



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 12, 2022 – 07:05 pm BST

PDB ID : 7Q3C
Title : Ribonucleotide Reductase AaR2 protein from Aquifex aeolicus
Authors : Scaletti, E.R.; Rehling, D.; Stenmark, P.
Deposited on : 2021-10-27
Resolution : 2.15 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 : 2.27
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0267
CCP4 : 7.1.010 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.27

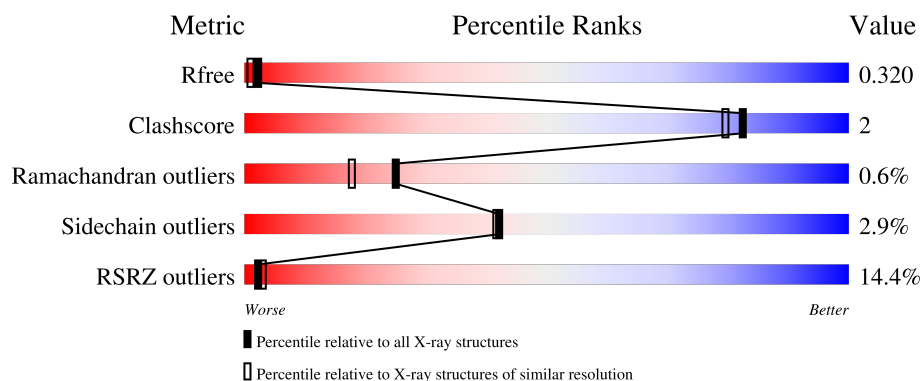
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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	AAA	350	<div> <div>13%</div> <div>83%</div> <div>8%</div> <div>9%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2761 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 Ribonucleoside-diphosphate reductase subunit beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	320	Total	C	N	O	S	0	1	0
			2676	1741	434	494	7			

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	AAA	1	Total	Fe	0	0
			1	1		

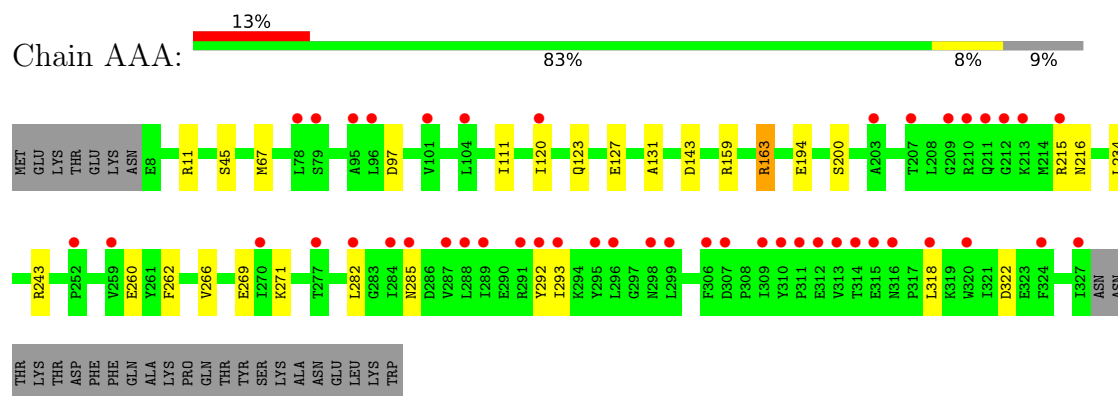
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	AAA	84	Total	O	0	0
			84	84		

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: Ribonucleoside-diphosphate reductase subunit beta



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	69.18Å 69.18Å 177.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.97 – 2.15 48.92 – 2.15	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.97-2.15) 100.0 (48.92-2.15)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.247 , 0.315 0.249 , 0.320	Depositor DCC
R_{free} test set	1203 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	54.3	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	(Not available) , (Not available)	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2761	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.41% 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FE

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	AAA	0.64	0/2742	0.70	0/3717

There are no bond length outliers.

There are no bond angle outliers.

