



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 2, 2021 – 10:37 AM EST

PDB ID : 5R0V
Title : PanDDA analysis group deposition – Auto-refined data of Aar2/RNaseH for ground state model 09, DMSO-free
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Deposited on : 2020-02-12
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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.17.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.17.1

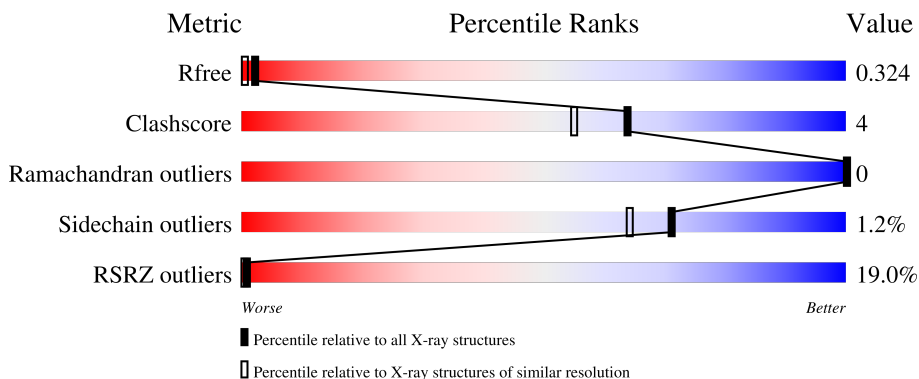
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.81 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}	130704	7484 (1.84-1.80)
Clashscore	141614	8401 (1.84-1.80)
Ramachandran outliers	138981	8290 (1.84-1.80)
Sidechain outliers	138945	8290 (1.84-1.80)
RSRZ outliers	127900	7371 (1.84-1.80)

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	
2	B	308	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4673 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	2568	1645	420	483	20	0	9	0

There are 20 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	LEU	conflict	UNP P32357
B	167	SER	LYS	conflict	UNP P32357
B	170	SER	LEU	conflict	UNP P32357
B	?	-	GLN	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357
B	?	-	ALA	deletion	UNP P32357
B	?	-	GLY	deletion	UNP P32357
B	?	-	SER	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357

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Chain	Residue	Modelled	Actual	Comment	Reference
B	?	-	MET	deletion	UNP P32357
B	?	-	GLU	deletion	UNP P32357
B	?	-	ALA	deletion	UNP P32357
B	?	-	LYS	deletion	UNP P32357
B	?	-	ASN	deletion	UNP P32357
B	?	-	GLU	deletion	UNP P32357
B	?	-	ASP	deletion	UNP P32357

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	63	Total O 63 63	0	0
3	B	40	Total O 40 40	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	88.79Å 82.18Å 92.92Å 90.00° 108.16° 90.00°	Depositor
Resolution (Å)	22.68 – 1.81 44.70 – 1.81	Depositor EDS
% Data completeness (in resolution range)	99.4 (22.68-1.81) 99.4 (44.70-1.81)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.98 (at 1.81Å)	Xtrriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.283 , 0.320 0.288 , 0.324	Depositor DCC
R_{free} test set	2100 reflections (3.62%)	wwPDB-VP
Wilson B-factor (Å ²)	43.4	Xtrriage
Anisotropy	0.344	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4673	wwPDB-VP
Average B, all atoms (Å ²)	81.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.69% 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	A	0.79	3/2049 (0.1%)	0.87	5/2775 (0.2%)
2	B	0.74	3/2638 (0.1%)	0.76	1/3563 (0.0%)
All	All	0.76	6/4687 (0.1%)	0.81	6/6338 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1880	PHE	CE2-CZ	-9.19	1.19	1.37
1	A	1880	PHE	CG-CD1	-6.16	1.29	1.38
2	B	115	TYR	CD2-CE2	-5.62	1.30	1.39
2	B	204	TYR	CD2-CE2	-5.59	1.30	1.39
1	A	1878	CYS	CB-SG	-5.49	1.72	1.81
2	B	62	CYS	CB-SG	-5.45	1.73	1.81

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1949	LEU	CB-CG-CD2	-6.29	100.30	111.00
2	B	176	LEU	CB-CG-CD1	-6.19	100.48	111.00
1	A	1854	ASP	CB-CG-OD2	6.05	123.75	118.30
1	A	1928	GLU	OE1-CD-OE2	-5.69	116.47	123.30
1	A	1905	LEU	CB-CG-CD2	-5.41	101.80	111.00
1	A	1997	ASP	CB-CG-OD2	-5.02	113.78	118.30

There are no chirality outliers.

