



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:57 AM UTC

PDB ID : 8BOR / pdb_00008bor
Title : Photosensory module from DrBphP without PHY tongue
Authors : Kurttila, M.; Takala, H.; Ihalainen, J.A.
Deposited on : 2022-11-15
Resolution : 2.30 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
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

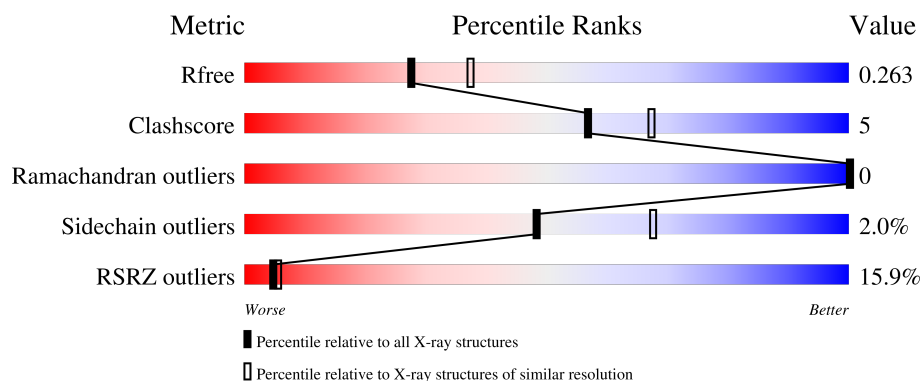
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	496	18% 81% 11% 8%
1	B	496	12% 82% 10% 7%
1	C	496	19% 83% 10% 7%
1	D	496	10% 86% 8% 6%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 14714 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
1	A	457	Total	C	N	O	S	0	3	0
			3513	2241	619	641	12			
1	B	462	Total	C	N	O	S	0	2	0
			3524	2243	623	647	11			
1	C	463	Total	C	N	O	S	0	1	0
			3538	2254	623	649	12			
1	D	466	Total	C	N	O	S	0	1	0
			3541	2253	628	649	11			

