



# Full wwPDB X-ray Structure Validation Report ⓘ

Jun 22, 2016 – 04:30 AM EDT

PDB ID : 5E37  
Title : Redox protein from Chlamydomonas reinhardtii  
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Deposited on : 2015-10-02  
Resolution : 1.60 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : unknown  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20027790  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20027790

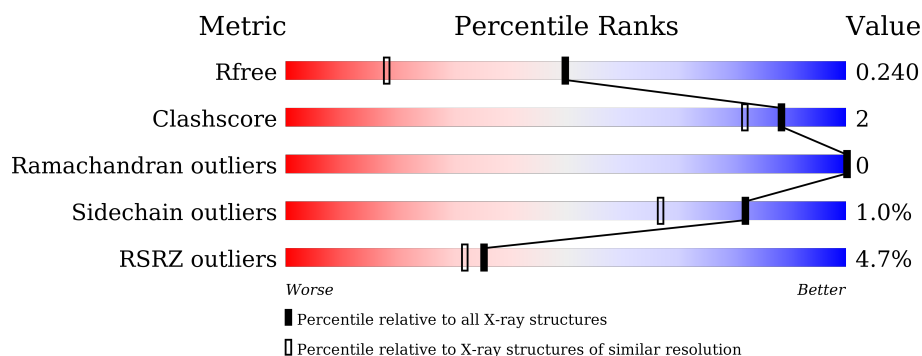
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.60 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))
$R_{free}$	91344	2475 (1.60-1.60)
Clashscore	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)
RSRZ outliers	91569	2479 (1.60-1.60)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	356	<div> <div>4%</div> <div> <div></div> <div>81%</div> <div>6% • 12%</div> </div> </div>
1	C	356	<div> <div>4%</div> <div> <div></div> <div>77%</div> <div>10% • 13%</div> </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
2	CA	C	404	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5719 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 EF-Hand domain-containing thioredoxin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	0	0
			2432	1540	397	476	19			
1	C	310	Total	C	N	O	S	0	1	0
			2430	1540	400	471	19			

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	Ca	0	0
			4	4		
2	C	4	Total	Ca	0	0
			4	4		

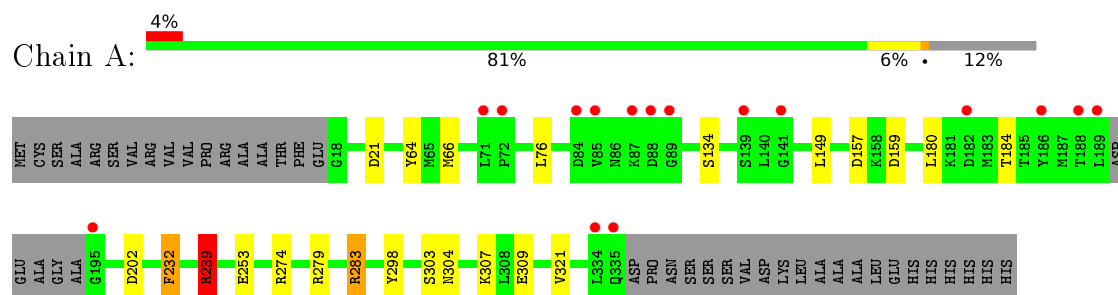
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	410	Total	O	0	0
			410	410		
3	C	439	Total	O	0	0
			439	439		

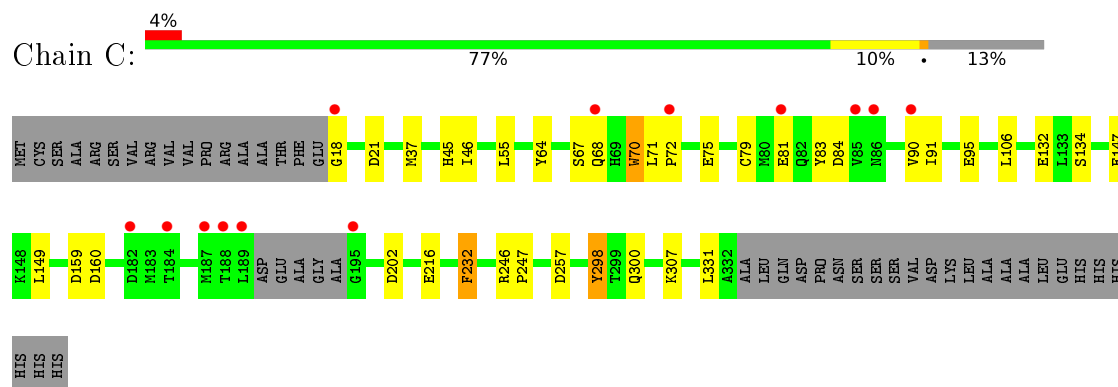
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: EF-Hand domain-containing thioredoxin



