



Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 4Q0R / pdb_00004q0r
Title : The catalytic core of Rad2 (complex I)
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Deposited on : 2014-04-02
Resolution : 2.75 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
Mogul	:	2022.3.0, CSD as543be (2022)
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

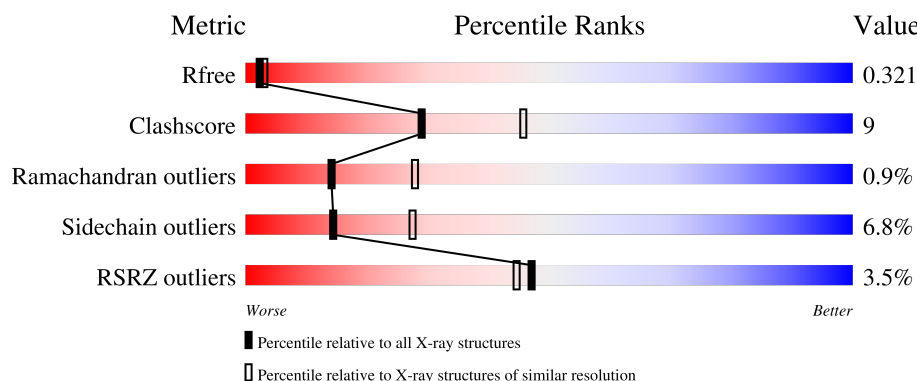
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	367	
1	B	367	
2	D	15	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4556 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 DNA repair protein RAD2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	293	Total	C	N	O	S	Se	0	1	0
			2291	1480	373	426	2	10			
1	B	277	Total	C	N	O	S	Se	0	0	0
			2104	1367	334	391	2	10			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	SER	-	expression tag	UNP P07276
A	1	MSE	-	expression tag	UNP P07276
B	0	SER	-	expression tag	UNP P07276
B	1	MSE	-	expression tag	UNP P07276

- Molecule 2 is a DNA chain called DNA (5'-D(*CP*TP*GP*AP*GP*TP*CP*AP*GP*AP*GP*CP*AP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	4	Total	C	N	O	P	0	0	0
			82	39	15	24	4			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	59	Total	O	0	0
			59	59		
3	B	20	Total	O	0	0
			20	20		



DC	DT	DG	DA	G7	T8	C9	A10	DG	DA	DG	DG	DC	DA	DA	DA
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4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	94.08Å 94.08Å 155.26Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	39.41 – 2.75 39.41 – 2.75	Depositor EDS
% Data completeness (in resolution range)	97.8 (39.41-2.75) 98.8 (39.41-2.75)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.05 (at 2.77Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.245 , 0.310 0.266 , 0.321	Depositor DCC
R_{free} test set	1089 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	58.4	Xtriage
Anisotropy	0.418	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 23.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4556	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.06% 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	0.58	0/2335	0.91	2/3145 (0.1%)
1	B	0.59	1/2140 (0.0%)	0.90	4/2890 (0.1%)
2	D	0.48	0/91	0.51	0/138
All	All	0.58	1/4566 (0.0%)	0.90	6/6173 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	980	ILE	CA-CB	6.39	1.57	1.54

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	11	GLY	N-CA-C	5.75	124.08	112.34
1	B	80	VAL	CA-C-N	5.64	126.89	119.84
1	B	80	VAL	C-N-CA	5.64	126.89	119.84
1	B	873	GLY	CA-C-N	5.33	124.78	119.24
1	B	873	GLY	C-N-CA	5.33	124.78	119.24
1	A	80	VAL	N-CA-C	5.08	113.46	107.77

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	2291	0	2165	47	0
1	B	2104	0	1966	28	0
2	D	82	0	46	2	0
3	A	59	0	0	5	0
3	B	20	0	0	0	0
All	All	4556	0	4177	75	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 9.

