Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 08:36 PM UTC

PDB ID	:	5QZX / pdb_00005qzx
Title	:	PanDDA analysis group deposition – Auto-refined data of Aar2/RNaseH for ground state model 48
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Deposited on	:	2020-02-12
Resolution	:	1.55 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:

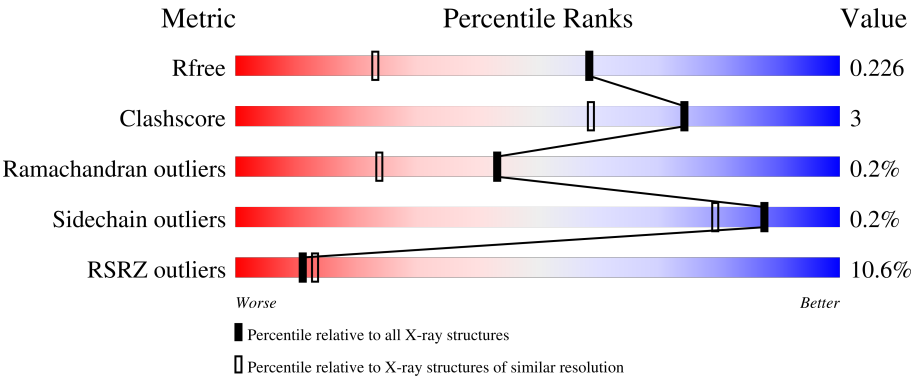
MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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	2145 (1.56-1.56)
Clashscore	190562	2189 (1.56-1.56)
Ramachandran outliers	187476	2153 (1.56-1.56)
Sidechain outliers	187428	2150 (1.56-1.56)
RSRZ outliers	180081	2146 (1.56-1.56)

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 <=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	9% 87% 5% 8%
2	B	308	11% 88% 9% .

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4697 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
1	A	237	Total	C	N	O	S	0	12	0
			2002	1283	335	372	12			

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
2	B	300	Total	C	N	O	S	0	9	0
			2580	1654	421	485	20			

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

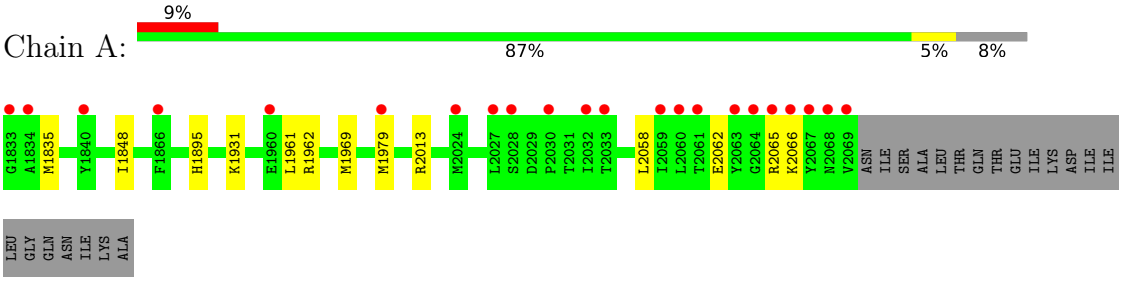
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	57	Total O 57 57	0	0
3	B	58	Total O 58 58	0	0

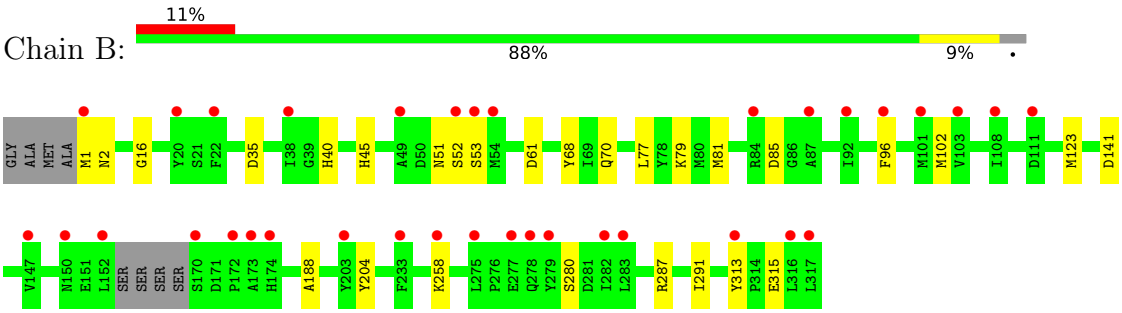
3 Residue-property plots (i)

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: Pre-mRNA-splicing factor 8



• Molecule 2: A1 cistron-splicing factor AAR2



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	89.22Å 81.75Å 93.19Å 90.00° 108.28° 90.00°	Depositor
Resolution (Å)	44.74 – 1.55 44.74 – 1.55	Depositor EDS
% Data completeness (in resolution range)	99.2 (44.74-1.55) 99.8 (44.74-1.55)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.96 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.8.0238, PHENIX 1.16.3549	Depositor
R, R_{free}	0.219 , 0.221 0.230 , 0.226	Depositor DCC
R_{free} test set	4635 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	34.7	Xtriage
Anisotropy	0.206	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 28.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	4697	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 5.40% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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	1.06	1/2049 (0.0%)	1.22	0/2775
2	B	1.03	0/2651	1.26	4/3581 (0.1%)
All	All	1.04	1/4700 (0.0%)	1.24	4/6356 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1895	HIS	CE1-NE2	5.30	1.37	1.32

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	85	ASP	CA-C-N	5.46	126.16	119.99
2	B	85	ASP	C-N-CA	5.46	126.16	119.99
2	B	141	ASP	CA-C-N	5.11	127.37	120.38
2	B	141	ASP	C-N-CA	5.11	127.37	120.38

There are no chirality outliers.

