



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 04:53 PM UTC

PDB ID : 2F8T / pdb_00002f8t
Title : Crystal structure of Aa-Ago with externally-bound siRNA
Authors : Yuan, Y.R.; Chen, H.Y.; Patel, D.J.
Deposited on : 2005-12-03
Resolution : 3.10 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

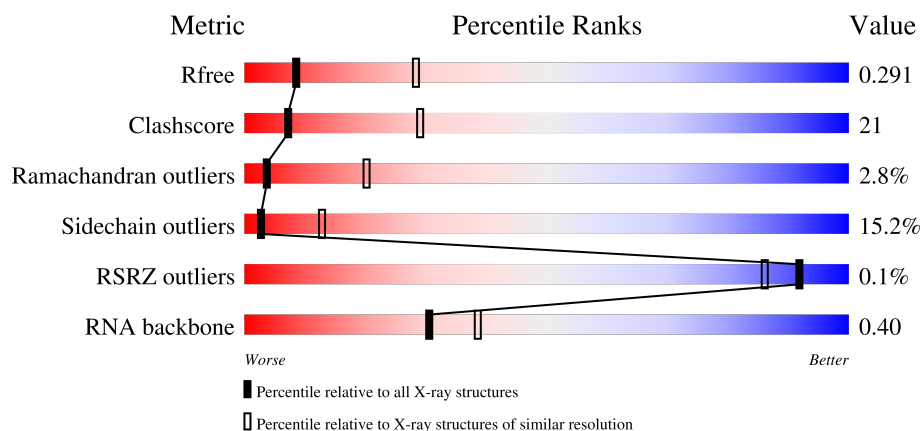
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)
RNA backbone	3983	1022 (3.32-2.88)

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	C	26	31% 35% 19% 15%
1	D	26	15% 38% 35% 8%
2	A	706	54% 37% 8%
2	B	706	49% 40% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12841 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 RNA chain called 26-MER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	26	Total	C	N	O	P	0	0	0
			551	246	94	185	26			
1	D	24	Total	C	N	O	P	0	0	0
			511	228	90	169	24			

- Molecule 2 is a protein called Argonaute protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	704	Total	C	N	O	S	0	0	0
			5864	3818	993	1042	11			
2	B	704	Total	C	N	O	S	0	0	0
			5864	3818	993	1042	11			

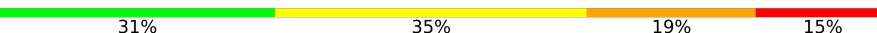
- Molecule 3 is water.

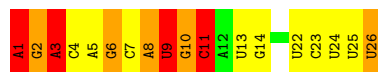
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	C	2	Total	O	0	0
			2	2		
3	D	2	Total	O	0	0
			2	2		
3	A	24	Total	O	0	0
			24	24		
3	B	23	Total	O	0	0
			23	23		

3 Residue-property plots

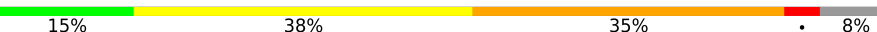
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 26-MER

Chain C: 