- Molecule 1: EF-Hand domain-containing thioredoxin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.96 Å   55.19 Å   89.58 Å 90.00°   101.00°   90.00°	Depositor
Resolution (Å)	50.00 – 1.60 43.97 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.00-1.60) 99.7 (43.97-1.60)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.68 (at 1.60 Å)	Xtriage
Refinement program	REFMAC 5.8.0103	Depositor
R, $R_{free}$	0.199 , 0.233 0.207 , 0.240	Depositor DCC
$R_{free}$ test set	5367 reflections (5.20%)	DCC
Wilson B-factor (Å <sup>2</sup> )	17.4	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 50.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5719	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 40.58 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.6775e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup> Intensities estimated from amplitudes.

<sup>2</sup> Theoretical 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: CA

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	1.16	3/2473 (0.1%)	1.19	18/3332 (0.5%)
1	C	1.22	6/2471 (0.2%)	1.20	13/3323 (0.4%)
All	All	1.19	9/4944 (0.2%)	1.19	31/6655 (0.5%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	70	TRP	CB-CG	-9.65	1.32	1.50
1	C	147	GLU	CD-OE1	8.17	1.34	1.25
1	C	64	TYR	CE1-CZ	-7.22	1.29	1.38
1	C	18	GLY	N-CA	5.63	1.54	1.46
1	C	132	GLU	CD-OE1	-5.59	1.19	1.25
1	A	64	TYR	CG-CD2	-5.57	1.31	1.39
1	A	253	GLU	CD-OE2	5.54	1.31	1.25
1	A	64	TYR	CB-CG	5.33	1.59	1.51
1	C	216	GLU	CD-OE1	-5.15	1.20	1.25

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	21	ASP	CB-CG-OD2	-9.26	109.97	118.30
1	A	298	TYR	CB-CG-CD2	-8.80	115.72	121.00
1	C	84	ASP	CB-CG-OD2	-8.64	110.52	118.30
1	A	21	ASP	CB-CG-OD2	-8.16	110.95	118.30
1	C	202	ASP	CB-CG-OD2	-8.07	111.03	118.30
1	A	159	ASP	CB-CG-OD1	7.58	125.12	118.30
1	C	21	ASP	CB-CG-OD1	7.24	124.82	118.30
1	A	232	PHE	CB-CG-CD1	7.02	125.71	120.80
1	A	274	ARG	NE-CZ-NH1	6.95	123.78	120.30
1	A	202	ASP	CB-CG-OD2	-6.93	112.07	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	21	ASP	CB-CG-OD1	6.90	124.51	118.30
1	C	232	PHE	CB-CG-CD1	6.88	125.62	120.80
1	A	279	ARG	NE-CZ-NH2	-6.52	117.04	120.30
1	C	257	ASP	CB-CG-OD1	6.52	124.17	118.30
1	C	84	ASP	CB-CG-OD1	6.40	124.06	118.30
1	C	246	ARG	NE-CZ-NH1	6.35	123.47	120.30
1	C	159	ASP	CB-CG-OD1	6.24	123.91	118.30
1	A	298	TYR	CB-CG-CD1	6.18	124.71	121.00
1	C	160	ASP	CB-CG-OD1	5.86	123.57	118.30
1	C	298	TYR	CB-CG-CD2	-5.76	117.55	121.00
1	A	298	TYR	OH-CZ-CE2	-5.69	104.75	120.10
1	A	232	PHE	CB-CG-CD2	-5.60	116.88	120.80
1	A	157	ASP	CB-CG-OD1	5.54	123.29	118.30
1	A	239	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	A	321	VAL	CA-C-N	-5.42	105.37	116.20
1	C	64	TYR	CB-CG-CD2	-5.23	117.86	121.00
1	A	309	GLU	OE1-CD-OE2	5.22	129.56	123.30
1	C	232	PHE	CB-CG-CD2	-5.20	117.16	120.80
1	A	303	SER	CA-CB-OG	-5.12	97.38	111.20
1	A	283	ARG	NE-CZ-NH1	5.09	122.85	120.30
1	A	76	LEU	CB-CG-CD2	-5.01	102.48	111.00

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	2432	0	2362	5	1
1	C	2430	0	2385	17	1
2	A	4	0	0	0	0
2	C	4	0	0	0	0
3	A	410	0	0	2	0
3	C	439	0	0	4	0
All	All	5719	0	4747	22	1