All (75) 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:D:8:DT:H2''	2:D:9:DC:H5''	1.64	0.78
1:B:777:GLU:HG3	1:B:781:ARG:HE	1.62	0.64
1:B:884:PHE:O	1:B:886:ASN:N	2.31	0.64
1:A:876:SER:OG	2:D:8:DT:OP1	2.08	0.63
1:A:957:LEU:H	1:A:957:LEU:HD23	1.65	0.62
1:B:3:VAL:HA	1:B:862:GLY:O	2.02	0.60
1:A:963:PHE:HD1	1:A:964:MSE:HE3	1.68	0.57
1:A:26:ARG:NH1	1:A:804:ASN:O	2.31	0.57
1:A:956:ASP:CG	1:A:959:MSE:HE3	2.28	0.57
1:A:963:PHE:CD1	1:A:964:MSE:HE3	2.40	0.56
1:A:5:SER:HB3	1:A:874:PRO:HG3	1.88	0.56
1:B:863:SER:HB2	1:B:865:TYR:H	1.73	0.54
1:B:19:LEU:HD21	1:B:835:VAL:HG13	1.89	0.54
1:B:818:LEU:HD13	1:B:938:TYR:CE2	2.43	0.53
1:A:59:PHE:CZ	1:A:979:LEU:HD21	2.44	0.53
1:A:926:ASP:HB2	3:A:1042:HOH:O	2.08	0.52
1:B:889:ASN:OD1	1:B:889:ASN:N	2.41	0.52
1:A:3:VAL:HA	1:A:862:GLY:O	2.10	0.51
1:A:826:LYS:O	1:A:835:VAL:HG22	2.11	0.51
1:B:965:LYS:HA	1:B:970:TRP:H	1.76	0.51
1:A:871:GLY:O	1:A:917:LYS:HD3	2.11	0.51
1:B:913:ASP:N	1:B:913:ASP:OD1	2.44	0.51
1:A:781:ARG:HG2	1:A:959:MSE:HB3	1.93	0.50
1:A:886:ASN:ND2	1:A:889:ASN:OD1	2.44	0.50
1:A:14:ALA:HB2	1:A:838:TYR:CZ	2.47	0.50
1:A:926:ASP:HB3	1:A:928:ASP:H	1.77	0.49
1:B:980:ILE:O	1:B:982:LEU:N	2.40	0.49
1:A:802:GLN:HG3	3:A:1006:HOH:O	2.12	0.49
1:B:14:ALA:HB1	1:B:836:GLU:HG2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:908:ASN:HB2	1:A:911:GLU:OE2	2.13	0.49
1:A:769:MSE:O	1:A:772:ILE:HG22	2.13	0.49
1:A:956:ASP:OD2	1:A:959:MSE:HE3	2.13	0.49
1:A:856:GLU:HB3	1:A:887:LEU:HG	1.94	0.48
1:B:891:LYS:HG3	1:B:925:LEU:HB3	1.95	0.48
1:A:851:ARG:HA	1:A:854:MSE:HE3	1.96	0.48
1:A:776:GLN:HB3	1:A:786:TYR:CZ	2.49	0.47
1:A:801:LEU:HD22	1:A:809:ILE:HG13	1.96	0.47
1:A:80:VAL:HG13	1:A:84:LYS:HD3	1.98	0.46
1:A:815:ASP:OD2	1:A:864:ASP:HB2	2.16	0.46
1:B:70:ILE:O	1:B:72:PRO:HD3	2.16	0.46
1:A:34:TRP:O	1:A:37:GLN:HG3	2.16	0.46
1:A:853:ASN:ND2	3:A:1026:HOH:O	2.49	0.46
1:B:962:SER:O	1:B:966:THR:HG23	2.16	0.45
1:B:980:ILE:C	1:B:982:LEU:H	2.24	0.45
1:A:3:VAL:HA	1:A:863:SER:HB3	1.97	0.45
1:B:30:ASP:HB2	1:B:794:GLU:OE2	2.16	0.45
1:B:3:VAL:HB	1:B:6:PHE:HB2	1.99	0.45
1:B:6:PHE:HD2	1:B:814:SER:HB2	1.82	0.45
1:A:872:MSE:HE1	1:A:914:LEU:HD11	1.99	0.44
1:B:5:SER:HB3	1:B:874:PRO:HG3	1.99	0.44
1:B:859:GLN:HE21	1:B:934:VAL:HB	1.82	0.44
1:A:947:THR:O	1:A:947:THR:OG1	2.36	0.43
1:A:947:THR:O	1:A:949:PRO:HD3	2.18	0.43
1:A:35:ILE:CD1	1:A:772:ILE:HG13	2.49	0.43
1:A:6:PHE:HD2	1:A:814:SER:HB2	1.84	0.43
1:A:935:TYR:CZ	1:A:939:MSE:HG3	2.54	0.43
1:A:768:THR:C	1:A:770:ASP:H	2.27	0.43
1:A:54:HIS:HD2	1:A:771:MSE:CB	2.32	0.42
1:A:59:PHE:HZ	1:A:979:LEU:HD21	1.82	0.42
1:A:82:VAL:O	1:A:86:GLU:HG3	2.19	0.42
3:A:1031:HOH:O	1:B:18:ARG:HG2	2.19	0.42
1:A:2:GLY:HA3	1:A:814:SER:OG	2.19	0.42
1:B:968:LEU:HD23	1:B:970:TRP:NE1	2.35	0.41
1:A:62:ILE:HD11	1:A:779:LEU:HD23	2.02	0.41
1:A:4:HIS:O	1:A:861:LEU:HB3	2.21	0.41
1:B:3:VAL:HB	1:B:6:PHE:CB	2.50	0.41
1:A:836:GLU:OE2	1:B:832:LYS:HE3	2.20	0.41
1:A:852:LYS:HD3	3:A:1049:HOH:O	2.21	0.41
1:B:894:TYR:HD2	1:B:925:LEU:HD22	1.84	0.41
1:B:968:LEU:HD23	1:B:970:TRP:HE1	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:5:SER:O	1:A:8:ASP:HB2	2.21	0.40
1:B:873:GLY:H	1:B:876:SER:HB2	1.85	0.40
1:B:894:TYR:CD2	1:B:925:LEU:HD22	2.56	0.40
1:A:890:PHE:HE1	1:A:914:LEU:HD21	1.86	0.40
1:A:957:LEU:H	1:A:957:LEU:CD2	2.30	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	288/367 (78%)	269 (93%)	18 (6%)	1 (0%)	36	56
1	B	269/367 (73%)	249 (93%)	16 (6%)	4 (2%)	8	15
All	All	557/734 (76%)	518 (93%)	34 (6%)	5 (1%)	14	28

