2:C:3:A:H2'	2:C:4:C:C6	2.53	0.44
2:H:11:G:C3'	8:H:241:HOH:O	2.66	0.44
1:A:154:LEU:HD21	1:A:287:MET:HE2	1.99	0.43
1:A:245:LYS:HE3	2:C:8:A:H1'	1.99	0.43
1:F:653:LEU:HD13	1:F:688:PRO:HB2	2.00	0.43
1:F:706:LEU:HD21	1:F:713:LYS:HA	1.99	0.43
1:A:62:HIS:HD2	8:A:1187:HOH:O	2.01	0.43
1:F:527:GLY:HA3	1:F:741:THR:HB	2.00	0.43
1:F:576:ARG:HB3	1:F:616:ILE:CD1	2.49	0.43
3:G:2:DA:H2''	3:G:3:DA:C8	2.53	0.42
1:F:620:CYS:O	1:F:624:ILE:HG23	2.18	0.42
1:A:620:CYS:O	1:A:624:ILE:HG23	2.20	0.42
1:A:662:ARG:HD3	8:A:1097:HOH:O	2.20	0.42
1:F:501:ILE:HD11	1:F:733:THR:HG22	2.01	0.42
1:A:27:PHE:CZ	1:A:288:GLY:HA3	2.54	0.42
1:A:675:ARG:HG3	1:A:681:ARG:CZ	2.50	0.42
1:F:513:THR:HG22	1:F:514:VAL:H	1.80	0.42
4:J:-15:DT:O2	4:J:-15:DT:O4'	2.37	0.42
1:A:507:LEU:O	1:A:508:ASN:HB2	2.19	0.42
1:A:196:GLY:HA2	1:A:214:ARG:HG2	2.01	0.42
1:F:275:ARG:HB3	1:F:697:PRO:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:513:THR:CG2	1:F:514:VAL:H	2.32	0.42
1:A:29:VAL:HG21	1:A:170:VAL:HG23	2.01	0.41
1:F:587:LYS:HD2	1:F:592:TRP:CE3	2.55	0.41
1:F:622:ARG:HG3	1:F:623:GLU:N	2.36	0.41
1:F:421:ARG:HA	1:F:421:ARG:HD3	1.70	0.41
1:F:640:ASP:O	4:J:-7:DC:H3'	2.20	0.41
1:F:567:GLU:HA	8:F:1008:HOH:O	2.20	0.41
1:A:571:TYR:CD1	1:A:571:TYR:C	2.99	0.41
1:A:651[B]:ARG:HA	1:A:651[B]:ARG:HD2	1.71	0.40
1:A:747:ARG:NH1	8:A:908:HOH:O	2.46	0.40
1:A:331:PHE:CE1	1:A:345:GLY:HA3	2.56	0.40
1:A:606:ARG:HH11	1:A:606:ARG:CG	2.35	0.40
1:F:199:VAL:CG1	1:F:239:ALA:HB1	2.51	0.40
1:F:665:VAL:O	1:F:666:PHE:HB2	2.20	0.40
2:H:3:A:H2'	2:H:4:C:C6	2.57	0.40
1:F:421:ARG:HD2	1:F:456:LEU:HD21	2.03	0.40
7:H:102:MPD:H12	7:H:102:MPD:O4	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)
8:A:1227:HOH:O	8:F:1174:HOH:O[2_455]	1.26	0.94

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	757/791 (96%)	744 (98%)	12 (2%)	1 (0%)	48	38
1	F	756/791 (96%)	740 (98%)	15 (2%)	1 (0%)	48	38
All	All	1513/1582 (96%)	1484 (98%)	27 (2%)	2 (0%)	48	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	508	ASN
1	F	768	ILE

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	613/672 (91%)	566 (92%)	47 (8%)	12	1
1	F	591/672 (88%)	553 (94%)	38 (6%)	16	3
All	All	1204/1344 (90%)	1119 (93%)	85 (7%)	13	2