There are 224 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	initiating methionine	UNP Q9RZA4
A	-12	ALA	-	expression tag	UNP Q9RZA4
A	-11	SER	-	expression tag	UNP Q9RZA4
A	-10	MET	-	expression tag	UNP Q9RZA4
A	-9	THR	-	expression tag	UNP Q9RZA4
A	-8	GLY	-	expression tag	UNP Q9RZA4
A	-7	GLY	-	expression tag	UNP Q9RZA4
A	-6	GLN	-	expression tag	UNP Q9RZA4
A	-5	GLN	-	expression tag	UNP Q9RZA4
A	-4	MET	-	expression tag	UNP Q9RZA4
A	-3	GLY	-	expression tag	UNP Q9RZA4
A	-2	ARG	-	expression tag	UNP Q9RZA4
A	-1	GLY	-	expression tag	UNP Q9RZA4
A	0	SER	-	expression tag	UNP Q9RZA4
A	?	-	LEU	deletion	UNP Q9RZA4
A	?	-	GLU	deletion	UNP Q9RZA4
A	?	-	VAL	deletion	UNP Q9RZA4
A	?	-	ALA	deletion	UNP Q9RZA4
A	?	-	TRP	deletion	UNP Q9RZA4
A	?	-	GLY	deletion	UNP Q9RZA4
A	?	-	GLY	deletion	UNP Q9RZA4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	?	-	ALA	deletion	UNP Q9RZA4
A	?	-	THR	deletion	UNP Q9RZA4
A	?	-	PRO	deletion	UNP Q9RZA4
A	?	-	ASP	deletion	UNP Q9RZA4
A	?	-	GLN	deletion	UNP Q9RZA4
A	?	-	ALA	deletion	UNP Q9RZA4
A	?	-	LYS	deletion	UNP Q9RZA4
A	?	-	ASP	deletion	UNP Q9RZA4
A	?	-	ASP	deletion	UNP Q9RZA4
A	?	-	LEU	deletion	UNP Q9RZA4
A	?	-	GLY	deletion	UNP Q9RZA4
A	?	-	PRO	deletion	UNP Q9RZA4
A	?	-	ARG	deletion	UNP Q9RZA4
A	?	-	HIS	deletion	UNP Q9RZA4
A	?	-	SER	deletion	UNP Q9RZA4
A	?	-	PHE	deletion	UNP Q9RZA4
A	?	-	ASP	deletion	UNP Q9RZA4
A	?	-	THR	deletion	UNP Q9RZA4
A	?	-	TYR	deletion	UNP Q9RZA4
A	?	-	LEU	deletion	UNP Q9RZA4
A	?	-	GLU	deletion	UNP Q9RZA4
A	?	-	GLU	deletion	UNP Q9RZA4
A	?	-	LYS	deletion	UNP Q9RZA4
A	?	-	ARG	deletion	UNP Q9RZA4
A	474	GLY	-	linker	UNP Q9RZA4
A	475	GLY	-	linker	UNP Q9RZA4
A	476	GLY	-	linker	UNP Q9RZA4
A	477	SER	-	linker	UNP Q9RZA4
A	503	GLU	-	expression tag	UNP Q9RZA4
A	504	HIS	-	expression tag	UNP Q9RZA4
A	505	HIS	-	expression tag	UNP Q9RZA4
A	506	HIS	-	expression tag	UNP Q9RZA4
A	507	HIS	-	expression tag	UNP Q9RZA4
A	508	HIS	-	expression tag	UNP Q9RZA4
A	509	HIS	-	expression tag	UNP Q9RZA4
B	-13	MET	-	initiating methionine	UNP Q9RZA4
B	-12	ALA	-	expression tag	UNP Q9RZA4
B	-11	SER	-	expression tag	UNP Q9RZA4
B	-10	MET	-	expression tag	UNP Q9RZA4
B	-9	THR	-	expression tag	UNP Q9RZA4
B	-8	GLY	-	expression tag	UNP Q9RZA4
B	-7	GLY	-	expression tag	UNP Q9RZA4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-6	GLN	-	expression tag	UNP Q9RZA4
B	-5	GLN	-	expression tag	UNP Q9RZA4
B	-4	MET	-	expression tag	UNP Q9RZA4
B	-3	GLY	-	expression tag	UNP Q9RZA4
B	-2	ARG	-	expression tag	UNP Q9RZA4
B	-1	GLY	-	expression tag	UNP Q9RZA4
B	0	SER	-	expression tag	UNP Q9RZA4
B	?	-	LEU	deletion	UNP Q9RZA4
B	?	-	GLU	deletion	UNP Q9RZA4
B	?	-	VAL	deletion	UNP Q9RZA4
B	?	-	ALA	deletion	UNP Q9RZA4
B	?	-	TRP	deletion	UNP Q9RZA4
B	?	-	GLY	deletion	UNP Q9RZA4
B	?	-	GLY	deletion	UNP Q9RZA4
B	?	-	ALA	deletion	UNP Q9RZA4
B	?	-	THR	deletion	UNP Q9RZA4
B	?	-	PRO	deletion	UNP Q9RZA4
B	?	-	ASP	deletion	UNP Q9RZA4
B	?	-	GLN	deletion	UNP Q9RZA4
B	?	-	ALA	deletion	UNP Q9RZA4
B	?	-	LYS	deletion	UNP Q9RZA4
B	?	-	ASP	deletion	UNP Q9RZA4
B	?	-	ASP	deletion	UNP Q9RZA4
B	?	-	LEU	deletion	UNP Q9RZA4
B	?	-	GLY	deletion	UNP Q9RZA4
B	?	-	PRO	deletion	UNP Q9RZA4
B	?	-	ARG	deletion	UNP Q9RZA4
B	?	-	HIS	deletion	UNP Q9RZA4
B	?	-	SER	deletion	UNP Q9RZA4
B	?	-	PHE	deletion	UNP Q9RZA4
B	?	-	ASP	deletion	UNP Q9RZA4
B	?	-	THR	deletion	UNP Q9RZA4
B	?	-	TYR	deletion	UNP Q9RZA4
B	?	-	LEU	deletion	UNP Q9RZA4
B	?	-	GLU	deletion	UNP Q9RZA4
B	?	-	GLU	deletion	UNP Q9RZA4
B	?	-	LYS	deletion	UNP Q9RZA4
B	?	-	ARG	deletion	UNP Q9RZA4
B	474	GLY	-	linker	UNP Q9RZA4
B	475	GLY	-	linker	UNP Q9RZA4
B	476	GLY	-	linker	UNP Q9RZA4
B	477	SER	-	linker	UNP Q9RZA4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	503	GLU	-	expression tag	UNP Q9RZA4
B	504	HIS	-	expression tag	UNP Q9RZA4
B	505	HIS	-	expression tag	UNP Q9RZA4
B	506	HIS	-	expression tag	UNP Q9RZA4
B	507	HIS	-	expression tag	UNP Q9RZA4
B	508	HIS	-	expression tag	UNP Q9RZA4
B	509	HIS	-	expression tag	UNP Q9RZA4
C	-13	MET	-	initiating methionine	UNP Q9RZA4
C	-12	ALA	-	expression tag	UNP Q9RZA4
C	-11	SER	-	expression tag	UNP Q9RZA4
C	-10	MET	-	expression tag	UNP Q9RZA4
C	-9	THR	-	expression tag	UNP Q9RZA4
C	-8	GLY	-	expression tag	UNP Q9RZA4
C	-7	GLY	-	expression tag	UNP Q9RZA4
C	-6	GLN	-	expression tag	UNP Q9RZA4
C	-5	GLN	-	expression tag	UNP Q9RZA4
C	-4	MET	-	expression tag	UNP Q9RZA4
C	-3	GLY	-	expression tag	UNP Q9RZA4
C	-2	ARG	-	expression tag	UNP Q9RZA4
C	-1	GLY	-	expression tag	UNP Q9RZA4
C	0	SER	-	expression tag	UNP Q9RZA4
C	?	-	LEU	deletion	UNP Q9RZA4
C	?	-	GLU	deletion	UNP Q9RZA4
C	?	-	VAL	deletion	UNP Q9RZA4
C	?	-	ALA	deletion	UNP Q9RZA4
C	?	-	TRP	deletion	UNP Q9RZA4
C	?	-	GLY	deletion	UNP Q9RZA4
C	?	-	GLY	deletion	UNP Q9RZA4
C	?	-	ALA	deletion	UNP Q9RZA4
C	?	-	THR	deletion	UNP Q9RZA4
C	?	-	PRO	deletion	UNP Q9RZA4
C	?	-	ASP	deletion	UNP Q9RZA4
C	?	-	GLN	deletion	UNP Q9RZA4
C	?	-	ALA	deletion	UNP Q9RZA4
C	?	-	LYS	deletion	UNP Q9RZA4
C	?	-	ASP	deletion	UNP Q9RZA4
C	?	-	ASP	deletion	UNP Q9RZA4
C	?	-	LEU	deletion	UNP Q9RZA4
C	?	-	GLY	deletion	UNP Q9RZA4
C	?	-	PRO	deletion	UNP Q9RZA4
C	?	-	ARG	deletion	UNP Q9RZA4
C	?	-	HIS	deletion	UNP Q9RZA4