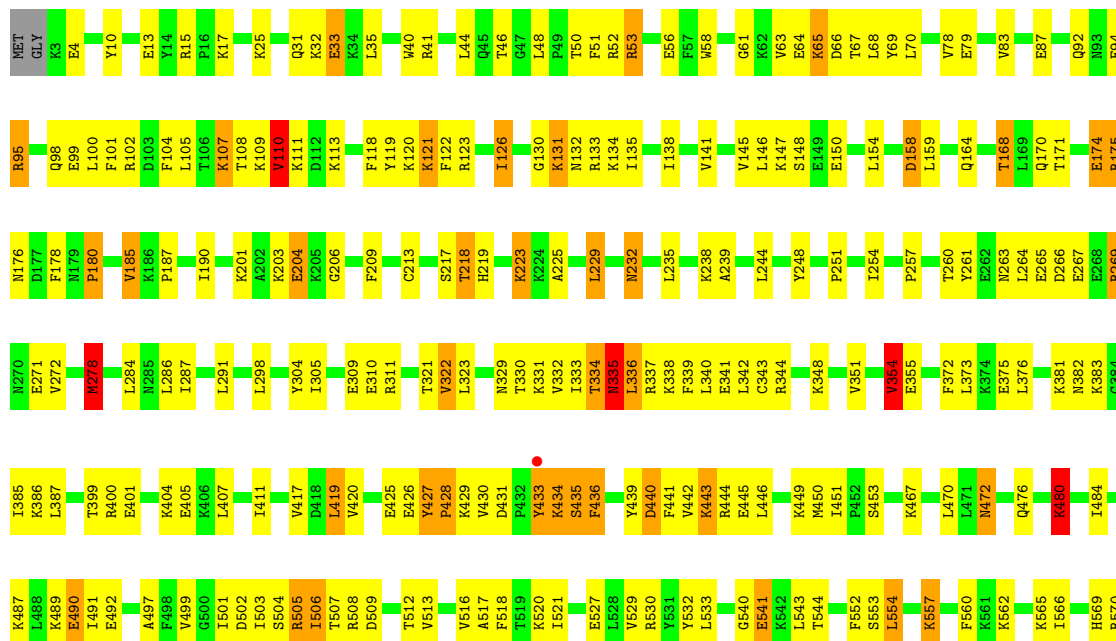
• Molecule 1: 26-MER

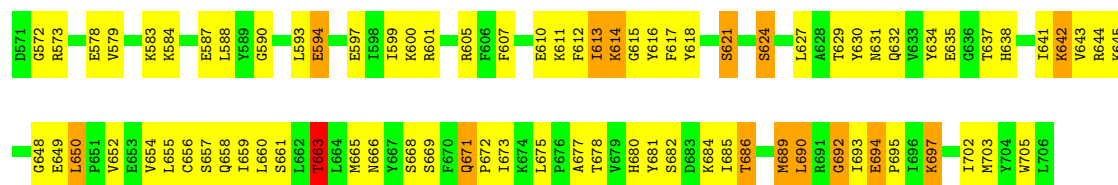
Chain D: 



• Molecule 2: Argonaute protein

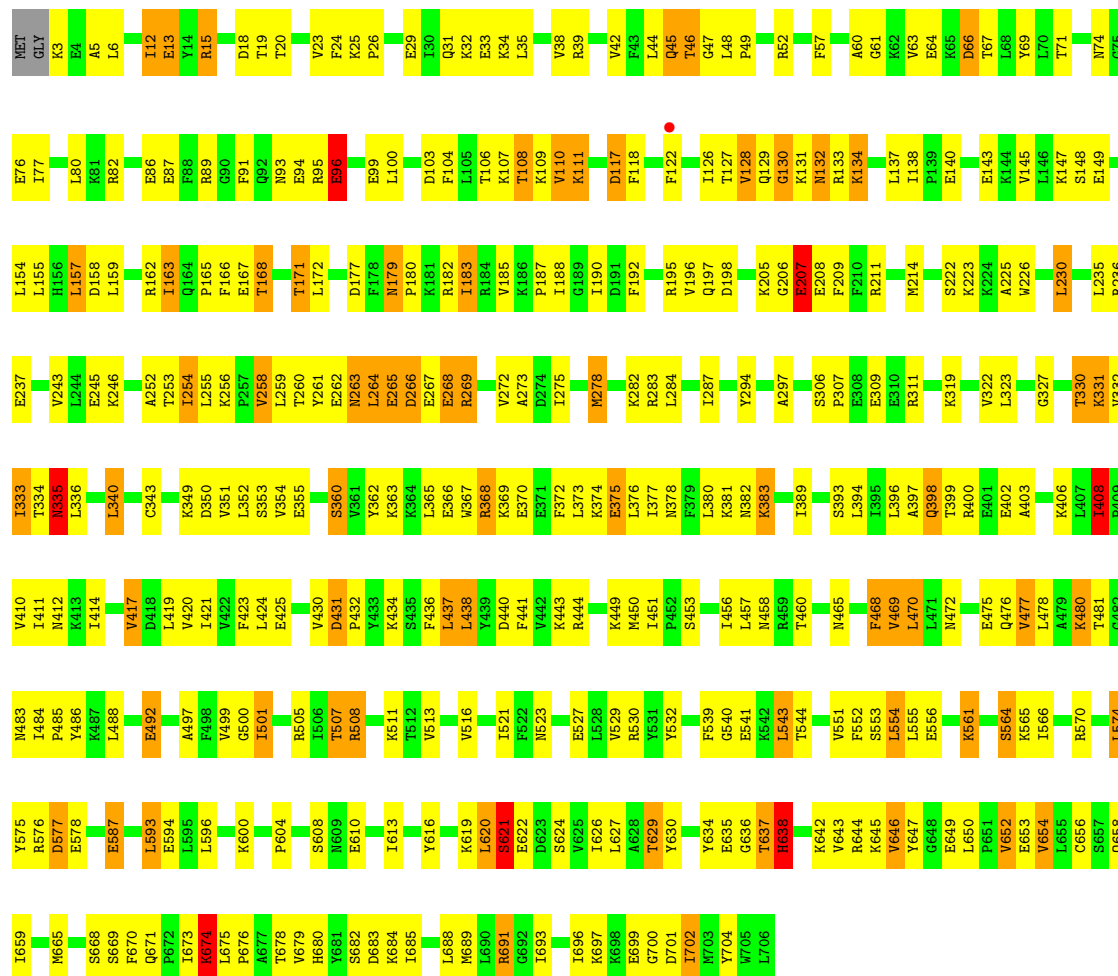
Chain A: 





