



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 12:15 PM UTC

PDB ID : 4IJG / pdb_00004ijg
Title : Crystal structure of monomeric bacteriophytochrome
Authors : Auldridge, M.E.
Deposited on : 2012-12-21
Resolution : 1.70 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

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A user guide is available at

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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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

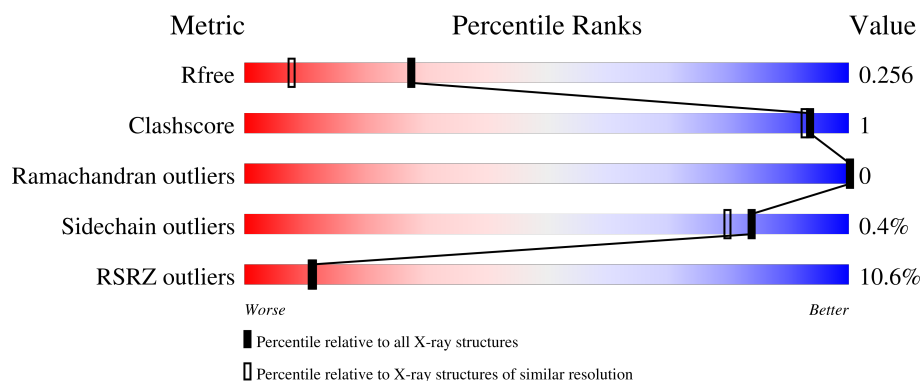
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	5551 (1.70-1.70)
Clashscore	190562	5924 (1.70-1.70)
Ramachandran outliers	187476	5846 (1.70-1.70)
Sidechain outliers	187428	5846 (1.70-1.70)
RSRZ outliers	180081	5554 (1.70-1.70)

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	341	10% 88% 9%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2689 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 Bacteriophytochrome.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	311	2411	1533	425	443	10	0	7	0

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	expression tag	UNP Q9RZA4
A	-10	ALA	-	expression tag	UNP Q9RZA4
A	-9	SER	-	expression tag	UNP Q9RZA4
A	-8	MET	-	expression tag	UNP Q9RZA4
A	-7	THR	-	expression tag	UNP Q9RZA4
A	-6	GLY	-	expression tag	UNP Q9RZA4
A	-5	GLY	-	expression tag	UNP Q9RZA4
A	-4	GLN	-	expression tag	UNP Q9RZA4
A	-3	GLN	-	expression tag	UNP Q9RZA4
A	-2	MET	-	expression tag	UNP Q9RZA4
A	-1	GLY	-	expression tag	UNP Q9RZA4
A	0	ARG	-	expression tag	UNP Q9RZA4
A	1	GLY	-	expression tag	UNP Q9RZA4
A	2	SER	-	expression tag	UNP Q9RZA4
A	?	-	PRO	deletion	UNP Q9RZA4
A	?	-	LEU	deletion	UNP Q9RZA4
A	145	SER	PHE	conflict	UNP Q9RZA4
A	311	GLU	LEU	conflict	UNP Q9RZA4
A	314	GLU	LEU	conflict	UNP Q9RZA4
A	322	LEU	ALA	conflict	UNP Q9RZA4
A	323	GLU	ASP	conflict	UNP Q9RZA4
A	324	HIS	-	expression tag	UNP Q9RZA4
A	325	HIS	-	expression tag	UNP Q9RZA4
A	326	HIS	-	expression tag	UNP Q9RZA4
A	327	HIS	-	expression tag	UNP Q9RZA4
A	328	HIS	-	expression tag	UNP Q9RZA4
A	329	HIS	-	expression tag	UNP Q9RZA4