All (85) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	LEU
1	A	35	VAL
1	A	42	LEU
1	A	61	VAL
1	A	68	LEU
1	A	144	ASN
1	A	183	SER
1	A	199	VAL
1	A	205	GLN
1	A	209	ARG
1	A	214	ARG
1	A	235	ASN
1	A	245	LYS
1	A	365	LEU
1	A	389	SER
1	A	400	ASP
1	A	401	LEU
1	A	415	GLU
1	A	421	ARG
1	A	426	THR
1	A	455	ARG

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Mol	Chain	Res	Type
1	A	460	ARG
1	A	488	GLU
1	A	516	HIS
1	A	520	ILE
1	A	522	ASP
1	A	526	VAL
1	A	543	ARG
1	A	556	SER
1	A	562	VAL
1	A	564	LYS
1	A	567	GLU
1	A	569	GLU
1	A	576	ARG
1	A	593	ARG
1	A	606	ARG
1	A	622	ARG
1	A	624	ILE
1	A	626	SER
1	A	630	ILE
1	A	644	VAL
1	A	653	LEU
1	A	675	ARG
1	A	682	LEU
1	A	683	LEU
1	A	733	THR
1	A	747	ARG
1	F	22	LEU
1	F	35	VAL
1	F	42	LEU
1	F	61	VAL
1	F	68	LEU
1	F	85	ARG
1	F	101	ASP
1	F	115	LYS
1	F	183	SER
1	F	195	ARG
1	F	199	VAL
1	F	214	ARG
1	F	223	VAL
1	F	315	GLU
1	F	330	VAL
1	F	332	ASP

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Mol	Chain	Res	Type
1	F	365	LEU
1	F	401	LEU
1	F	406	LYS
1	F	421	ARG
1	F	495	THR
1	F	526	VAL
1	F	528	MET
1	F	534	SER
1	F	556	SER
1	F	562	VAL
1	F	564	LYS
1	F	566	CYS
1	F	567	GLU
1	F	587	LYS
1	F	622	ARG
1	F	624	ILE
1	F	630	ILE
1	F	636	THR
1	F	644	VAL
1	F	675	ARG
1	F	682	LEU
1	F	683	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	62	HIS
1	A	124	GLN
1	A	205	GLN
1	A	270	GLN
1	A	348	GLN
1	A	508	ASN
1	A	542	GLN
1	F	359	ASN
1	F	508	ASN
1	F	516	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	C	17/18 (94%)	1 (5%)	0
2	H	17/18 (94%)	3 (17%)	1 (5%)
All	All	34/36 (94%)	4 (11%)	1 (2%)

All (4) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	C	12	G
2	H	11	G
2	H	12	G
2	H	17	G

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	H	11	G

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 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	ACT	F	801	-	3,3,3	0.87	0	3,3,3	0.64	0
5	ACT	A	802	-	3,3,3	0.79	0	3,3,3	0.71	0
5	ACT	A	801	-	3,3,3	0.83	0	3,3,3	0.85	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	MPD	H	102	-	7,7,7	0.33	0	9,10,10	0.50	0
5	ACT	F	802	-	3,3,3	0.78	0	3,3,3	0.87	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	MPD	H	102	-	-	3/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	H	102	MPD	C1-C2-C3-C4
7	H	102	MPD	O2-C2-C3-C4
7	H	102	MPD	CM-C2-C3-C4

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	H	102	MPD	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	758/791 (95%)	0.49	47 (6%)	26	30	22, 38, 72, 102	2 (0%)
1	F	758/791 (95%)	0.71	54 (7%)	22	25	22, 40, 64, 96	0
2	C	18/18 (100%)	-0.40	0	100	100	24, 30, 55, 83	0
2	H	18/18 (100%)	-0.35	0	100	100	27, 32, 45, 52	0
3	G	19/19 (100%)	-0.14	0	100	100	30, 38, 94, 115	0
4	J	21/21 (100%)	0.01	1 (4%)	35	41	29, 39, 87, 104	0
All	All	1592/1658 (96%)	0.56	102 (6%)	25	29	22, 39, 70, 115	2 (0%)