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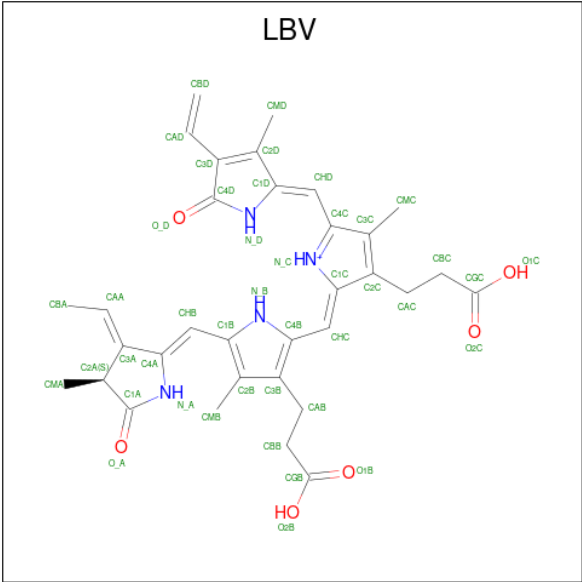
Chain	Residue	Modelled	Actual	Comment	Reference
C	?	-	SER	deletion	UNP Q9RZA4
C	?	-	PHE	deletion	UNP Q9RZA4
C	?	-	ASP	deletion	UNP Q9RZA4
C	?	-	THR	deletion	UNP Q9RZA4
C	?	-	TYR	deletion	UNP Q9RZA4
C	?	-	LEU	deletion	UNP Q9RZA4
C	?	-	GLU	deletion	UNP Q9RZA4
C	?	-	GLU	deletion	UNP Q9RZA4
C	?	-	LYS	deletion	UNP Q9RZA4
C	?	-	ARG	deletion	UNP Q9RZA4
C	474	GLY	-	linker	UNP Q9RZA4
C	475	GLY	-	linker	UNP Q9RZA4
C	476	GLY	-	linker	UNP Q9RZA4
C	477	SER	-	linker	UNP Q9RZA4
C	503	GLU	-	expression tag	UNP Q9RZA4
C	504	HIS	-	expression tag	UNP Q9RZA4
C	505	HIS	-	expression tag	UNP Q9RZA4
C	506	HIS	-	expression tag	UNP Q9RZA4
C	507	HIS	-	expression tag	UNP Q9RZA4
C	508	HIS	-	expression tag	UNP Q9RZA4
C	509	HIS	-	expression tag	UNP Q9RZA4
D	-13	MET	-	initiating methionine	UNP Q9RZA4
D	-12	ALA	-	expression tag	UNP Q9RZA4
D	-11	SER	-	expression tag	UNP Q9RZA4
D	-10	MET	-	expression tag	UNP Q9RZA4
D	-9	THR	-	expression tag	UNP Q9RZA4
D	-8	GLY	-	expression tag	UNP Q9RZA4
D	-7	GLY	-	expression tag	UNP Q9RZA4
D	-6	GLN	-	expression tag	UNP Q9RZA4
D	-5	GLN	-	expression tag	UNP Q9RZA4
D	-4	MET	-	expression tag	UNP Q9RZA4
D	-3	GLY	-	expression tag	UNP Q9RZA4
D	-2	ARG	-	expression tag	UNP Q9RZA4
D	-1	GLY	-	expression tag	UNP Q9RZA4
D	0	SER	-	expression tag	UNP Q9RZA4
D	?	-	LEU	deletion	UNP Q9RZA4
D	?	-	GLU	deletion	UNP Q9RZA4
D	?	-	VAL	deletion	UNP Q9RZA4
D	?	-	ALA	deletion	UNP Q9RZA4
D	?	-	TRP	deletion	UNP Q9RZA4
D	?	-	GLY	deletion	UNP Q9RZA4
D	?	-	GLY	deletion	UNP Q9RZA4

