



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:10 PM UTC

PDB ID : 5QYZ / pdb_00005qyz
Title : PanDDA analysis group deposition – Auto-refined data of Aar2/RNaseH for ground state model 15
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Deposited on : 2020-02-12
Resolution : 1.37 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
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

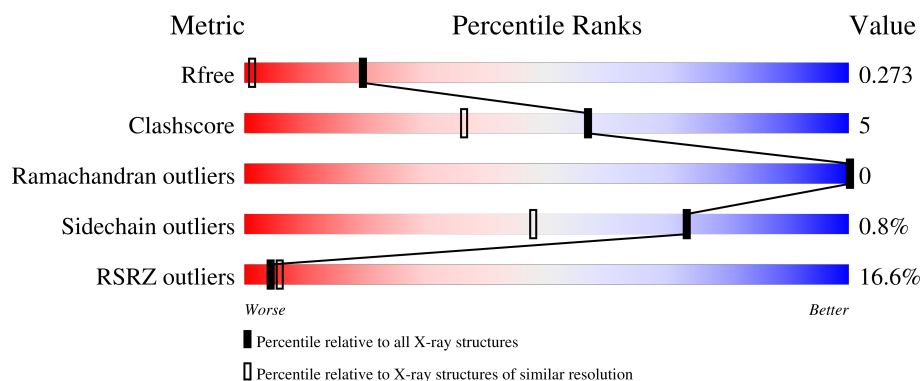
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.37 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	4403 (1.40-1.36)
Clashscore	190562	4528 (1.40-1.36)
Ramachandran outliers	187476	4459 (1.40-1.36)
Sidechain outliers	187428	4458 (1.40-1.36)
RSRZ outliers	180081	4399 (1.40-1.36)

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	258	15% 84% 6% • 8%
2	B	308	16% 87% 10% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4772 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 Pre-mRNA-splicing factor 8.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	237	2002	1283	335	372	12	0	12	0

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1833	GLY	-	expression tag	UNP P33334
A	1834	ALA	-	expression tag	UNP P33334
A	1835	MET	-	expression tag	UNP P33334

- Molecule 2 is a protein called A1 cistron-splicing factor AAR2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	300	2580	1654	421	485	20	0	9	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	GLY	-	expression tag	UNP P32357
B	-2	ALA	-	expression tag	UNP P32357
B	-1	MET	-	expression tag	UNP P32357
B	0	ALA	-	expression tag	UNP P32357
B	166	SER	-	linker	UNP P32357
B	167	SER	-	linker	UNP P32357
B	168	SER	-	linker	UNP P32357
B	169	SER	-	linker	UNP P32357
B	170	SER	-	linker	UNP P32357

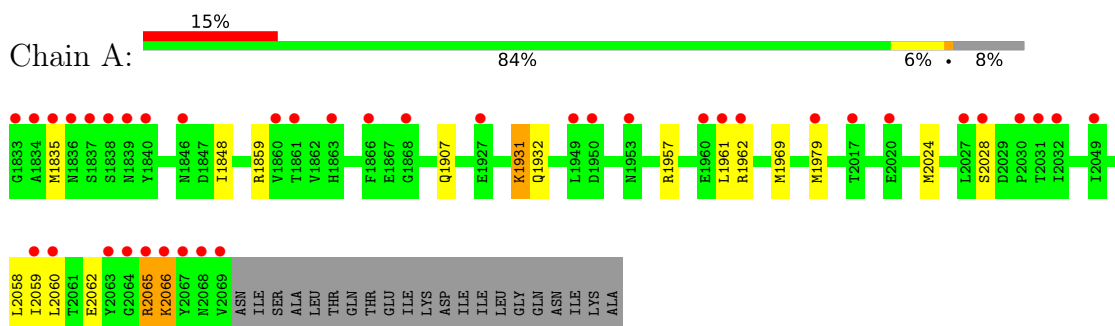
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	93	Total 93	O 93	0	0
3	B	97	Total 97	O 97	0	0

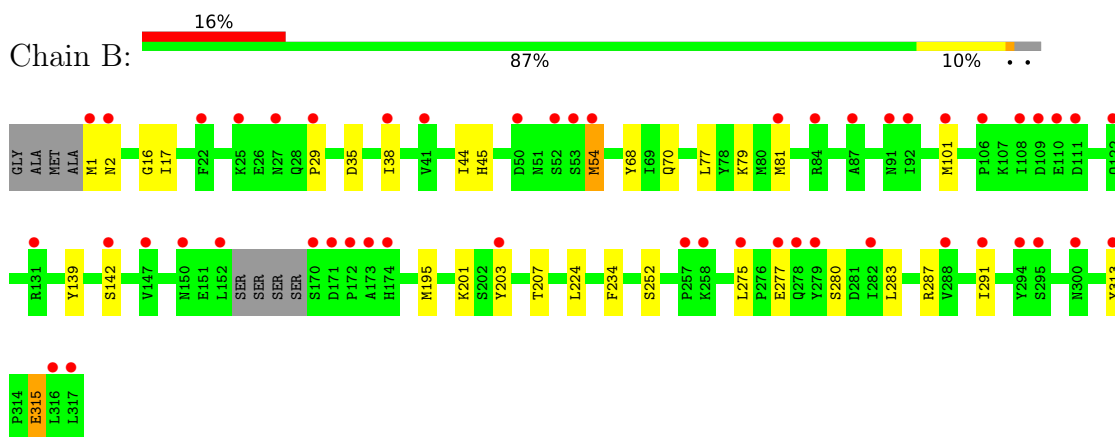
3 Residue-property plots

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: Pre-mRNA-splicing factor 8