There are no planarity outliers.

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	2002	0	2029	16	0
2	B	2568	0	2441	22	0
3	A	63	0	0	0	0
3	B	40	0	0	3	0
All	All	4673	0	4470	38	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 4.

All (38) 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:HE1	1.70	0.72
1:A:2062:GLU:O	1:A:2066:LYS:HG2	1.92	0.69
1:A:2062:GLU:HB3	1:A:2066:LYS:HE3	1.81	0.62
1:A:1962:ARG:O	1:A:2013:ARG:NH1	2.33	0.61
1:A:2047:GLN:O	1:A:2051:ILE:HG12	2.04	0.57
2:B:1:MET:N	3:B:401:HOH:O	2.40	0.54
1:A:1989:PHE:CD2	1:A:2039:LEU:HD21	2.44	0.53
2:B:214:PHE:O	2:B:215:LYS:HB2	2.09	0.51
2:B:16:GLY:HA3	2:B:45:HIS:CE1	2.47	0.49
1:A:1941:LEU:O	1:A:1945:GLU:HB2	2.12	0.49
2:B:190:ARG:HG3	2:B:203[B]:TYR:CZ	2.48	0.49
2:B:1:MET:HB3	2:B:35:ASP:HA	1.95	0.48
2:B:170:SER:HA	3:B:422:HOH:O	2.13	0.47
2:B:70:GLN:CB	2:B:81:MET:HE1	2.41	0.47
2:B:237:TYR:CE2	2:B:241:LEU:HD11	2.49	0.47
1:A:1875:ILE:HD12	1:A:1979[B]:MET:HE3	1.96	0.47
2:B:258:LYS:HD2	2:B:258:LYS:H	1.78	0.47
2:B:51:ASN:C	2:B:53:SER:H	2.19	0.46
2:B:251:CYS:O	2:B:296:SER:HB2	2.14	0.46
1:A:2058:LEU:C	1:A:2058:LEU:HD23	2.37	0.46
1:A:1843:LEU:HA	1:A:1849:LYS:HD2	1.99	0.45
2:B:17:ILE:O	2:B:17:ILE:HG23	2.16	0.45
2:B:61:ASP:OD1	2:B:61:ASP:C	2.55	0.45
1:A:1974:LEU:O	1:A:1977:VAL:HG12	2.17	0.45
1:A:1967:ALA:O	1:A:1970:SER:HB2	2.17	0.44
2:B:70:GLN:HB3	2:B:81:MET:CE	2.44	0.44
2:B:301:SER:O	2:B:302:LEU:HD23	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:141:ASP:OD1	2:B:144:MET:HG3	2.18	0.43
2:B:69:ILE:HD13	2:B:80:MET:HA	2.01	0.43
1:A:1927:GLU:OE2	1:A:1927:GLU:N	2.37	0.42
1:A:1949:LEU:HA	1:A:1949:LEU:HD23	1.78	0.42
1:A:2064:GLY:O	1:A:2068:ASN:N	2.53	0.42
1:A:1848:ILE:H	1:A:1931[A]:LYS:HZ2	1.67	0.41
1:A:1910:LYS:O	1:A:1913:THR:HB	2.20	0.41
2:B:129:ILE:HA	2:B:177[B]:ASN:HB3	2.01	0.41
2:B:5:PRO:HA	2:B:32:GLY:HA3	2.03	0.41
2:B:54[A]:MET:HE3	3:B:435:HOH:O	2.20	0.40
2:B:242:GLN:O	2:B:246:MET:HG3	2.21	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%)	242 (98%)	6 (2%)	0	100	100
2	B	305/308 (99%)	291 (95%)	14 (5%)	0	100	100
All	All	553/566 (98%)	533 (96%)	20 (4%)	0	100	100

There are no Ramachandran outliers to report.

