



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 20, 2026 – 03:50 PM UTC

PDB ID : 6IUS / pdb\_00006ius  
Title : A higher kcat Rubisco  
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Deposited on : 2018-11-30  
Resolution : 2.12 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

<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>.

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

MolProbity : 4-5-2 with Phenix2.0  
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

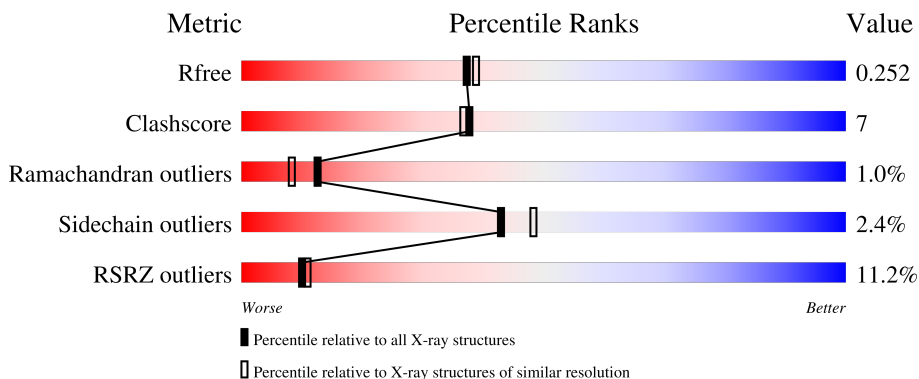
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.12 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	8290 (2.14-2.10)
Clashscore	190562	8817 (2.14-2.10)
Ramachandran outliers	187476	8738 (2.14-2.10)
Sidechain outliers	187428	8739 (2.14-2.10)
RSRZ outliers	180081	8294 (2.14-2.10)

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  $\geq 3$ , 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	461	
1	B	461	
1	C	461	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10515 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 Ribulose-1,5-bisphosphate carboxylase/oxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	426	Total 3294	C 2103	N 553	O 620	S 18	0	0	0
1	B	419	Total 3244	C 2072	N 545	O 610	S 17	0	0	0
1	C	422	Total 3271	C 2086	N 553	O 615	S 17	0	1	0

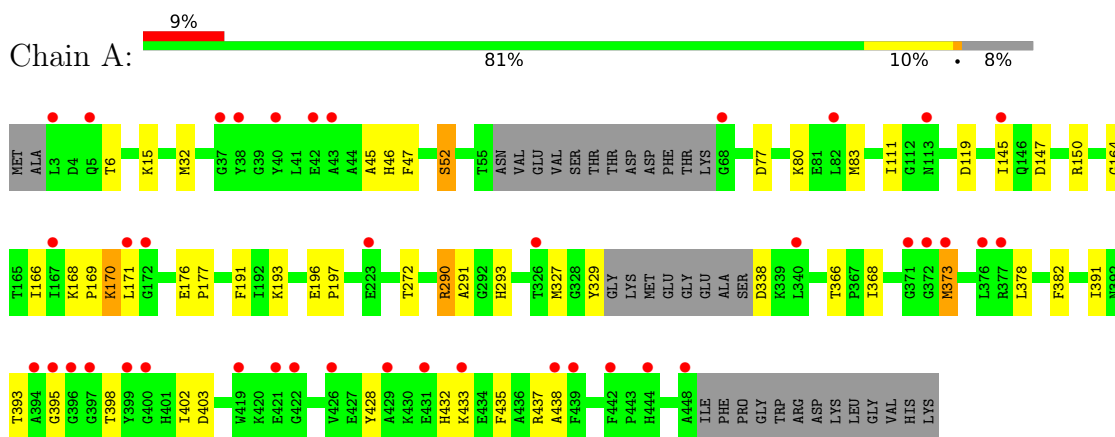
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	237	Total 237	O 237	0	0
2	B	225	Total 225	O 225	0	0
2	C	244	Total 244	O 244	0	0