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	AAA	2676	0	2574	13	0
2	AAA	1	0	0	0	0
3	AAA	84	0	0	4	0
All	All	2761	0	2574	13	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 2.

All (13) 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:AAA:163:ARG:NH2	1:AAA:269:GLU:OE1	2.30	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AAA:143:ASP:HB3	3:AAA:574:HOH:O	2.11	0.50
1:AAA:123:GLN:NE2	1:AAA:194:GLU:OE2	2.47	0.47
1:AAA:159:ARG:NH1	3:AAA:506:HOH:O	2.48	0.46
1:AAA:200:SER:CB	1:AAA:269:GLU:OE2	2.64	0.45
1:AAA:97:ASP:OD1	1:AAA:127:GLU:OE1	2.35	0.44
1:AAA:123:GLN:OE1	1:AAA:234:LEU:HD23	2.17	0.44
1:AAA:266:VAL:HG13	1:AAA:293:ILE:HG22	2.02	0.42
1:AAA:292:TYR:CD1	1:AAA:318:LEU:HD12	2.55	0.42
1:AAA:111:ILE:HD12	1:AAA:120:ILE:CD1	2.50	0.41
1:AAA:131:ALA:HB2	3:AAA:526:HOH:O	2.20	0.41
1:AAA:67:MET:SD	1:AAA:216:ASN:HB3	2.60	0.41
1:AAA:271:LYS:C	3:AAA:506:HOH:O	2.59	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	AAA	319/350 (91%)	308 (97%)	9 (3%)	2 (1%)	25	18

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	AAA	260	GLU
1	AAA	282	LEU

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	AAA	281/323 (87%)	273 (97%)	8 (3%)	43	44

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

Mol	Chain	Res	Type
1	AAA	11	ARG
1	AAA	45	SER
1	AAA	163	ARG
1	AAA	215	ARG
1	AAA	243	ARG
1	AAA	262	PHE
1	AAA	285	ASN
1	AAA	322	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

Of 1 ligands modelled in this entry, 1 is 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

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	AAA	320/350 (91%)	0.99	46 (14%) 2 3	36, 64, 96, 122	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	313	VAL	6.2
1	AAA	288	LEU	6.0
1	AAA	292	TYR	5.2
1	AAA	311	PRO	5.0
1	AAA	284	ILE	5.0
1	AAA	310	TYR	4.7
1	AAA	320	TRP	4.6
1	AAA	309	ILE	4.6
1	AAA	287	VAL	4.4
1	AAA	285	ASN	4.3
1	AAA	299	LEU	4.1
1	AAA	318	LEU	4.1
1	AAA	291	ARG	4.0
1	AAA	295	TYR	3.9
1	AAA	282	LEU	3.8
1	AAA	315	GLU	3.8
1	AAA	252	PRO	3.8
1	AAA	78	LEU	3.5
1	AAA	215	ARG	3.4
1	AAA	270	ILE	3.4
1	AAA	307	ASP	3.3
1	AAA	212	GLY	3.2
1	AAA	314	THR	3.2
1	AAA	209	GLY	3.2
1	AAA	259	VAL	3.1
1	AAA	210	ARG	2.9
1	AAA	120	ILE	2.9

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Mol	Chain	Res	Type	RSRZ
1	AAA	327	ILE	2.8
1	AAA	203	ALA	2.8
1	AAA	298	ASN	2.7
1	AAA	289	ILE	2.7
1	AAA	324	PHE	2.7
1	AAA	213	LYS	2.6
1	AAA	316	ASN	2.6
1	AAA	277	THR	2.5
1	AAA	312	GLU	2.5
1	AAA	211	GLN	2.5
1	AAA	296	LEU	2.5
1	AAA	101	VAL	2.4
1	AAA	96	LEU	2.4
1	AAA	207	THR	2.3
1	AAA	293	ILE	2.3
1	AAA	95	ALA	2.2
1	AAA	79	SER	2.1
1	AAA	306	PHE	2.0
1	AAA	104	LEU	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 monosaccharides in this entry.