• Molecule 2: A1 cistron-splicing factor AAR2



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	88.22Å 81.79Å 93.82Å 90.00° 108.46° 90.00°	Depositor
Resolution (Å)	44.60 – 1.37 44.60 – 1.37	Depositor EDS
% Data completeness (in resolution range)	99.5 (44.60-1.37) 99.5 (44.60-1.37)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.96 (at 1.37Å)	Xtriage
Refinement program	REFMAC 5.8.0238, PHENIX 1.16.3549	Depositor
R, R_{free}	0.223 , 0.224 (Not available) , 0.273	Depositor DCC
R_{free} test set	6474 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	26.5	Xtriage
Anisotropy	0.102	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 26.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4772	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.44% 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

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	1.00	1/2049 (0.0%)	1.17	2/2775 (0.1%)
2	B	1.05	2/2651 (0.1%)	1.20	0/3581
All	All	1.03	3/4700 (0.1%)	1.19	2/6356 (0.0%)

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
2	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	38	ILE	C-O	7.55	1.33	1.24
1	A	2059	ILE	C-O	-6.48	1.16	1.24
2	B	139	TYR	C-O	5.33	1.30	1.24

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2065	ARG	CB-CG-CD	6.50	126.24	111.30
1	A	2066	LYS	CB-CA-C	-5.18	102.74	110.88

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	54[A]	MET	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	2002	0	2029	18	0
2	B	2580	0	2450	25	0
3	A	93	0	0	4	0
3	B	97	0	0	6	0
All	All	4772	0	4479	42	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 5.

All (42) 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:B:101:MET:HG3	3:B:490:HOH:O	1.05	1.20
2:B:54[A]:MET:HE2	3:B:495:HOH:O	1.63	0.99
2:B:101:MET:CG	3:B:490:HOH:O	1.78	0.87
2:B:1:MET:SD	3:B:486:HOH:O	2.33	0.84
1:A:2028:SER:CB	3:A:2186:HOH:O	2.26	0.83
1:A:2062:GLU:O	1:A:2066:LYS:HG2	1.84	0.78
2:B:1:MET:HB3	2:B:35:ASP:HA	1.67	0.76
2:B:101:MET:SD	3:B:490:HOH:O	2.34	0.76
1:A:2028:SER:HA	3:A:2186:HOH:O	1.89	0.72
1:A:1848:ILE:H	1:A:1931[A]:LYS:HZ2	1.40	0.69
1:A:1835:MET:HE1	1:A:1961:LEU:HD12	1.75	0.67
1:A:2028:SER:CA	3:A:2186:HOH:O	2.44	0.64
2:B:280:SER:HB3	2:B:313:TYR:CE1	2.33	0.63
1:A:2062:GLU:HB3	1:A:2066:LYS:HE3	1.83	0.60
2:B:277:GLU:CD	2:B:277:GLU:H	2.10	0.60
2:B:287:ARG:O	2:B:291:ILE:HD13	2.04	0.57
1:A:2065:ARG:HH11	1:A:2065:ARG:HG2	1.69	0.57
1:A:1907:GLN:HG2	2:B:195:MET:HE3	1.90	0.53
1:A:1969:MET:HA	1:A:1969:MET:HE2	1.91	0.52
1:A:1979[B]:MET:HA	1:A:1979[B]:MET:HE2	1.90	0.52
2:B:275:LEU:CD2	2:B:283:LEU:HD13	2.40	0.51
1:A:2058:LEU:C	1:A:2058:LEU:HD23	2.35	0.51
2:B:77:LEU:HD21	2:B:79:LYS:HE3	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:280:SER:CB	2:B:313:TYR:CE1	2.97	0.48
2:B:142:SER:OG	2:B:234:PHE:O	2.26	0.47
1:A:1962:ARG:HD3	1:A:1962:ARG:H	1.80	0.46
2:B:54[A]:MET:HG3	3:B:484:HOH:O	2.15	0.46
2:B:315:GLU:H	2:B:315:GLU:CD	2.24	0.46
2:B:203[A]:TYR:CZ	2:B:207:THR:HG21	2.51	0.46
1:A:2028:SER:HB3	3:A:2186:HOH:O	2.04	0.45
2:B:17:ILE:HD13	2:B:44[B]:ILE:CG1	2.47	0.45
2:B:16:GLY:HA3	2:B:45:HIS:CE1	2.53	0.44
2:B:201:LYS:HG3	2:B:252:SER:HB2	2.00	0.44
2:B:224:LEU:C	2:B:224:LEU:HD23	2.44	0.43
2:B:275:LEU:HD23	2:B:275:LEU:HA	1.93	0.42
1:A:1859:ARG:HH12	1:A:1979[A]:MET:CE	2.32	0.42
2:B:29:PRO:HB2	2:B:101:MET:SD	2.61	0.41
2:B:68:TYR:CE2	2:B:81:MET:HE3	2.56	0.41
1:A:1932:GLN:OE1	1:A:1957:ARG:NH1	2.47	0.41
2:B:70:GLN:HB3	2:B:81:MET:HE2	2.02	0.41
1:A:2024[A]:MET:HE3	1:A:2024[A]:MET:HB3	1.95	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	248/258 (96%)	244 (98%)	4 (2%)	0	100	100
2	B	306/308 (99%)	295 (96%)	11 (4%)	0	100	100
All	All	554/566 (98%)	539 (97%)	15 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	226/233 (97%)	223 (99%)	3 (1%)	61	29
2	B	287/284 (101%)	285 (99%)	2 (1%)	76	52
All	All	513/517 (99%)	508 (99%)	5 (1%)	73	40