-
- The chemical structure, labeled LBV, is a complex molecule featuring a central core with multiple substituents and functional groups. The structure is composed of several fused and linked rings, including a central benzene ring and a five-membered ring containing a nitrogen atom (N₁). The molecule is substituted with various groups, including a carboxylic acid group (CO₂H) at the bottom, a carboxylic acid group (CO₂H) on the right, and a carboxylic acid group (CO₂H) on the left. The structure also features a central nitrogen atom (N₁) and a central carbon atom (C₁). The molecule is labeled with various atoms and groups, including C₁, C₂, C₃, C₄, C₅, C₆, C₇, C₈, C₉, C₁₀, C₁₁, C₁₂, C₁₃, C₁₄, C₁₅, C₁₆, C₁₇, C₁₈, C₁₉, C₂₀, C₂₁, C₂₂, C₂₃, C₂₄, C₂₅, C₂₆, C₂₇, C₂₈, C₂₉, C₃₀, C₃₁, C₃₂, C₃₃, C₃₄, C₃₅, C₃₆, C₃₇, C₃₈, C₃₉, C₄₀, C₄₁, C₄₂, C₄₃, C₄₄, C₄₅, C₄₆, C₄₇, C₄₈, C₄₉, C₅₀, C₅₁, C₅₂, C₅₃, C₅₄, C₅₅, C₅₆, C₅₇, C₅₈, C₅₉, C₆₀, C₆₁, C₆₂, C₆₃, C₆₄, C₆₅, C₆₆, C₆₇, C₆₈, C₆₉, C₇₀, C₇₁, C₇₂, C₇₃, C₇₄, C₇₅, C₇₆, C₇₇, C₇₈, C₇₉, C₈₀, C₈₁, C₈₂, C₈₃, C₈₄, C₈₅, C₈₆, C₈₇, C₈₈, C₈₉, C₉₀, C₉₁, C₉₂, C₉₃, C₉₄, C₉₅, C₉₆, C₉₇, C₉₈, C₉₉, C₁₀₀, C₁₀₁, C₁₀₂, C₁₀₃, C₁₀₄, C₁₀₅, C₁₀₆, C₁₀₇, C₁₀₈, C₁₀₉, C₁₁₀, C₁₁₁, C₁₁₂, C₁₁₃, C₁₁₄, C₁₁₅, C₁₁₆, C₁₁₇, C₁₁₈, C₁₁₉, C₁₂₀, C₁₂₁, C₁₂₂, C₁₂₃, C₁₂₄, C₁₂₅, C₁₂₆, C₁₂₇, C₁₂₈, C₁₂₉, C₁₃₀, C₁₃₁, C₁₃₂, C₁₃₃, C₁₃₄, C₁₃₅, C₁₃₆, C₁₃₇, C₁₃₈, C₁₃₉, C₁₄₀, C₁₄₁, C₁₄₂, C₁₄₃, C₁₄₄, C₁₄₅, C₁₄₆, C₁₄₇, C₁₄₈, C₁₄₉, C₁₅₀, C₁₅₁, C₁₅₂, C₁₅₃, C₁₅₄, C₁₅₅, C₁₅₆, C₁₅₇, C₁₅₈, C₁₅₉, C₁₆₀, C₁₆₁, C₁₆₂, C₁₆₃, C₁₆₄, C₁₆₅, C₁₆₆, C₁₆₇, C₁₆₈, C₁₆₉, C₁₇₀, C₁₇₁, C₁₇₂, C₁₇₃, C₁₇₄, C₁₇₅, C₁₇₆, C₁₇₇, C₁₇₈, C₁₇₉, C₁₈₀, C₁₈₁, C₁₈₂, C₁₈₃, C₁₈₄, C₁₈₅, C₁₈₆, C₁₈₇, C₁₈₈, C₁₈₉, C₁₉₀, C₁₉₁, C₁₉₂, C₁₉₃, C₁₉₄, C₁₉₅, C₁₉₆, C₁₉₇, C₁₉₈, C₁₉₉, C₂₀₀, C₂₀₁, C₂₀₂, C₂₀₃, C₂₀₄, C₂₀₅, C₂₀₆, C₂₀₇, C₂₀₈, C₂₀₉, C₂₁₀, C₂₁₁, C₂₁₂, C₂₁₃, C₂₁₄, C₂₁₅, C₂₁₆, C₂₁₇, C₂₁₈, C₂₁₉, C₂₂₀, C₂₂₁, C₂₂₂, C₂₂₃, C₂₂₄, C₂₂₅, C₂₂₆, C₂₂₇, C₂₂₈, C₂₂₉, C₂₃₀, C₂₃₁, C₂₃₂, C₂₃₃, C₂₃₄, C₂₃₅, C₂₃₆, C₂₃₇, C₂₃₈, C₂₃₉, C₂₄₀, C₂₄₁, C₂₄₂, C₂₄₃, C₂₄₄, C₂₄₅, C₂₄₆, C₂₄₇, C₂₄₈, C₂₄₉, C₂₅₀, C₂₅₁, C₂₅₂, C₂₅₃, C₂₅₄, C₂₅₅, C₂₅₆, C₂₅₇, C₂₅₈, C₂₅₉, C₂₆₀, C₂₆₁, C₂₆₂, C₂₆₃, C₂₆₄, C₂₆₅, C₂₆₆, C₂₆₇, C₂₆₈, C₂₆₉, C₂₇₀, C₂₇₁, C₂₇₂, C₂₇₃, C₂₇₄, C₂₇₅, C₂₇₆, C₂₇₇, C₂₇₈, C₂₇₉, C₂₈₀, C₂₈₁, C₂₈₂, C₂₈₃, C₂₈₄, C₂₈₅, C₂₈₆, C₂₈₇, C₂₈₈, C₂₈₉, C₂₉₀, C₂₉₁, C₂₉₂, C₂₉₃, C₂₉₄, C₂₉₅, C₂₉₆, C₂₉₇, C₂₉₈, C₂₉₉, C₃₀₀, C₃₀₁, C₃₀₂, C₃₀₃, C₃₀₄, C₃₀₅, C₃₀₆, C₃₀₇, C₃₀₈, C₃₀₉, C₃₁₀, C₃₁₁, C₃₁₂, C₃₁₃, C₃₁₄, C₃₁₅, C₃₁₆, C₃₁₇, C₃₁₈, C₃₁₉, C₃₂₀, C₃₂₁, C₃₂₂, C₃₂₃, C₃₂₄, C₃₂₅, C₃₂₆, C₃₂₇, C₃₂₈, C₃₂₉, C₃₃₀, C₃₃₁, C₃₃₂, C₃₃₃, C₃₃₄, C₃₃₅, C₃₃₆, C₃₃₇, C₃₃₈, C₃₃₉, C₃₄₀, C₃₄₁, C₃₄₂, C₃₄₃, C₃₄₄, C₃₄₅, C₃₄₆, C₃₄₇, C₃₄₈, C₃₄₉, C₃₅₀, C₃₅₁, C₃₅₂, C₃₅₃, C₃₅₄, C₃₅₅, C₃₅₆, C₃₅₇, C₃₅₈, C₃₅₉, C₃₆₀, C₃₆₁, C₃₆₂, C₃₆₃, C₃₆₄, C₃₆₅, C₃₆₆, C₃₆₇, C₃₆₈, C₃₆₉, C₃₇₀, C₃₇₁, C₃₇₂, C₃₇₃, C₃₇₄, C₃₇₅, C₃₇₆, C₃₇₇, C₃₇₈, C₃₇₉, C₃₈₀, C₃₈₁, C₃₈₂, C₃₈₃, C₃₈₄, C₃₈₅, C₃₈₆, C₃₈₇, C₃₈₈, C₃₈₉, C₃₉₀, C₃₉₁, C₃₉₂, C₃₉₃, C₃₉₄, C₃₉₅, C₃₉₆, C₃₉₇, C₃₉₈, C₃₉₉, C₄₀₀, C<