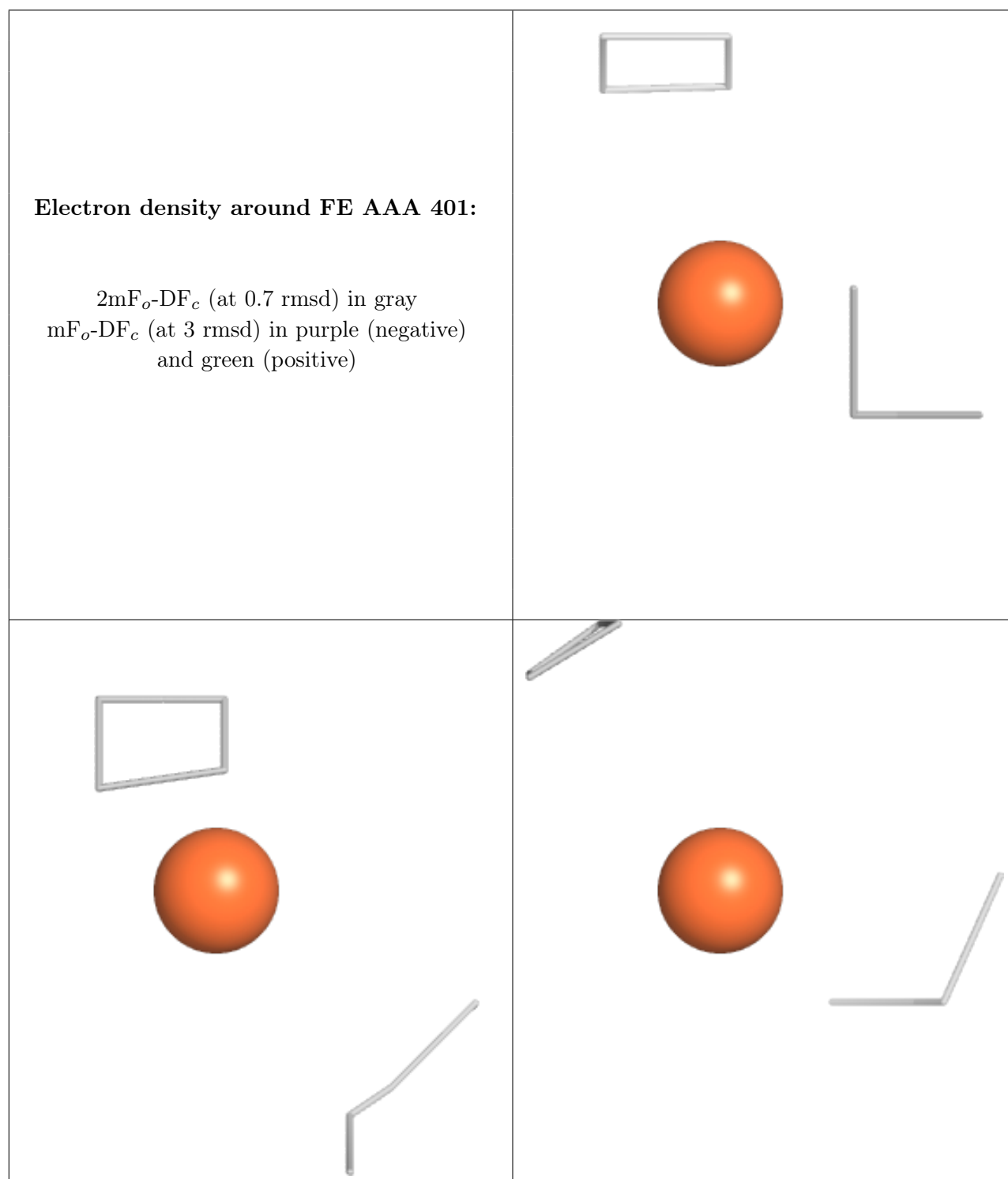
6.4 Ligands [i](#)

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(\AA^2)	Q<0.9
2	FE	AAA	401	1/1	0.95	0.04	80,80,80,80	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.