



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 12, 2021 – 01:48 PM EDT

PDB ID : 1YKW
Title : Crystal Structure of a Novel RuBisCO-Like Protein from the Green Sulfur Bacterium *Chlorobium tepidum*
Authors : Li, H.; Sawaya, M.R.; Tabita, F.R.; Eisenberg, D.
Deposited on : 2005-01-18
Resolution : 2.00 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	:	2.23.2
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)	:	2.23.2

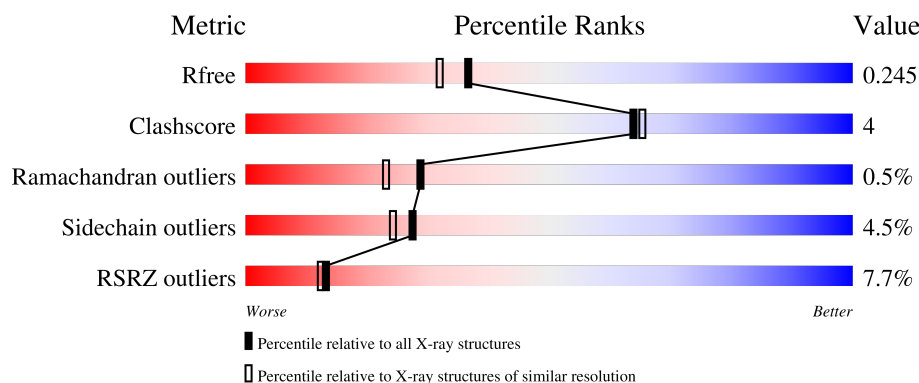
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


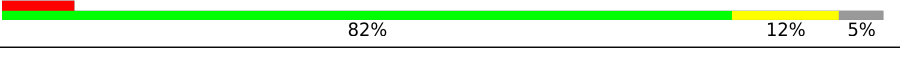
The reported resolution of this entry is 2.00 Å.

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 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	435	
1	B	435	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6736 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 RuBisCO-like protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	414	Total	C	N	O	S	0	0	0
			3205	2048	537	602	18			
1	B	412	Total	C	N	O	S	0	2	0
			3214	2052	542	602	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	328	VAL	MET	engineered mutation	UNP Q8KBL4
B	328	VAL	MET	engineered mutation	UNP Q8KBL4


- Molecule 2 is water.

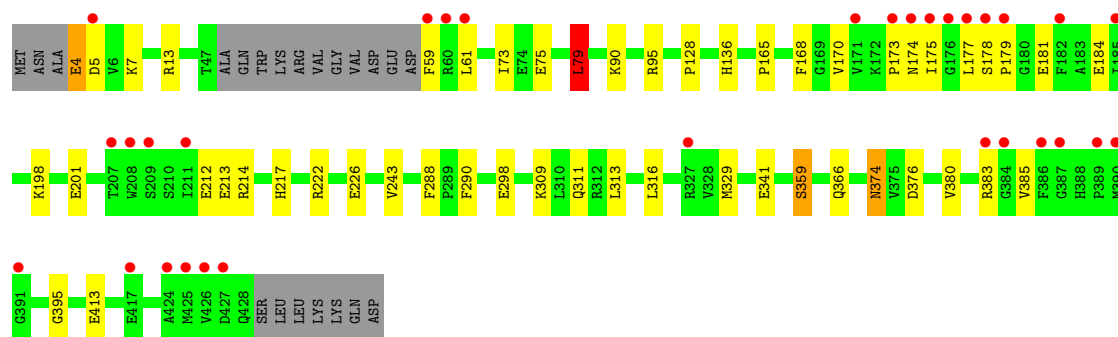
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	156	Total	O	0	0
			156	156		
2	B	161	Total	O	0	0
			161	161		

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.

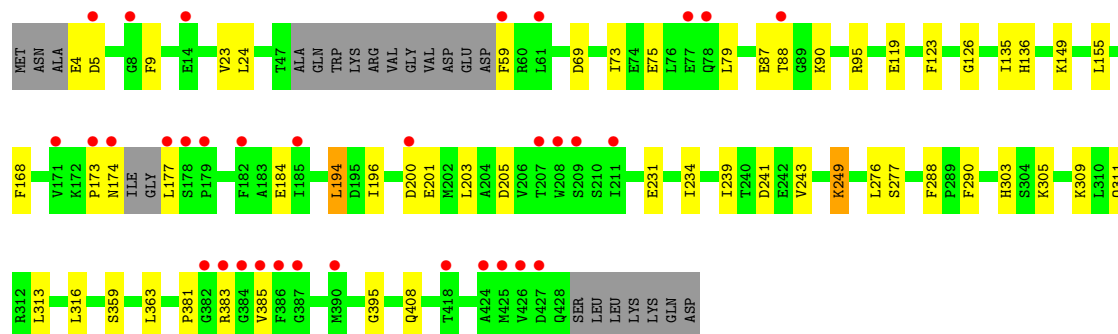
- Molecule 1: RuBisCO-like protein

Chain A: 