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	188	LEU	3.7
1	F	659	SER	3.5
1	F	658	GLY	3.5
1	F	323	LEU	3.4
1	A	187	LEU	3.3
1	A	257	GLY	3.3
1	F	258	HIS	3.1
1	A	300	ASP	3.1
1	F	566	CYS	3.1
1	F	520	ILE	3.1
1	A	117	ARG	3.1
1	F	317	GLN	3.0
1	F	661	ALA	3.0
1	F	538	THR	2.9
1	A	190	ALA	2.9
1	A	334	THR	2.9
1	A	232	ALA	2.9
1	A	658	GLY	2.9
1	F	541	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	218	ILE	2.8
1	A	191	GLY	2.8
1	F	206	PRO	2.8
1	F	188	LEU	2.8
1	F	77	GLY	2.8
1	F	260	TYR	2.7
1	F	656	TYR	2.7
1	F	540	LYS	2.7
1	F	567	GLU	2.7
1	F	588	LYS	2.7
1	F	657	LYS	2.7
1	F	321	VAL	2.7
1	F	20	ARG	2.7
1	F	257	GLY	2.6
1	F	439	ALA	2.6
1	A	234	VAL	2.6
1	A	660	THR	2.6
1	A	223	VAL	2.6
1	F	162	VAL	2.6
1	A	258	HIS	2.5
1	A	260	TYR	2.5
1	A	64	PHE	2.5
1	F	190	ALA	2.5
1	A	222	MET	2.5
1	A	63	TRP	2.5
1	A	183	SER	2.5
1	A	162	VAL	2.5
1	F	223	VAL	2.5
1	F	660	THR	2.5
1	F	277	GLY	2.5
1	F	438	GLY	2.5
1	A	86	ALA	2.5
1	F	440	GLY	2.4
1	F	652	GLY	2.4
1	A	185	ARG	2.4
1	A	42	LEU	2.4
1	F	700	THR	2.4
1	F	234	VAL	2.4
1	F	341	TYR	2.3
1	A	252	LEU	2.3
1	F	64	PHE	2.3
1	F	528	MET	2.3

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Mol	Chain	Res	Type	RSRZ
1	F	337	LYS	2.3
1	F	42	LEU	2.3
1	F	63	TRP	2.2
1	F	624	ILE	2.2
1	A	200	VAL	2.2
1	A	209	ARG	2.2
1	A	194	LEU	2.2
1	A	94	TRP	2.2
1	F	553	GLY	2.2
1	F	494	TYR	2.2
4	J	5	DC	2.2
1	A	212	LEU	2.2
1	A	440	GLY	2.2
1	F	655	ALA	2.2
1	A	541	ARG	2.2
1	F	772	TRP	2.1
1	F	300	ASP	2.1
1	F	554	ASP	2.1
1	A	335	GLY	2.1
1	F	664	GLY	2.1
1	A	528	MET	2.1
1	A	519	ALA	2.1
1	F	132	ALA	2.1
1	F	94	TRP	2.1
1	A	76	GLY	2.1
1	F	35	VAL	2.1
1	A	661	ALA	2.1
1	A	656	TYR	2.1
1	A	421	ARG	2.1
1	F	421	ARG	2.1
1	F	536	SER	2.1
1	A	538	THR	2.1
1	A	551	PHE	2.0
1	A	37	VAL	2.0
1	A	184	LEU	2.0
1	A	589	ARG	2.0
1	A	585	GLU	2.0
1	A	192	ILE	2.0
1	A	438	GLY	2.0
1	F	513	THR	2.0
1	F	37	VAL	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](#)

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	ACT	F	801	4/4	0.45	0.20	71,74,77,77	0
5	ACT	A	801	4/4	0.85	0.20	61,63,63,63	0
5	ACT	F	802	4/4	0.86	0.16	64,64,65,66	0
7	MPD	H	102	8/8	0.88	0.14	54,56,57,58	0
5	ACT	A	802	4/4	0.91	0.09	61,65,67,68	0
6	MG	C	101	1/1	0.98	0.04	21,21,21,21	0
6	MG	H	101	1/1	0.99	0.03	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.