• Molecule 2: Argonaute protein

Chain B: 49% 40% 10% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.20Å 118.08Å 98.50Å 90.00° 99.57° 90.00°	Depositor
Resolution (Å)	40.00 – 3.10 40.00 – 3.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.00-3.10) 99.9 (40.00-3.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.71 (at 3.12Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.204 , 0.297 0.201 , 0.291	Depositor DCC
R_{free} test set	1672 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	58.5	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 54.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12841	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.23% 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 i

5.1 Standard geometry i

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	C	1.21	6/614 (1.0%)	1.39	5/952 (0.5%)
1	D	0.85	1/570 (0.2%)	1.39	7/884 (0.8%)
2	A	0.88	4/5979 (0.1%)	1.13	16/8027 (0.2%)
2	B	0.85	1/5979 (0.0%)	1.12	22/8027 (0.3%)
All	All	0.88	12/13142 (0.1%)	1.15	50/17890 (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	C	0	1
2	A	0	1
2	B	0	4
All	All	0	6

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	642	LYS	CG-CD	8.12	1.76	1.52
1	C	2	G	N9-C8	7.60	1.53	1.37
1	C	2	G	C1'-N9	6.07	1.57	1.48
1	D	15	C	C1'-N1	5.89	1.57	1.48
1	C	2	G	C5-C4	5.78	1.50	1.38

The worst 5 of 50 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	638	HIS	N-CA-C	7.77	119.95	108.60
1	C	13	U	P-O3'-C3'	-7.65	108.72	120.20
2	B	620	LEU	N-CA-C	-7.53	101.15	110.41
2	A	344	ARG	CA-C-N	7.46	127.42	120.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	344	ARG	C-N-CA	7.46	127.42	120.03

There are no chirality outliers.

5 of 6 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	335	ASN	Peptide
2	B	507	THR	Peptide
2	B	576	ARG	Peptide
2	B	621	SER	Peptide
1	C	1	A	Sidechain

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	C	551	0	279	24	0
1	D	511	0	259	31	0
2	A	5864	0	6128	221	0
2	B	5864	0	6128	266	0
3	A	24	0	0	3	0
3	B	23	0	0	1	0
3	C	2	0	0	0	0
3	D	2	0	0	1	0
All	All	12841	0	12794	529	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 21.

The worst 5 of 529 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)
2:A:642:LYS:CD	2:A:642:LYS:CG	1.76	1.61
2:B:532:TYR:CZ	2:B:554:LEU:HD21	1.70	1.27
2:B:532:TYR:CE1	2:B:554:LEU:HD21	1.78	1.18

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:66:ASP:HB3	3:A:708:HOH:O	1.51	1.11
2:B:122:PHE:CZ	2:B:275:ILE:HG21	1.87	1.09

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	
2	A	702/706 (99%)	605 (86%)	77 (11%)	20 (3%)	4	20
2	B	702/706 (99%)	605 (86%)	78 (11%)	19 (3%)	4	20
All	All	1404/1412 (99%)	1210 (86%)	155 (11%)	39 (3%)	4	20

5 of 39 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	434	LYS
2	A	435	SER
2	B	130	GLY
2	B	132	ASN
2	B	265	GLU

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	
2	A	646/647 (100%)	553 (86%)	93 (14%)	3	14
2	B	646/647 (100%)	543 (84%)	103 (16%)	2	12
All	All	1292/1294 (100%)	1096 (85%)	196 (15%)	3	13

5 of 196 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	143	GLU
2	B	351	VAL
2	B	163	ILE
2	B	258	VAL
2	B	383	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 16 such sidechains are listed below:

Mol	Chain	Res	Type
2	B	523	ASN
2	B	335	ASN
2	B	93	ASN
2	B	263	ASN
2	A	680	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	C	25/26 (96%)	6 (24%)	0
1	D	23/26 (88%)	6 (26%)	1 (4%)
All	All	48/52 (92%)	12 (25%)	1 (2%)

5 of 12 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	C	3	A
1	C	6	G
1	C	9	U
1	C	10	G
1	C	11	C

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	D	23	C

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å²)	Q<0.9	
1	C	26/26 (100%)	-0.18	0	100	100	31, 58, 79, 82	0
1	D	24/26 (92%)	-0.26	0	100	100	26, 59, 89, 91	0
2	A	704/706 (99%)	-0.37	1 (0%)	92	86	27, 54, 82, 96	0
2	B	704/706 (99%)	-0.33	1 (0%)	92	86	30, 55, 78, 93	0
All	All	1458/1464 (99%)	-0.35	2 (0%)	92	86	26, 54, 80, 96	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	433	TYR	2.3
2	B	122	PHE	2.3

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.