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			43	33	4	6		

- GOL
-
- The diagram shows the skeletal structure of 1,2,3-propanetriol (glycerol). The carbon chain is represented by a zigzag line with three vertices labeled C1, C2, and C3 in green. C1 is at the left end, C2 is in the middle, and C3 is at the right end. Each carbon is bonded to a hydroxyl group (OH) in red. The OH group on C1 is labeled O1 in green below it. The OH group on C2 is labeled O2 in green below it. The OH group on C3 is labeled O3 in green to its right. The entire structure is enclosed in a black rectangular box.

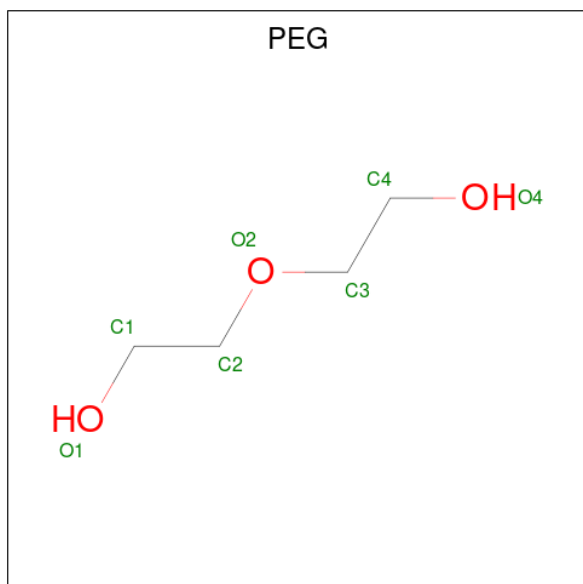
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 5 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			7	4	3		