- Molecule 1: RuBisCO-like protein

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	67.35Å 78.45Å 90.37Å 90.00° 99.95° 90.00°	Depositor
Resolution (Å)	87.71 – 2.00 50.65 – 2.01	Depositor EDS
% Data completeness (in resolution range)	97.9 (87.71-2.00) 97.4 (50.65-2.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.23 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.201 , 0.245 0.201 , 0.245	Depositor DCC
R_{free} test set	2956 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	22.6	Xtriage
Anisotropy	0.911	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 54.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6736	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.27% 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	0.60	0/3281	0.78	1/4450 (0.0%)
1	B	0.59	0/3290	0.77	1/4460 (0.0%)
All	All	0.60	0/6571	0.77	2/8910 (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	194	LEU	CB-CG-CD1	-5.92	100.94	111.00
1	A	79	LEU	CA-CB-CG	5.16	127.16	115.30

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	3205	0	3163	27	0
1	B	3214	0	3166	29	0
2	A	156	0	0	5	0
2	B	161	0	0	4	0
All	All	6736	0	6329	53	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 4.

All (53) 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:136[A]:HIS:ND1	2:B:480:HOH:O	2.03	0.91
1:B:205:ASP:HB2	1:B:249:LYS:HE3	1.62	0.81
1:B:303:HIS:HD2	1:B:305:LYS:H	1.29	0.80
1:B:231:GLU:OE1	2:B:548:HOH:O	1.99	0.80
1:B:311:GLN:HB3	1:B:316:LEU:HD12	1.66	0.78
1:A:73:ILE:HD11	1:A:95:ARG:HG3	1.73	0.71
1:A:243:VAL:HG11	1:B:243:VAL:HG11	1.73	0.70
1:A:136:HIS:ND1	2:A:490:HOH:O	2.24	0.70
1:B:136[A]:HIS:CE1	2:B:480:HOH:O	2.43	0.69
1:B:75:GLU:OE2	1:B:90:LYS:HD2	1.95	0.67
1:B:203:LEU:O	1:B:203:LEU:HG	1.97	0.65
1:B:73:ILE:HD11	1:B:95:ARG:HG3	1.80	0.64
1:A:309:LYS:HE2	1:A:341:GLU:OE1	1.99	0.63
1:B:173:PRO:HD2	1:B:177:LEU:HD11	1.82	0.60
1:A:217:HIS:ND1	2:A:533:HOH:O	2.31	0.59
1:A:359:SER:HA	2:A:517:HOH:O	2.01	0.59
1:A:79:LEU:HD21	1:A:128:PRO:HG2	1.85	0.59
1:B:174:ASN:ND2	1:B:200:ASP:OD1	2.36	0.57
1:A:222:ARG:O	1:A:226:GLU:HG3	2.04	0.57
1:B:149:LYS:HD2	1:B:277:SER:O	2.05	0.56
1:A:170:VAL:HG22	1:A:198:LYS:HD3	1.88	0.56
1:A:173:PRO:HG2	1:A:177:LEU:HD11	1.87	0.55
1:A:385:VAL:HA	1:A:395:GLY:HA3	1.88	0.54
1:A:311:GLN:HB3	1:A:316:LEU:HD12	1.89	0.54
1:A:179:PRO:HG2	1:A:213:GLU:HG3	1.89	0.53
1:A:5:ASP:HB3	1:A:59:PHE:CE2	2.43	0.53
1:B:205:ASP:HB2	1:B:249:LYS:CE	2.34	0.53
1:A:329:MET:HE3	1:B:123:PHE:HA	1.90	0.53
1:B:87:GLU:HG3	1:B:126:GLY:HA2	1.90	0.52
1:B:196:ILE:HG12	1:B:234:ILE:HB	1.91	0.52
1:B:9:PHE:CZ	1:B:59:PHE:HB3	2.46	0.51
1:A:374:ASN:HD22	1:A:376:ASP:H	1.60	0.49
1:A:5:ASP:HB3	1:A:59:PHE:HE2	1.79	0.47
1:A:4:GLU:HA	1:A:4:GLU:OE1	2.14	0.47
1:A:168:PHE:HB2	1:A:380:VAL:HG22	1.96	0.47
1:B:385:VAL:HA	1:B:395:GLY:HA3	1.96	0.47
1:B:309:LYS:HG3	2:B:484:HOH:O	2.15	0.47
1:A:329:MET:CE	1:B:123:PHE:HA	2.45	0.46
1:A:174:ASN:HD22	1:A:175:ILE:HG12	1.81	0.46
1:B:241:ASP:OD2	1:B:249:LYS:HD2	2.16	0.46
1:A:174:ASN:ND2	1:A:175:ILE:HG12	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:366:GLN:NE2	2:A:521:HOH:O	2.49	0.44
1:B:168:PHE:O	1:B:381:PRO:HD2	2.17	0.44
1:A:136:HIS:CE1	2:A:490:HOH:O	2.66	0.43
1:B:205:ASP:CB	1:B:249:LYS:HE3	2.39	0.43
1:A:179:PRO:HB3	1:A:214:ARG:HA	2.02	0.42
1:A:75:GLU:OE2	1:A:90:LYS:HD3	2.20	0.42
1:A:178:SER:HB2	1:A:181:GLU:CD	2.40	0.41
1:B:24:LEU:HD23	1:B:135:ILE:HG23	2.02	0.41
1:B:309:LYS:HB2	1:B:309:LYS:HE2	1.93	0.41
1:B:381:PRO:HG2	1:B:385:VAL:HG21	2.01	0.41
1:B:23:VAL:HB	1:B:136[B]:HIS:HB3	2.02	0.41
1:B:239:ILE:HD11	1:B:276:LEU:HD21	2.03	0.41