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	11	GLY
1	B	885	GLY
1	B	926	ASP
1	B	81	PRO
1	B	981	PRO

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	232/313 (74%)	215 (93%)	17 (7%)	13	24
1	B	210/313 (67%)	197 (94%)	13 (6%)	16	31
All	All	442/626 (71%)	412 (93%)	30 (7%)	14	27

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	54	HIS
1	A	56	THR
1	A	82	VAL
1	A	775	VAL
1	A	777	GLU
1	A	788	THR
1	A	792	GLU
1	A	804	ASN
1	A	835	VAL
1	A	857	LEU
1	A	880	VAL
1	A	911	GLU
1	A	943	VAL
1	A	957	LEU
1	A	958	ASP
1	A	964	MSE
1	B	21	SER
1	B	32	SER
1	B	56	THR
1	B	87	THR
1	B	768	THR
1	B	832	LYS
1	B	866	THR
1	B	875	VAL
1	B	889	ASN
1	B	913	ASP
1	B	925	LEU
1	B	933	MSE
1	B	958	ASP

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

Mol	Chain	Res	Type
1	B	859	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 oligosaccharides 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

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	281/367 (76%)	0.44	7 (2%) 58 56	30, 48, 74, 79	2 (0%)
1	B	265/367 (72%)	0.57	12 (4%) 38 37	42, 55, 67, 70	0
2	D	4/15 (26%)	0.59	0 100 100	58, 59, 62, 67	0
All	All	550/749 (73%)	0.50	19 (3%) 47 44	30, 51, 68, 79	2 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	2	GLY	4.6
1	A	3	VAL	4.2
1	B	3	VAL	3.8
1	A	36	TYR	3.7
1	B	4	HIS	3.5
1	B	865	TYR	3.4
1	B	899	PHE	3.1
1	B	768	THR	3.1
1	A	4	HIS	2.9
1	B	0	SER	2.8
1	A	957	LEU	2.7
1	B	87	THR	2.5
1	A	35	ILE	2.5
1	B	963	PHE	2.4
1	B	58	PHE	2.3
1	B	918	LEU	2.3
1	B	928	ASP	2.2
1	A	54	HIS	2.2
1	B	833	ASN	2.1

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.