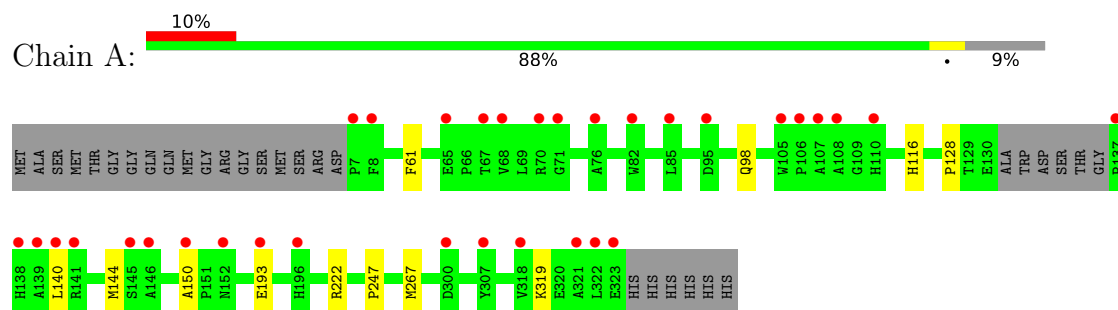
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	211	Total	O	0	0
			211	211		

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: Bacteriophytochrome



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	94.56Å 54.82Å 69.77Å 90.00° 91.80° 90.00°	Depositor
Resolution (Å)	23.71 – 1.70 23.71 – 1.70	Depositor EDS
% Data completeness (in resolution range)	95.2 (23.71-1.70) 95.2 (23.71-1.70)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.48 (at 1.71Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.205 , 0.227 (Not available) , 0.256	Depositor DCC
R_{free} test set	1903 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	20.5	Xtriage
Anisotropy	0.166	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 31.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.010 for -1/2*h+3/2*k,1/2*h+1/2*k,-l 0.005 for -1/2*h-3/2*k,-1/2*h+1/2*k,-l 0.012 for 1/2*h+3/2*k,1/2*h-1/2*k,-l 0.011 for 1/2*h-3/2*k,-1/2*h-1/2*k,-l 0.018 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2689	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.25% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, GOL, LBV, PO4

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	0.45	0/2493	0.77	0/3408

There are no bond length outliers.

There are no bond angle outliers.

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	2411	0	2420	7	0
2	A	43	0	33	0	0
3	A	12	0	16	1	0
4	A	5	0	0	1	0
5	A	7	0	10	0	0
6	A	211	0	0	0	0
All	All	2689	0	2479	7	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 1.

All (7) 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:A:247:PRO:HB2	3:A:403:GOL:H12	1.79	0.64
1:A:61:PHE:HZ	1:A:128:PRO:HD3	1.68	0.58
1:A:140:LEU:O	1:A:144:MET:HG2	2.13	0.49
1:A:98:GLN:HG2	1:A:116[A]:HIS:NE2	2.28	0.48
1:A:150:ALA:O	1:A:319:LYS:HE3	2.15	0.47
1:A:267[A]:MET:HE2	1:A:267[A]:MET:HB3	1.64	0.47
1:A:193:GLU:HG2	4:A:404:PO4:O1	2.21	0.41

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	314/341 (92%)	313 (100%)	1 (0%)	0	100	100

There are no Ramachandran outliers to report.

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	257/273 (94%)	256 (100%)	1 (0%)	84	80

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	222	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	HIS
1	A	98	GLN
1	A	282	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](#)

5 ligands are modelled in this entry.

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	$\# Z > 2$
3	GOL	A	403	-	5,5,5	0.32	0	5,5,5	0.39	0
2	LBV	A	401	1	46,46,46	2.99	18 (39%)	58,67,67	1.59	14 (24%)
3	GOL	A	402	-	5,5,5	0.15	0	5,5,5	0.56	0
5	PEG	A	405	-	6,6,6	0.44	0	5,5,5	0.30	0
4	PO4	A	404	-	4,4,4	0.94	0	6,6,6	0.49	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	PEG	A	405	-	-	2/4/4/4	-
2	LBV	A	401	1	-	6/26/74/74	0/4/4/4
3	GOL	A	402	-	-	4/4/4/4	-
3	GOL	A	403	-	-	4/4/4/4	-

