



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 1, 2026 – 09:20 PM UTC

PDB ID : 6D95 / pdb_00006d95
Title : Ternary RsAgo Complex with Guide RNA Paired and Target DNA containing A8-A8' Non-Canonical Pair
Authors : Liu, Y.; Esyunina, D.; Olovnikov, I.; Teplova, M.; Patel, D.J.
Deposited on : 2018-04-27
Resolution : 1.85 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

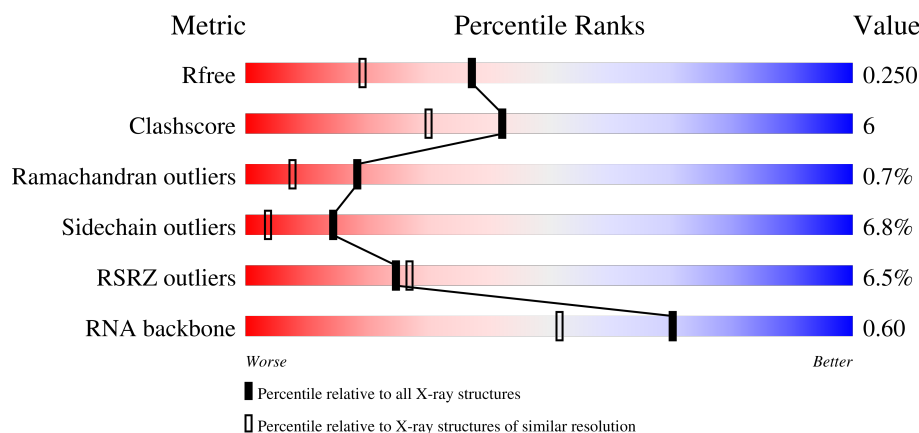
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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	3428 (1.86-1.86)
Clashscore	190562	3579 (1.86-1.86)
Ramachandran outliers	187476	3553 (1.86-1.86)
Sidechain outliers	187428	3553 (1.86-1.86)
RSRZ outliers	180081	3429 (1.86-1.86)
RNA backbone	3983	1030 (2.26-1.46)

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	758	5% 83% 15%
1	F	758	8% 86% 11%
2	B	18	61% 33% 6%
2	H	18	56% 33% 11%

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Mol	Chain	Length	Quality of chain
3	C	24	71%8%21%
3	J	24	67%21%12%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13786 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	0	0
			5761	3673	1023	1049	16			
1	F	755	Total	C	N	O	S	0	0	0
			5725	3649	1013	1047	16			

- 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	B	18	Total	C	N	O	P	0	0	0
			386	172	70	126	18			
2	H	18	Total	C	N	O	P	0	2	0
			429	191	78	140	20			

- Molecule 3 is a DNA chain called DNA 24-Mer.

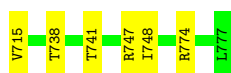
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	19	Total	C	N	O	P	0	0	0
			387	184	71	113	19			
3	J	21	Total	C	N	O	P	0	0	0
			429	204	78	126	21			

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Mg	0	0
			2	2		
4	H	2	Total	Mg	0	0
			2	2		

- Molecule 5 is water.

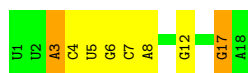
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	277	Total 277	O 277	0	0
5	B	48	Total 48	O 48	0	0
5	F	218	Total 218	O 218	0	0
5	C	43	Total 43	O 43	0	0
5	H	42	Total 42	O 42	0	0
5	J	37	Total 37	O 37	0	0



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



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



- Molecule 3: DNA 24-Mer



- Molecule 3: DNA 24-Mer



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.20Å 119.98Å 118.69Å 90.00° 95.51° 90.00°	Depositor
Resolution (Å)	37.44 – 1.85 37.44 – 1.85	Depositor EDS
% Data completeness (in resolution range)	98.3 (37.44-1.85) 98.3 (37.44-1.85)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 1.85Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.210 , 0.250 0.214 , 0.250	Depositor DCC
R_{free} test set	7974 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	38.8	Xtriage
Anisotropy	0.350	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 34.1	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	13786	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.24% 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	0.87	4/5890 (0.1%)	0.99	9/8016 (0.1%)
1	F	0.77	1/5852 (0.0%)	0.95	5/7966 (0.1%)
2	B	0.63	0/431	1.00	2/668 (0.3%)
2	H	0.48	0/479	0.90	2/743 (0.3%)
3	C	0.45	0/433	0.86	0/665
3	J	0.48	0/480	0.88	0/738
All	All	0.79	5/13565 (0.0%)	0.96	18/18796 (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	2