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1931[A]	LYS
1	A	1931[B]	LYS
1	A	2060	LEU
2	B	2	ASN
2	B	315	GLU

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

Mol	Chain	Res	Type
2	B	2	ASN
2	B	47	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	237/258 (91%)	0.97	39 (16%)	4 6	8, 24, 46, 71	12 (5%)
2	B	300/308 (97%)	1.16	50 (16%)	4 6	9, 27, 50, 91	9 (3%)
All	All	537/566 (94%)	1.08	89 (16%)	4 6	8, 25, 47, 91	21 (3%)

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2069	VAL	8.2
2	B	22	PHE	7.9
2	B	152	LEU	6.8
2	B	1	MET	6.8
1	A	1834	ALA	6.3
2	B	170	SER	6.2
2	B	54[A]	MET	5.7
2	B	173	ALA	5.7
1	A	2065	ARG	5.6
2	B	279	TYR	5.3
2	B	172	PRO	5.1
2	B	203[A]	TYR	4.9
2	B	108	ILE	4.7
2	B	101	MET	4.7
2	B	317	LEU	4.6
1	A	2017[A]	THR	4.6
1	A	1835	MET	4.6
1	A	2060	LEU	4.5
1	A	1979[A]	MET	4.4
1	A	1833	GLY	4.4
2	B	41[A]	VAL	4.3
1	A	1837	SER	4.3
2	B	174	HIS	4.3
2	B	316	LEU	4.2

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Mol	Chain	Res	Type	RSRZ
2	B	38	ILE	4.2
1	A	1838	SER	4.1
1	A	1866	PHE	4.0
2	B	282	ILE	3.8
1	A	1840	TYR	3.8
2	B	313	TYR	3.6
1	A	1839	ASN	3.6
1	A	2063	TYR	3.5
1	A	2030	PRO	3.5
1	A	2027	LEU	3.5
1	A	2068	ASN	3.4
1	A	1836	ASN	3.4
2	B	50	ASP	3.2
2	B	92	ILE	3.1
2	B	147	VAL	3.1
1	A	2028	SER	3.0
2	B	275	LEU	3.0
2	B	150	ASN	3.0
2	B	84	ARG	2.9
2	B	258	LYS	2.9
2	B	2	ASN	2.9
2	B	122[A]	GLN	2.8
2	B	106	PRO	2.8
2	B	171	ASP	2.8
2	B	295	SER	2.8
1	A	1846	ASN	2.7
2	B	53	SER	2.7
2	B	277	GLU	2.7
1	A	1962	ARG	2.7
2	B	87	ALA	2.6
2	B	142	SER	2.6
2	B	110	GLU	2.6
2	B	278	GLN	2.5
1	A	1961	LEU	2.5
1	A	2066	LYS	2.5
1	A	2031	THR	2.5
2	B	257	PRO	2.5
1	A	1960[A]	GLU	2.4
2	B	109	ASP	2.4
2	B	131	ARG	2.3
2	B	81	MET	2.3
1	A	2064	GLY	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	1860	VAL	2.3
2	B	27	ASN	2.2
2	B	111	ASP	2.2
2	B	291	ILE	2.2
1	A	2067	TYR	2.2
1	A	1950	ASP	2.2
1	A	2020	GLU	2.1
1	A	2032	ILE	2.1
1	A	1861	THR	2.1
1	A	2049	ILE	2.1
1	A	2059	ILE	2.1
2	B	294	TYR	2.1
2	B	288	VAL	2.1
2	B	29	PRO	2.1
1	A	1953	ASN	2.1
1	A	1949	LEU	2.1
1	A	1927	GLU	2.1
2	B	25	LYS	2.1
2	B	52	SER	2.0
2	B	91	ASN	2.0
2	B	300	ASN	2.0
1	A	1868	GLY	2.0
1	A	1863	HIS	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.