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	LBV	C2A-C1A	-10.59	1.40	1.52
2	A	401	LBV	C2A-C3A	-7.26	1.42	1.51
2	A	401	LBV	C1C-C2C	-5.71	1.36	1.45
2	A	401	LBV	CBD-CAD	5.04	1.54	1.30
2	A	401	LBV	C4C-C3C	-5.03	1.34	1.45
2	A	401	LBV	C1D-C2D	-4.85	1.36	1.45
2	A	401	LBV	CAD-C3D	-4.40	1.35	1.47
2	A	401	LBV	C4A-C3A	-4.38	1.37	1.45
2	A	401	LBV	C1A-N_A	-3.10	1.33	1.37
2	A	401	LBV	C3D-C4D	-3.03	1.34	1.47
2	A	401	LBV	CMB-C2B	2.79	1.56	1.50
2	A	401	LBV	C4B-C3B	-2.75	1.36	1.42
2	A	401	LBV	CMD-C2D	2.45	1.55	1.50
2	A	401	LBV	CMC-C3C	2.39	1.55	1.50
2	A	401	LBV	C1B-C2B	-2.37	1.35	1.41
2	A	401	LBV	CAC-C2C	2.16	1.56	1.51
2	A	401	LBV	CBC-CGC	2.07	1.55	1.50
2	A	401	LBV	C1D-N_D	-2.04	1.34	1.37

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	LBV	C2C-C1C-N_C	-3.29	105.39	110.04
2	A	401	LBV	CHD-C4C-C3C	3.12	132.89	124.87
2	A	401	LBV	C3D-C4D-N_D	3.02	109.93	106.13
2	A	401	LBV	CHB-C1B-N_B	2.97	131.99	125.29
2	A	401	LBV	CHC-C4B-N_B	2.74	131.48	125.29
2	A	401	LBV	O_A-C1A-C2A	-2.69	123.89	126.74
2	A	401	LBV	C1D-N_D-C4D	-2.59	107.48	110.66
2	A	401	LBV	CHB-C1B-C2B	-2.41	121.62	127.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	LBV	C1C-C2C-C3C	2.28	109.19	106.73
2	A	401	LBV	CHC-C1C-N_C	2.24	129.31	124.60
2	A	401	LBV	CHC-C4B-C3B	-2.23	122.43	127.22
2	A	401	LBV	C1C-N_C-C4C	2.21	110.57	106.52
2	A	401	LBV	CAC-C2C-C3C	-2.10	123.94	127.87
2	A	401	LBV	CHD-C4C-N_C	-2.09	120.44	124.95

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	LBV	C3C-C4C-CHD-C1D
2	A	401	LBV	N_C-C4C-CHD-C1D
3	A	402	GOL	C1-C2-C3-O3
3	A	402	GOL	O2-C2-C3-O3
3	A	403	GOL	O1-C1-C2-C3
3	A	403	GOL	C1-C2-C3-O3
3	A	402	GOL	O1-C1-C2-C3
3	A	403	GOL	O1-C1-C2-O2
3	A	403	GOL	O2-C2-C3-O3
5	A	405	PEG	O1-C1-C2-O2
5	A	405	PEG	C4-C3-O2-C2
3	A	402	GOL	O1-C1-C2-O2
2	A	401	LBV	CAB-CBB-CGB-O2B
2	A	401	LBV	CAB-CBB-CGB-O1B
2	A	401	LBV	CAC-CBC-CGC-O1C
2	A	401	LBV	CAC-CBC-CGC-O2C