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	205	GLN	C-N	6.41	1.41	1.33
1	A	206	PRO	N-CD	5.30	1.55	1.47
1	F	117	ARG	C-O	-5.29	1.21	1.23
1	A	545	VAL	C-O	5.07	1.29	1.23
1	A	311	VAL	N-CA	-5.04	1.40	1.46

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	567	GLU	N-CA-C	8.75	123.26	112.58
1	A	630	ILE	CB-CA-C	-8.56	97.86	110.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	630	ILE	CB-CA-C	-7.08	100.05	110.81
1	A	510	THR	CA-C-N	7.01	126.83	119.05
1	A	510	THR	C-N-CA	7.01	126.83	119.05

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	F	565	GLU	Peptide
1	F	566	CYS	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	5761	0	5592	81	0
1	F	5725	0	5533	61	0
2	B	386	0	195	3	0
2	H	429	0	217	9	0
3	C	387	0	214	4	0
3	J	429	0	237	7	0
4	B	2	0	0	0	0
4	H	2	0	0	0	0
5	A	277	0	0	5	0
5	B	48	0	0	3	0
5	C	43	0	0	0	0
5	F	218	0	0	9	0
5	H	42	0	0	3	0
5	J	37	0	0	0	0
All	All	13786	0	11988	159	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 6.

The worst 5 of 159 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:F:528:MET:HE1	1:F:571:TYR:HB2	1.34	1.03
1:A:205:GLN:HG2	1:A:208:GLU:CB	2.00	0.91
1:F:67:LYS:CB	5:F:980:HOH:O	2.21	0.88
1:F:513:THR:HG23	1:F:557:TYR:O	1.75	0.87
3:J:5:DC:H2'	3:J:5:DC:O2	1.78	0.83

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	756/758 (100%)	728 (96%)	23 (3%)	5 (1%)	18	8
1	F	751/758 (99%)	715 (95%)	31 (4%)	5 (1%)	18	8
All	All	1507/1516 (99%)	1443 (96%)	54 (4%)	10 (1%)	18	8

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	220	ASP
1	A	567	GLU
1	F	219	SER
1	F	259	ASN
1	F	300	ASP

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	580/642 (90%)	538 (93%)	42 (7%)	13	3
1	F	575/642 (90%)	538 (94%)	37 (6%)	16	4
All	All	1155/1284 (90%)	1076 (93%)	79 (7%)	14	4

5 of 79 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	406	LYS
1	F	630	ILE
1	F	452	ASP
1	F	562	VAL
1	F	683	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	F	686	ASN
1	F	465	HIS
1	F	235	ASN
1	A	591	ASN
1	F	359	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	17/18 (94%)	2 (11%)	0
2	H	15/18 (83%)	1 (6%)	0
All	All	32/36 (88%)	3 (9%)	0

All (3) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	12	G
2	B	18	A
2	H	12	G

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 4 ligands modelled in this entry, 4 are 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 [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	758/758 (100%)	0.56	40 (5%)	32	35	23, 45, 74, 120	2 (0%)
1	F	755/758 (99%)	0.72	64 (8%)	16	17	27, 47, 84, 108	1 (0%)
2	B	18/18 (100%)	-0.39	0	100	100	30, 33, 56, 86	0
2	H	18/18 (100%)	-0.25	0	100	100	20, 35, 51, 57	2 (11%)
3	C	19/24 (79%)	-0.17	0	100	100	28, 38, 101, 117	0
3	J	21/24 (87%)	-0.09	0	100	100	30, 45, 91, 103	0
All	All	1589/1600 (99%)	0.60	104 (6%)	25	27	20, 45, 79, 120	5 (0%)

The worst 5 of 104 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	188	LEU	5.1
1	A	551	PHE	4.8
1	F	255	LEU	4.7
1	A	205	GLN	4.5
1	F	551	PHE	4.0

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(Å ²)	Q<0.9
4	MG	H	102	1/1	0.97	0.05	48,48,48,48	0
4	MG	B	102	1/1	0.99	0.06	39,39,39,39	0
4	MG	H	101	1/1	1.00	0.02	30,30,30,30	0
4	MG	B	101	1/1	1.00	0.03	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.