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 (22) 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:C:45:HIS:CD2	1:C:90:VAL:HG11	2.21	0.75
1:C:81:GLU:CB	3:C:865:HOH:O	2.39	0.70
1:C:45:HIS:CG	1:C:90:VAL:HG11	2.32	0.65
1:C:300[B]:GLN:NE2	3:C:505:HOH:O	2.40	0.55
1:C:90:VAL:HG12	1:C:91:ILE:N	2.24	0.53
1:C:70:TRP:CZ2	1:C:106:LEU:HD11	2.48	0.48
1:C:71:LEU:HG	1:C:75:GLU:HB3	1.95	0.48
1:C:45:HIS:CD2	1:C:90:VAL:CG1	2.95	0.47
1:C:72:PRO:O	1:C:75:GLU:N	2.48	0.47
1:C:331:LEU:CG	3:C:501:HOH:O	2.62	0.46
1:A:239:ARG:NH1	3:A:508:HOH:O	2.49	0.46
1:C:68:GLN:HG2	3:C:624:HOH:O	2.16	0.45
1:A:283:ARG:NH1	3:A:509:HOH:O	2.50	0.44
1:C:300[B]:GLN:OE1	1:C:307:LYS:HE2	2.18	0.44
1:A:180:LEU:O	1:A:184:THR:HG23	2.19	0.43
1:C:83:TYR:O	1:C:95:GLU:HB3	2.19	0.42
1:C:298:TYR:OH	1:C:300[A]:GLN:OE1	2.32	0.42
1:A:134:SER:HB2	1:A:149:LEU:HD11	2.00	0.42
1:A:304:ASN:ND2	1:A:307:LYS:HE3	2.35	0.42
1:C:134:SER:HB2	1:C:149:LEU:HD11	2.01	0.41
1:C:46:ILE:HB	1:C:91:ILE:HB	2.01	0.41
1:C:55:LEU:HD11	1:C:79:CYS:SG	2.61	0.41

All (1) 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)
1:A:66:MET:O	1:C:37:MET:SD[1_656]	1.90	0.30

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	309/356 (87%)	306 (99%)	3 (1%)	0	100	100
1	C	307/356 (86%)	305 (99%)	2 (1%)	0	100	100
All	All	616/712 (86%)	611 (99%)	5 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

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	259/305 (85%)	257 (99%)	2 (1%)	86	75
1	C	261/305 (86%)	258 (99%)	3 (1%)	80	63
All	All	520/610 (85%)	515 (99%)	5 (1%)	82	67

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	232	PHE
1	A	239	ARG
1	C	67	SER
1	C	232	PHE
1	C	247	PRO

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

Mol	Chain	Res	Type
1	A	335	GLN

### 5.3.3 RNA ⓘ

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 8 ligands modelled in this entry, 8 are 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 [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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	313/356 (87%)	0.09	16 (5%) 32 29	11, 22, 48, 78	0
1	C	310/356 (87%)	0.06	13 (4%) 40 36	10, 21, 49, 70	0
All	All	623/712 (87%)	0.08	29 (4%) 35 32	10, 22, 49, 78	0

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	85	VAL	8.5
1	C	85	VAL	5.7
1	A	188	THR	4.4
1	A	84	ASP	4.1
1	C	189	LEU	4.0
1	C	188	THR	3.8
1	C	187	MET	3.6
1	A	141	GLY	3.6
1	C	90	VAL	3.3
1	C	18	GLY	3.3
1	A	195	GLY	3.2
1	A	71	LEU	3.1
1	C	195	GLY	3.1
1	A	87	LYS	3.1
1	A	335	GLN	3.1
1	A	189	LEU	3.0
1	C	81	GLU	2.8
1	C	182	ASP	2.7
1	A	334	LEU	2.6
1	C	184	THR	2.5
1	C	72	PRO	2.5
1	A	139	SER	2.4
1	C	68	GLN	2.4
1	A	89	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	86	ASN	2.3
1	A	72	PRO	2.2
1	A	88	ASP	2.1
1	A	182	ASP	2.1
1	A	186	TYR	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	CA	C	404	1/1	1.00	0.09	2.33	13,13,13,13	0
2	CA	A	404	1/1	0.99	0.08	1.38	15,15,15,15	0
2	CA	A	403	1/1	0.99	0.07	-0.13	15,15,15,15	0
2	CA	C	403	1/1	1.00	0.06	-0.80	13,13,13,13	0
2	CA	A	402	1/1	0.93	0.09	-1.32	36,36,36,36	0
2	CA	A	401	1/1	0.93	0.06	-1.97	33,33,33,33	0
2	CA	C	401	1/1	0.95	0.05	-2.23	28,28,28,28	0
2	CA	C	402	1/1	0.96	0.07	-2.60	41,41,41,41	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.