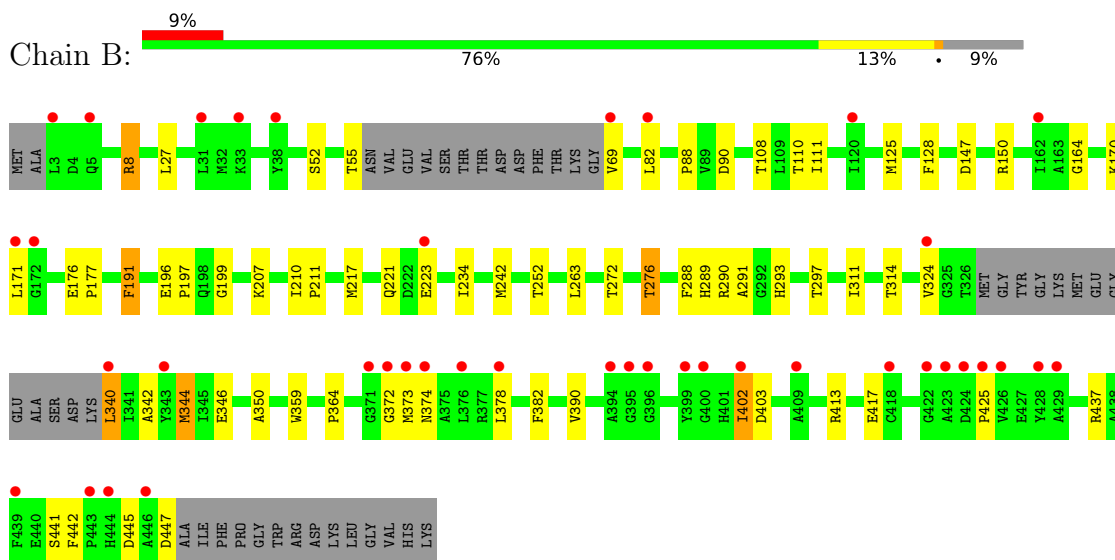
### 3 Residue-property plots [i](#)

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.

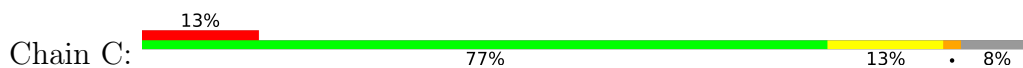
- Molecule 1: Ribulose-1,5-bisphosphate carboxylase/oxygenase

