



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 06:47 am BST

PDB ID : 4NCK
Title : Crystal Structure of Pyrococcus furiosus Rad50 R797G mutation
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Deposited on : 2013-10-24
Resolution : 1.99 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

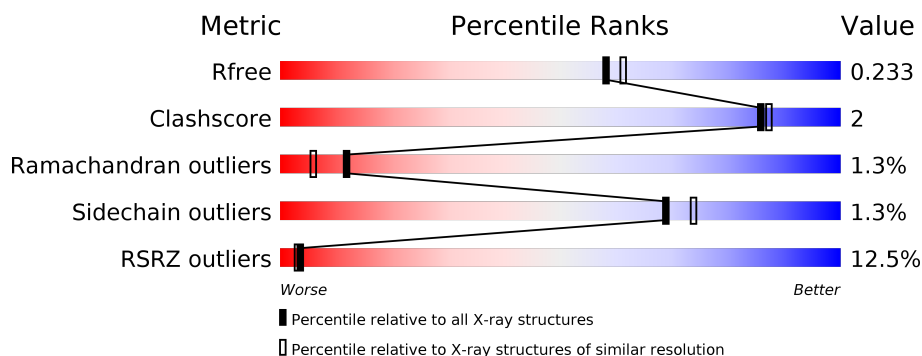
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.99 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 on 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	339	<div> <div>13%</div> <div>84%</div> <div>8%</div> <div>8%</div> </div>
1	B	339	<div> <div>10%</div> <div>86%</div> <div>•</div> <div>9%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	MG	B	903	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 9552 atoms, of which 4562 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 DNA double-strand break repair Rad50 ATPase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	313	Total	C	H	N	O	S	0	5	0
			4625	1512	2284	389	435	5			
1	B	309	Total	C	H	N	O	S	0	2	0
			4596	1491	2278	390	432	5			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	721	GLY	-	LINKER	UNP P58301
A	722	GLY	-	LINKER	UNP P58301
A	723	SER	-	LINKER	UNP P58301
A	724	GLY	-	LINKER	UNP P58301
A	725	GLY	-	LINKER	UNP P58301
A	797	GLY	ARG	ENGINEERED MUTATION	UNP P58301
B	721	GLY	-	LINKER	UNP P58301
B	722	GLY	-	LINKER	UNP P58301
B	723	SER	-	LINKER	UNP P58301
B	724	GLY	-	LINKER	UNP P58301
B	725	GLY	-	LINKER	UNP P58301
B	797	GLY	ARG	ENGINEERED MUTATION	UNP P58301

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Cl	0	0
			1	1		

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		

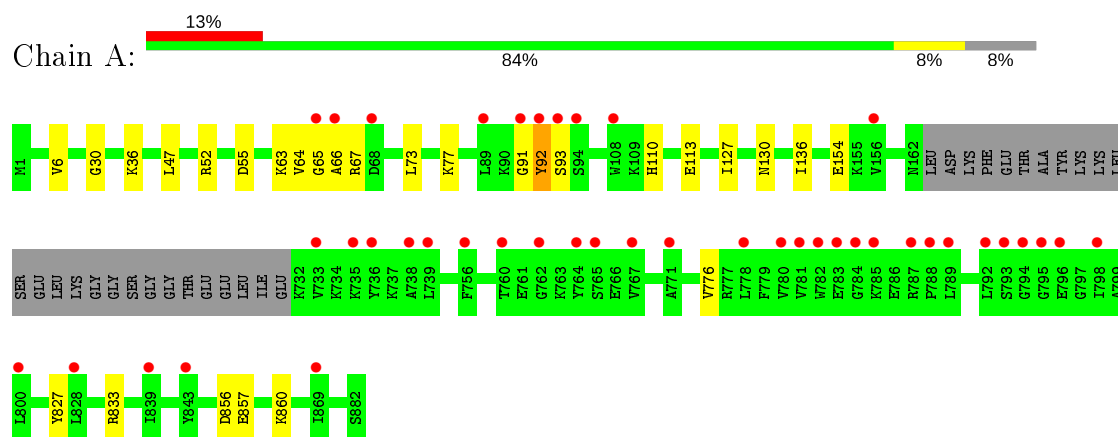
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	200	Total	O	0	0
			200	200		
5	B	124	Total	O	0	0
			124	124		