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%)	221 (98%)	5 (2%)	52	39
2	B	286/284 (101%)	283 (99%)	3 (1%)	76	70
All	All	512/517 (99%)	504 (98%)	8 (2%)	71	53

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1835	MET
1	A	1962	ARG
1	A	1979[A]	MET
1	A	1979[B]	MET
1	A	1979[C]	MET
2	B	17	ILE
2	B	77	LEU
2	B	258	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	47	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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 monosaccharides 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(Å ²)	Q<0.9
1	A	237/258 (91%)	1.14	47 (19%) 1 0	35, 70, 134, 177	0
2	B	300/308 (97%)	1.20	55 (18%) 1 0	39, 75, 143, 207	0
All	All	537/566 (94%)	1.17	102 (18%) 1 0	35, 72, 138, 207	0

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1	MET	10.7
2	B	53	SER	7.7
2	B	52	SER	7.3
2	B	108	ILE	7.3
2	B	96	PHE	5.8
2	B	174	HIS	5.1
1	A	2060	LEU	5.1
2	B	313	TYR	5.1
2	B	109	ASP	5.0
1	A	1979[A]	MET	4.9
1	A	1878	CYS	4.9
2	B	54[A]	MET	4.7
1	A	2027	LEU	4.5
1	A	2066	LYS	4.4
2	B	87	ALA	4.3
1	A	2063	TYR	4.3
2	B	277	GLU	4.3
1	A	1833	GLY	4.1
1	A	1996	LEU	4.1
1	A	1964	PRO	3.9
2	B	22	PHE	3.8
1	A	1988	LEU	3.8
2	B	281	ASP	3.7
2	B	172	PRO	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	1940	MET	3.6
2	B	280	SER	3.6
2	B	150	ASN	3.6
1	A	1840	TYR	3.6
2	B	175	SER	3.4
1	A	1934	ILE	3.4
2	B	111	ASP	3.4
2	B	171	ASP	3.3
1	A	1969	MET	3.2
2	B	279	TYR	3.1
1	A	1838	SER	3.1
2	B	308	ILE	3.1
2	B	86	GLY	3.1
2	B	122[A]	GLN	3.1
2	B	89	PHE	3.1
2	B	101	MET	3.0
2	B	317	LEU	3.0
1	A	1866	PHE	3.0
1	A	2055	MET	3.0
1	A	1965	PHE	2.9
2	B	110	GLU	2.9
2	B	99	ARG	2.9
1	A	1881	THR	2.9
2	B	181	ILE	2.9
1	A	1889	LEU	2.8
1	A	1961	LEU	2.8
2	B	92	ILE	2.8
2	B	121	VAL	2.8
2	B	254	ALA	2.7
2	B	213	ILE	2.7
1	A	1834	ALA	2.7
2	B	73	PRO	2.7
1	A	1905	LEU	2.7
2	B	36	ILE	2.6
1	A	1860	VAL	2.6
1	A	2056[A]	ARG	2.6
1	A	1963	LEU	2.6
2	B	30	PHE	2.6
1	A	2048	TRP	2.6
1	A	2039	LEU	2.5
2	B	189	ILE	2.5
1	A	2012	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	2026	LEU	2.5
2	B	294	TYR	2.5
2	B	170	SER	2.5
1	A	1962	ARG	2.5
2	B	84	ARG	2.5
2	B	29	PRO	2.4
1	A	1954	ILE	2.3
1	A	2044	THR	2.3
1	A	2046	GLU	2.3
1	A	1948	MET	2.3
1	A	2030	PRO	2.3
1	A	2015	LEU	2.3
2	B	316	LEU	2.3
2	B	182	ASN	2.2
1	A	2028	SER	2.2
2	B	193	HIS	2.2
2	B	146	THR	2.2
1	A	1836	ASN	2.2
2	B	117	LEU	2.2
2	B	176	LEU	2.2
1	A	2032	ILE	2.1
1	A	2024[A]	MET	2.1
2	B	203[A]	TYR	2.1
1	A	2034	ILE	2.1
1	A	2022	ALA	2.1
2	B	249	LEU	2.1
1	A	2061	THR	2.1
1	A	2051	ILE	2.1
2	B	91	ASN	2.1
2	B	100	GLN	2.1
2	B	123	MET	2.1
2	B	295	SER	2.1
1	A	1839	ASN	2.1
2	B	312	LYS	2.0
1	A	1967	ALA	2.0
2	B	38	ILE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates [i](#)

There are no monosaccharides 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.