There are no planarity outliers.

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	9	0
2	B	2580	0	2450	19	0
3	A	57	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	58	0	0	4	0
All	All	4697	0	4479	28	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 3.

All (28) 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:70:GLN:HB3	2:B:81:MET:HE2	1.67	0.75
1:A:1979[B]:MET:HA	1:A:1979[B]:MET:HE2	1.70	0.72
1:A:1835:MET:HE1	1:A:1961:LEU:HD12	1.72	0.71
2:B:280:SER:HB3	2:B:313:TYR:CE1	2.27	0.69
2:B:1:MET:HB3	2:B:35:ASP:HA	1.76	0.67
2:B:96:PHE:HB2	2:B:102:MET:HE3	1.78	0.65
2:B:70:GLN:HB3	2:B:81:MET:CE	2.32	0.59
2:B:258:LYS:HD2	2:B:258:LYS:H	1.67	0.58
2:B:287:ARG:O	2:B:291:ILE:HD13	2.05	0.56
1:A:1969:MET:HA	1:A:1969:MET:HE2	1.89	0.55
1:A:1848:ILE:H	1:A:1931[A]:LYS:HZ2	1.55	0.55
2:B:2:ASN:HB2	3:B:423:HOH:O	2.08	0.54
1:A:2058:LEU:C	1:A:2058:LEU:HD23	2.34	0.53
2:B:96:PHE:CB	2:B:102:MET:HE3	2.39	0.52
1:A:2062:GLU:HB3	1:A:2066:LYS:HE3	1.92	0.51
2:B:51:ASN:C	2:B:53:SER:H	2.18	0.51
2:B:68:TYR:CE2	2:B:81:MET:HE3	2.46	0.50
2:B:77:LEU:HD21	2:B:79:LYS:HE3	1.95	0.49
1:A:2062:GLU:O	1:A:2066:LYS:HG2	2.13	0.48
2:B:51:ASN:ND2	3:B:401:HOH:O	2.46	0.48
1:A:2065:ARG:O	1:A:2066:LYS:C	2.56	0.47
1:A:1962:ARG:O	1:A:2013:ARG:NH1	2.49	0.46
2:B:188:ALA:HA	2:B:204:TYR:CD1	2.51	0.45
2:B:40:HIS:HD2	3:B:427:HOH:O	1.99	0.45
2:B:280:SER:CB	2:B:313:TYR:CE1	3.01	0.42
2:B:61:ASP:HB2	2:B:123:MET:HG2	2.02	0.42
2:B:16:GLY:HA3	2:B:45:HIS:CE1	2.54	0.42
2:B:1:MET:N	3:B:402:HOH:O	2.52	0.41

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%)	242 (98%)	6 (2%)	0	100 100
2	B	306/308 (99%)	296 (97%)	9 (3%)	1 (0%)	36 18
All	All	554/566 (98%)	538 (97%)	15 (3%)	1 (0%)	43 24

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	52	SER

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	226/233 (97%)	226 (100%)	0	100 100
2	B	287/284 (101%)	286 (100%)	1 (0%)	86 77
All	All	513/517 (99%)	512 (100%)	1 (0%)	87 79

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
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	40	HIS
2	B	136	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, 95^{th} 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%)	0.73	22 (9%)	14	16	12, 35, 68, 90	12 (5%)
2	B	300/308 (97%)	0.86	35 (11%)	9	11	13, 38, 73, 96	9 (3%)
All	All	537/566 (94%)	0.80	57 (10%)	11	13	12, 37, 71, 96	21 (3%)

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	152	LEU	6.9
2	B	54[A]	MET	6.4
1	A	2069	VAL	5.7
2	B	22	PHE	5.6
1	A	1979[A]	MET	5.1
1	A	2060	LEU	4.7
2	B	170	SER	4.7
2	B	316	LEU	4.6
2	B	1	MET	4.4
2	B	101	MET	4.3
1	A	2068	ASN	4.2
2	B	174	HIS	4.1
1	A	2064	GLY	4.1
1	A	2065	ARG	3.8
2	B	53	SER	3.8
1	A	2063	TYR	3.8
2	B	279	TYR	3.8
1	A	2067	TYR	3.8
1	A	1866	PHE	3.6
2	B	203[A]	TYR	3.6
1	A	1833	GLY	3.6
2	B	173	ALA	3.6
2	B	52	SER	3.3
2	B	172	PRO	3.3

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Mol	Chain	Res	Type	RSRZ
2	B	313	TYR	3.2
2	B	317	LEU	3.2
2	B	49	ALA	3.1
1	A	2027	LEU	3.1
2	B	108	ILE	3.1
1	A	2061	THR	3.0
2	B	84	ARG	3.0
1	A	2030	PRO	2.9
1	A	2066	LYS	2.9
2	B	147	VAL	2.9
1	A	1834	ALA	2.8
2	B	38	ILE	2.8
2	B	150	ASN	2.7
2	B	283	LEU	2.7
2	B	282	ILE	2.7
2	B	277	GLU	2.6
1	A	2032	ILE	2.5
1	A	1840	TYR	2.4
2	B	258	LYS	2.4
2	B	87	ALA	2.4
2	B	275	LEU	2.3
2	B	96	PHE	2.3
2	B	278	GLN	2.3
1	A	2059	ILE	2.3
1	A	2028	SER	2.2
2	B	20	TYR	2.2
2	B	111	ASP	2.2
2	B	233	PHE	2.2
2	B	92	ILE	2.1
1	A	1960[A]	GLU	2.0
1	A	2033	THR	2.0
2	B	103	VAL	2.0
1	A	2024[A]	MET	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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.