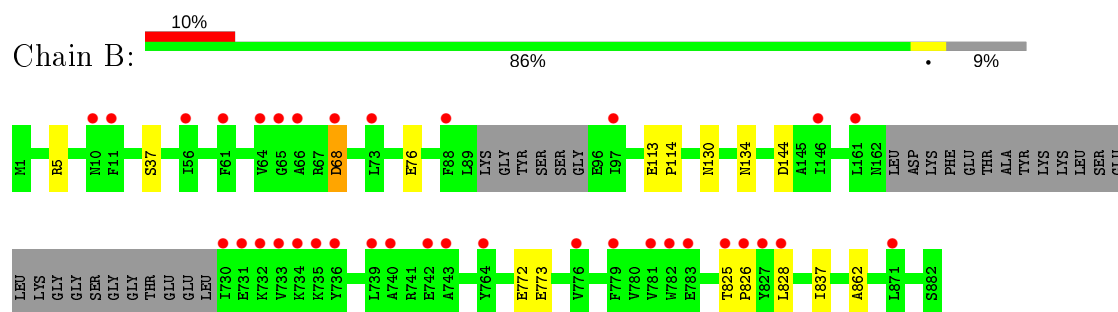
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: DNA double-strand break repair Rad50 ATPase



- Molecule 1: DNA double-strand break repair Rad50 ATPase



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.67Å 68.06Å 202.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.24 – 1.99 48.24 – 1.99	Depositor EDS
% Data completeness (in resolution range)	98.6 (48.24-1.99) 99.3 (48.24-1.99)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.63 (at 2.00Å)	Xtriage
Refinement program	PHENIX dev_1233	Depositor
R, R_{free}	0.181 , 0.230 0.183 , 0.233	Depositor DCC
R_{free} test set	2000 reflections (4.20%)	wwPDB-VP
Wilson B-factor (Å ²)	43.0	Xtriage
Anisotropy	0.281	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 66.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9552	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.83% 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: MG, SO4, CL

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	0.61	0/2390	0.69	0/3233
1	B	0.53	0/2358	0.63	2/3194 (0.1%)
All	All	0.57	0/4748	0.66	2/6427 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	5	ARG	NE-CZ-NH2	-6.79	116.91	120.30
1	B	5	ARG	NE-CZ-NH1	5.08	122.84	120.30

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	2341	2284	2290	15	0
1	B	2318	2278	2268	7	0
2	B	5	0	0	0	0
3	B	1	0	0	0	0
4	B	1	0	0	0	0
5	A	200	0	0	1	2
5	B	124	0	0	3	2

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4990	4562	4558	20	2

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 (20) 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:B:144:ASP:OD2	5:B:1045:HOH:O	2.12	0.67
1:A:6:VAL:HG23	1:A:73[A]:LEU:HD23	1.83	0.58
1:B:76:GLU:OE2	5:B:1032:HOH:O	2.18	0.57
1:A:77:LYS:NZ	5:A:924:HOH:O	2.35	0.53
1:A:6:VAL:HG23	1:A:73[A]:LEU:CD2	2.40	0.50
1:A:30:GLY:O	1:A:36:LYS:HE3	2.12	0.50
1:A:93:SER:OG	1:B:68:ASP:OD1	2.25	0.47
1:B:837:ILE:HG23	1:B:862:ALA:HB2	1.96	0.47
1:A:91:GLY:O	1:A:92:TYR:CB	2.62	0.47
1:A:63:LYS:O	1:A:65:GLY:N	2.48	0.47
1:A:52:ARG:NH2	1:A:55:ASP:OD2	2.49	0.45
1:A:857:GLU:OE1	1:A:860:LYS:NZ	2.31	0.45
1:A:833:ARG:NE	1:A:856:ASP:OD2	2.49	0.45
1:B:113:GLU:OE1	5:B:1030:HOH:O	2.21	0.43
1:A:30:GLY:O	1:A:36:LYS:CE	2.67	0.42
1:A:110:HIS:CD2	1:B:114:PRO:HG2	2.55	0.41
1:B:130:ASN:ND2	1:B:134:ASN:HD22	2.18	0.41
1:A:154:GLU:HA	1:A:776:VAL:HG23	2.01	0.41
1:A:127:ILE:HD11	1:A:136:ILE:CD1	2.50	0.41
1:A:47:LEU:HD21	1:A:73[B]:LEU:HD21	2.03	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:946:HOH:O	5:B:1085:HOH:O[1_655]	2.16	0.04
5:A:930:HOH:O	5:B:1087:HOH:O[1_655]	2.18	0.02