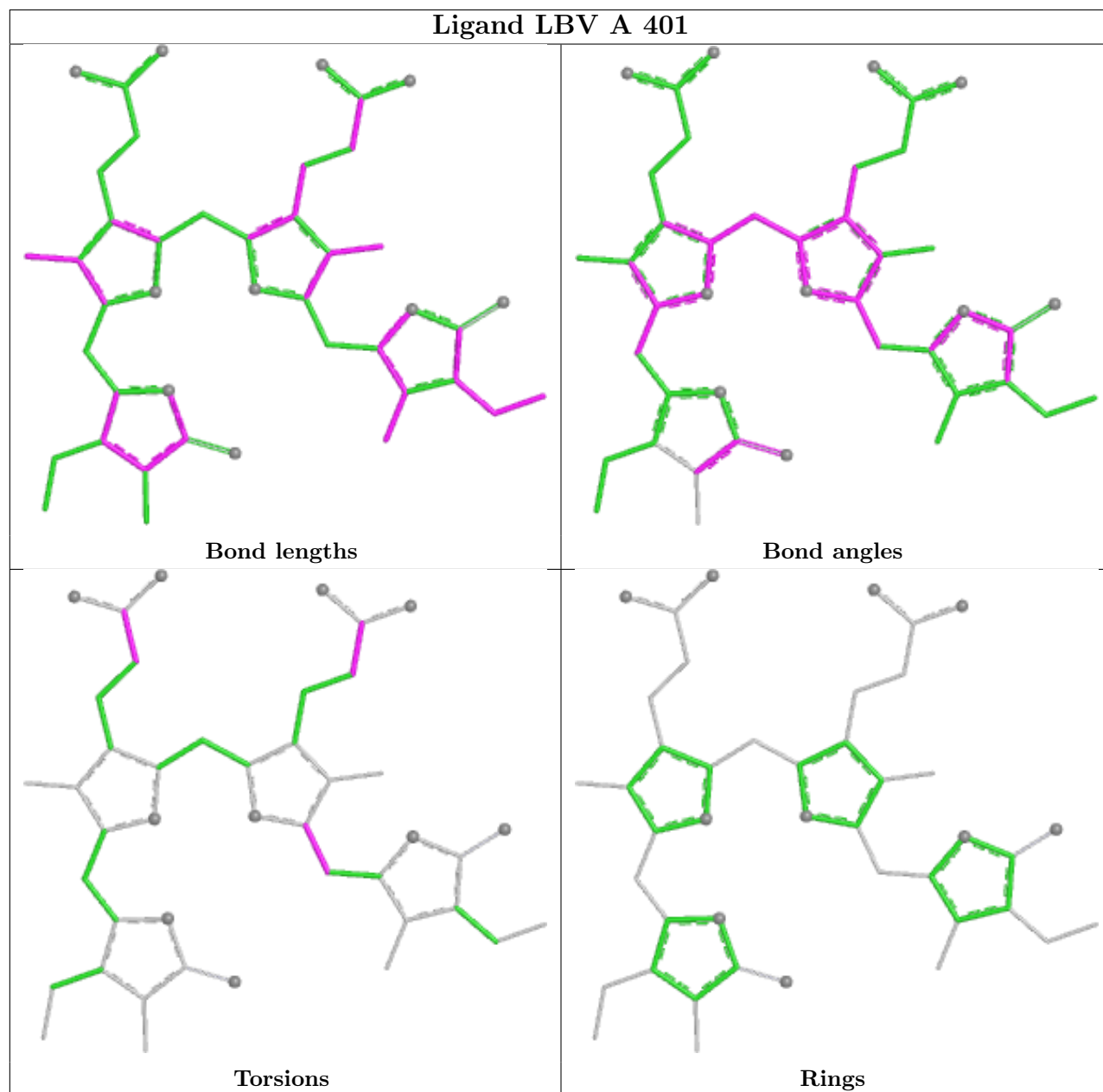
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	403	GOL	1	0
4	A	404	PO4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring

in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	311/341 (91%)	0.59	33 (10%)	11 11	10, 20, 45, 63	7 (2%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	322	LEU	5.1
1	A	138	HIS	4.8
1	A	67	THR	4.5
1	A	137	PRO	4.4
1	A	196	HIS	4.0
1	A	108	ALA	3.9
1	A	107	ALA	3.6
1	A	321	ALA	3.6
1	A	139	ALA	3.5
1	A	318	VAL	3.3
1	A	70	ARG	3.3
1	A	145	SER	3.2
1	A	307	TYR	3.1
1	A	105	TRP	3.1
1	A	110	HIS	2.9
1	A	323	GLU	2.9
1	A	76	ALA	2.7
1	A	8	PHE	2.7
1	A	95	ASP	2.7
1	A	150	ALA	2.7
1	A	106	PRO	2.7
1	A	7	PRO	2.6
1	A	193	GLU	2.6
1	A	146	ALA	2.6
1	A	141	ARG	2.5
1	A	68	VAL	2.4
1	A	152	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	65	GLU	2.3
1	A	85	LEU	2.3
1	A	300	ASP	2.2
1	A	140	LEU	2.1
1	A	71	GLY	2.1
1	A	82	TRP	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](#)