There are no symmetry-related clashes.

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	410/435 (94%)	396 (97%)	12 (3%)	2 (0%)	29	23
1	B	408/435 (94%)	393 (96%)	13 (3%)	2 (0%)	29	23
All	All	818/870 (94%)	789 (96%)	25 (3%)	4 (0%)	29	23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	290	PHE
1	B	290	PHE
1	A	288	PHE
1	B	288	PHE

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	343/361 (95%)	328 (96%)	15 (4%)	28	25
1	B	344/361 (95%)	328 (95%)	16 (5%)	26	22
All	All	687/722 (95%)	656 (96%)	31 (4%)	27	24

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	7	LYS
1	A	13	ARG
1	A	61	LEU
1	A	79	LEU
1	A	165	PRO
1	A	184	GLU
1	A	201	GLU
1	A	212	GLU
1	A	298	GLU
1	A	313	LEU
1	A	359	SER
1	A	374	ASN
1	A	383	ARG
1	A	413	GLU
1	B	4	GLU
1	B	5	ASP
1	B	69	ASP
1	B	79	LEU
1	B	88	THR
1	B	119	GLU
1	B	155	LEU
1	B	184	GLU
1	B	194	LEU
1	B	201	GLU
1	B	249	LYS
1	B	313	LEU

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Mol	Chain	Res	Type
1	B	359	SER
1	B	363	LEU
1	B	383	ARG
1	B	408	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	GLN
1	A	92	HIS
1	A	104	ASN
1	A	366	GLN
1	A	374	ASN
1	B	63	HIS
1	B	92	HIS
1	B	160	ASN
1	B	303	HIS
1	B	408	GLN
1	B	428	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	A	414/435 (95%)	0.31	31 (7%) 14 13	16, 25, 50, 62	0
1	B	412/435 (94%)	0.33	33 (8%) 12 11	16, 25, 50, 58	1 (0%)
All	All	826/870 (94%)	0.32	64 (7%) 13 12	16, 25, 50, 62	1 (0%)

All (64) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	387	GLY	7.5
1	A	208	TRP	5.8
1	A	173	PRO	5.7
1	A	175	ILE	5.6
1	B	209	SER	5.5
1	B	207	THR	5.5
1	A	59	PHE	5.5
1	B	179	PRO	5.2
1	A	61	LEU	5.1
1	A	387	GLY	4.9
1	B	59	PHE	4.8
1	B	61	LEU	4.8
1	A	174	ASN	4.6
1	B	208	TRP	4.6
1	A	177	LEU	4.4
1	B	174	ASN	4.3
1	B	171	VAL	4.2
1	B	386	PHE	4.1
1	B	384	GLY	4.1
1	A	179	PRO	4.0
1	A	182	PHE	3.9
1	B	173	PRO	3.8
1	B	427	ASP	3.7
1	A	384	GLY	3.7

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Mol	Chain	Res	Type	RSRZ
1	B	383	ARG	3.7
1	A	424	ALA	3.6
1	B	5	ASP	3.6
1	A	383	ARG	3.5
1	B	88	THR	3.4
1	A	209	SER	3.4
1	A	426	VAL	3.2
1	B	200	ASP	3.1
1	B	182	PHE	3.1
1	B	8	GLY	3.1
1	A	176	GLY	2.9
1	A	185	ILE	2.9
1	A	386	PHE	2.9
1	A	5	ASP	2.8
1	B	177	LEU	2.7
1	A	171	VAL	2.7
1	A	207	THR	2.6
1	A	211	ILE	2.5
1	B	425	MET	2.4
1	A	427	ASP	2.4
1	A	327	ARG	2.3
1	A	425	MET	2.2
1	B	385	VAL	2.2
1	B	382	GLY	2.2
1	B	77	GLU	2.2
1	B	211	ILE	2.2
1	A	391	GLY	2.2
1	B	424	ALA	2.2
1	B	78	GLN	2.1
1	B	418	THR	2.1
1	A	178	SER	2.1
1	A	417	GLU	2.1
1	B	426	VAL	2.1
1	A	389	PRO	2.1
1	A	60	ARG	2.1
1	B	178	SER	2.1
1	B	185	ILE	2.1
1	A	390	MET	2.1
1	B	390	MET	2.0
1	B	14	GLU	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.