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	314/339 (93%)	303 (96%)	7 (2%)	4 (1%)	12	6
1	B	305/339 (90%)	293 (96%)	8 (3%)	4 (1%)	12	6
All	All	619/678 (91%)	596 (96%)	15 (2%)	8 (1%)	12	6

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	64	VAL
1	A	92	TYR
1	B	773	GLU
1	B	826	PRO
1	A	66	ALA
1	A	67	ARG
1	B	772	GLU
1	B	828	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	A	225/293 (77%)	222 (99%)	3 (1%)	69	74
1	B	229/293 (78%)	226 (99%)	3 (1%)	69	74
All	All	454/586 (78%)	448 (99%)	6 (1%)	69	74

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	113	GLU
1	A	130	ASN
1	A	827	TYR
1	B	37	SER
1	B	68	ASP
1	B	825	THR

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

Mol	Chain	Res	Type
1	A	855	HIS
1	B	130	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	B	901	-	4,4,4	0.13	0	6,6,6	0.10	0

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 [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

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	313/339 (92%)	0.93	43 (13%) 3 2	25, 49, 110, 140	0
1	B	309/339 (91%)	0.79	35 (11%) 5 4	27, 59, 107, 137	0
All	All	622/678 (91%)	0.86	78 (12%) 3 3	25, 55, 109, 140	0

All (78) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	736	TYR	8.1
1	B	66	ALA	7.7
1	B	733	VAL	7.0
1	A	764	TYR	6.9
1	B	65	GLY	6.7
1	A	767	VAL	6.6
1	A	733	VAL	6.5
1	A	92	TYR	6.3
1	A	781	VAL	5.8
1	B	739	LEU	5.8
1	A	782	TRP	5.8
1	B	61	PHE	5.5
1	B	827	TYR	5.5
1	A	792	LEU	5.5
1	B	731	GLU	4.9
1	A	789	LEU	4.9
1	A	787	ARG	4.7
1	A	93	SER	4.7
1	B	730	ILE	4.7
1	B	64	VAL	4.6
1	A	784	GLY	4.3
1	A	843	TYR	4.3
1	A	735	LYS	4.2
1	B	782	TRP	4.2

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Mol	Chain	Res	Type	RSRZ
1	A	94	SER	4.2
1	A	739	LEU	4.1
1	B	88	PHE	4.1
1	A	762	GLY	4.0
1	A	795	GLY	3.9
1	A	771	ALA	3.9
1	B	734	LYS	3.7
1	A	756	PHE	3.7
1	A	765	SER	3.5
1	A	760	THR	3.4
1	B	779	PHE	3.4
1	A	788	PRO	3.4
1	A	66	ALA	3.3
1	B	828	LEU	3.3
1	A	785	LYS	3.3
1	A	738	ALA	3.2
1	A	794	GLY	3.2
1	B	740	ALA	3.2
1	B	825	THR	3.1
1	A	91	GLY	3.1
1	B	732	LYS	3.0
1	B	735	LYS	3.0
1	A	798	ILE	2.9
1	A	800	LEU	2.9
1	A	796	GLU	2.9
1	B	68	ASP	2.9
1	B	826	PRO	2.8
1	B	781	VAL	2.8
1	A	778	LEU	2.8
1	A	68	ASP	2.8
1	B	10	ASN	2.7
1	A	780	VAL	2.7
1	B	743	ALA	2.5
1	B	736	TYR	2.4
1	A	793	SER	2.4
1	A	828	LEU	2.3
1	B	764	TYR	2.3
1	B	783	GLU	2.3
1	B	146	ILE	2.3
1	B	871	LEU	2.2
1	A	156	VAL	2.2
1	A	839	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	869[A]	ILE	2.2
1	A	89	LEU	2.2
1	B	161	LEU	2.2
1	B	776	VAL	2.2
1	B	97	ILE	2.1
1	A	783	GLU	2.1
1	B	73	LEU	2.1
1	B	742	GLU	2.1
1	A	65	GLY	2.1
1	A	108	TRP	2.1
1	B	11	PHE	2.1
1	B	56	ILE	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 carbohydrates 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(Å ²)	Q<0.9
4	MG	B	903	1/1	0.75	0.67	80,80,80,80	0
2	SO4	B	901	5/5	0.81	0.16	118,119,120,121	0
3	CL	B	902	1/1	0.96	0.07	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.