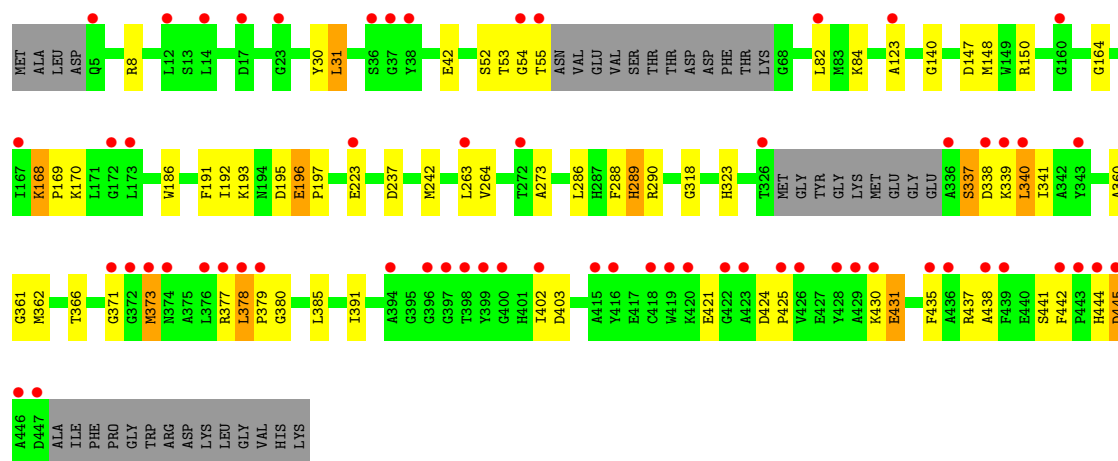


- Molecule 1: Ribulose-1,5-bisphosphate carboxylase/oxygenase



- Molecule 1: Ribulose-1,5-bisphosphate carboxylase/oxygenase





## 4 Data and refinement statistics i

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	166.50Å 107.84Å 112.31Å 90.00° 130.04° 90.00°	Depositor
Resolution (Å)	46.99 – 2.12 46.99 – 2.12	Depositor EDS
% Data completeness (in resolution range)	96.1 (46.99-2.12) 96.2 (46.99-2.12)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.48 (at 2.12Å)	Xtrriage
Refinement program	REFMAC 5.8.0238	Depositor
R, $R_{free}$	0.192 , 0.231 (Not available) , 0.252	Depositor DCC
$R_{free}$ test set	1926 reflections (2.23%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	33.4	Xtrriage
Anisotropy	0.069	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 41.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	0.000 for -h+k-l,-l,-k 0.000 for -h-k-l,l,k 0.003 for -h-2*l,-k,l	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	10515	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.49% of the height of the origin peak. No significant pseudotranslation is detected.*

<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](#)

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	1.09	0/3378	1.38	3/4573 (0.1%)
1	B	1.09	1/3327 (0.0%)	1.40	2/4506 (0.0%)
1	C	1.09	0/3354	1.44	7/4540 (0.2%)
All	All	1.09	1/10059 (0.0%)	1.41	12/13619 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	390	VAL	C-O	5.75	1.29	1.23

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	191	PHE	CA-CB-CG	6.64	120.44	113.80
1	A	47	PHE	CA-CB-CG	5.90	119.70	113.80
1	C	30	TYR	CB-CA-C	-5.71	97.44	109.38
1	C	169	PRO	CA-C-N	5.43	128.34	120.79
1	C	169	PRO	C-N-CA	5.43	128.34	120.79

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	3294	0	3176	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3244	0	3130	52	0
1	C	3271	0	3157	56	0
2	A	237	0	0	2	0
2	B	225	0	0	4	0
2	C	244	0	0	2	0
All	All	10515	0	9463	135	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 7.

The worst 5 of 135 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:344:MET:CE	1:B:359:TRP:HE1	1.53	1.20
1:B:344:MET:HE2	1:B:359:TRP:HE1	1.07	1.15
1:B:344:MET:HE2	1:B:359:TRP:NE1	1.71	1.05
1:B:342:ALA:O	1:B:346:GLU:HG2	1.73	0.89
1:C:338:ASP:O	1:C:341:ILE:HB	1.77	0.85

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	420/461 (91%)	403 (96%)	14 (3%)	3 (1%)	18 15
1	B	413/461 (90%)	388 (94%)	20 (5%)	5 (1%)	10 6
1	C	417/461 (90%)	400 (96%)	12 (3%)	5 (1%)	10 6
All	All	1250/1383 (90%)	1191 (95%)	46 (4%)	13 (1%)	12 8

5 of 13 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	54	GLY
1	C	373	MET
1	C	445	ASP
1	A	52	SER
1	A	373	MET

### 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	332/361 (92%)	328 (99%)	4 (1%)	63	71
1	B	328/361 (91%)	319 (97%)	9 (3%)	39	44
1	C	330/361 (91%)	319 (97%)	11 (3%)	33	36
All	All	990/1083 (91%)	966 (98%)	24 (2%)	43	48

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	168	LYS
1	C	286	LEU
1	C	223	GLU
1	C	289	HIS
1	B	223	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	384	ASN
1	C	289	HIS
1	B	392	ASN
1	B	259	GLN
1	C	146	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 [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, 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	426/461 (92%)	0.67	40 (9%) 14 15	22, 35, 68, 100	0
1	B	419/461 (90%)	0.78	40 (9%) 14 15	24, 37, 70, 94	0
1	C	422/461 (91%)	0.93	62 (14%) 6 6	16, 40, 69, 96	1 (0%)
All	All	1267/1383 (91%)	0.80	142 (11%) 10 11	16, 38, 69, 100	1 (0%)

The worst 5 of 142 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	446	ALA	6.4
1	C	376	LEU	6.2
1	C	443	PRO	6.0
1	B	446	ALA	5.4
1	C	442	PHE	5.1

### 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.