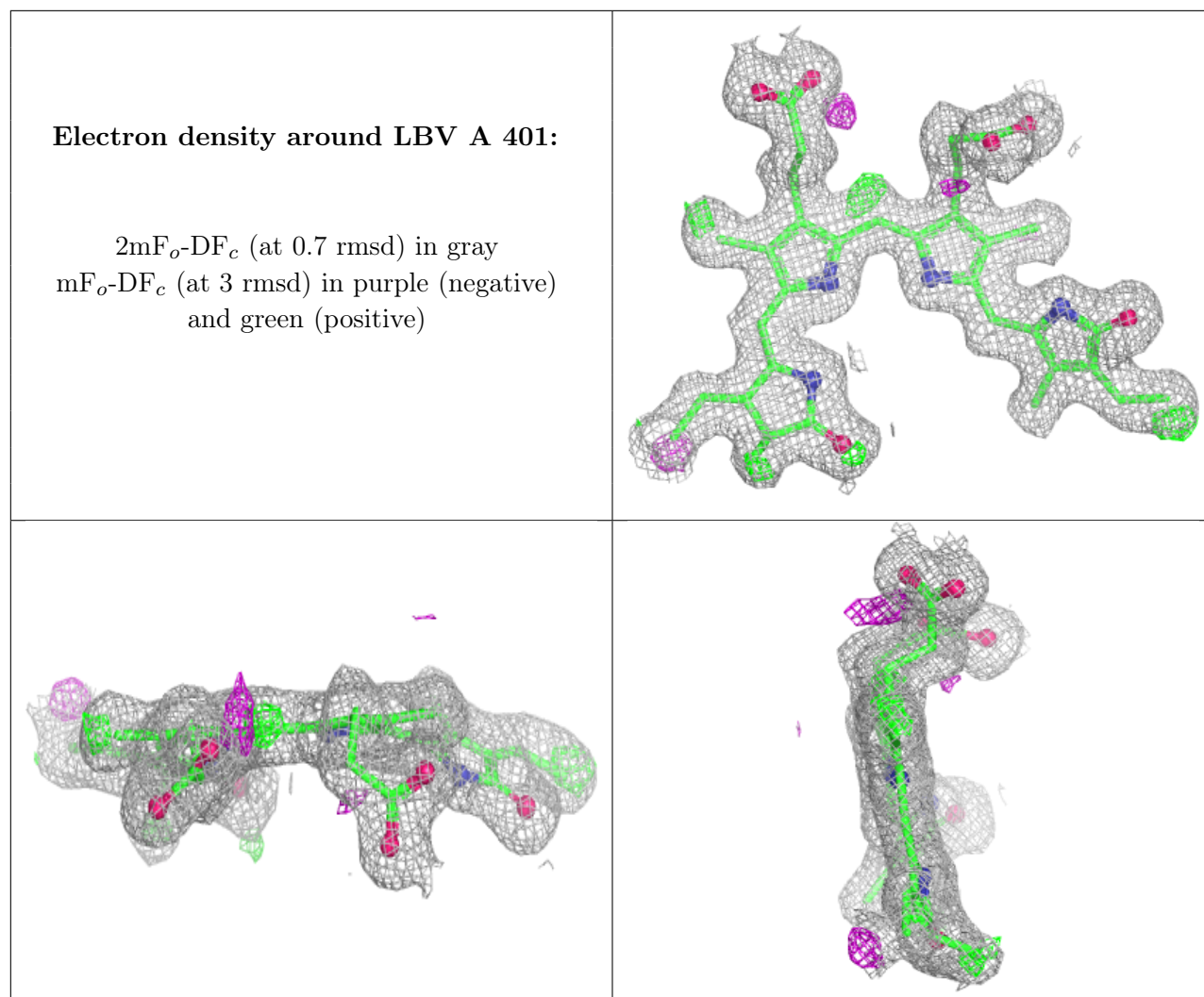
There are no oligosaccharides in this entry.

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
3	GOL	A	403	6/6	0.68	0.21	37,42,44,49	0
5	PEG	A	405	7/7	0.75	0.20	56,56,57,57	0
3	GOL	A	402	6/6	0.79	0.22	29,39,41,46	0
4	PO4	A	404	5/5	0.87	0.13	26,27,29,29	5
2	LBV	A	401	43/43	0.94	0.07	13,14,17,18	1

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers ⓘ

There are no such residues in this entry.