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Chain	Residue	Modelled	Actual	Comment	Reference
D	?	-	ALA	deletion	UNP Q9RZA4
D	?	-	THR	deletion	UNP Q9RZA4
D	?	-	PRO	deletion	UNP Q9RZA4
D	?	-	ASP	deletion	UNP Q9RZA4
D	?	-	GLN	deletion	UNP Q9RZA4
D	?	-	ALA	deletion	UNP Q9RZA4
D	?	-	LYS	deletion	UNP Q9RZA4
D	?	-	ASP	deletion	UNP Q9RZA4
D	?	-	ASP	deletion	UNP Q9RZA4
D	?	-	LEU	deletion	UNP Q9RZA4
D	?	-	GLY	deletion	UNP Q9RZA4
D	?	-	PRO	deletion	UNP Q9RZA4
D	?	-	ARG	deletion	UNP Q9RZA4
D	?	-	HIS	deletion	UNP Q9RZA4
D	?	-	SER	deletion	UNP Q9RZA4
D	?	-	PHE	deletion	UNP Q9RZA4
D	?	-	ASP	deletion	UNP Q9RZA4
D	?	-	THR	deletion	UNP Q9RZA4
D	?	-	TYR	deletion	UNP Q9RZA4
D	?	-	LEU	deletion	UNP Q9RZA4
D	?	-	GLU	deletion	UNP Q9RZA4
D	?	-	GLU	deletion	UNP Q9RZA4
D	?	-	LYS	deletion	UNP Q9RZA4
D	?	-	ARG	deletion	UNP Q9RZA4
D	474	GLY	-	linker	UNP Q9RZA4
D	475	GLY	-	linker	UNP Q9RZA4
D	476	GLY	-	linker	UNP Q9RZA4
D	477	SER	-	linker	UNP Q9RZA4
D	503	GLU	-	expression tag	UNP Q9RZA4
D	504	HIS	-	expression tag	UNP Q9RZA4
D	505	HIS	-	expression tag	UNP Q9RZA4
D	506	HIS	-	expression tag	UNP Q9RZA4
D	507	HIS	-	expression tag	UNP Q9RZA4
D	508	HIS	-	expression tag	UNP Q9RZA4
D	509	HIS	-	expression tag	UNP Q9RZA4

- Molecule 2 is 3-[2-[(Z)-[3-(2-carboxyethyl)-5-[(Z)-(4-ethenyl-3-methyl-5-oxidanylidene-pyrro-
l-2-ylidene)methyl]-4-methyl-pyrrol-1-ium -2-ylidene]methyl]-5-[(Z)-[(3E)-3-ethylidene-4-me-
thyl-5-oxidanylidene-pyrrolidin-2-ylidene]methyl]-4-methyl-1H-pyrrol-3-yl]propanoic acid
(CCD ID: LBV) (formula: C₃₃H₃₇N₄O₆) (labeled as "Ligand of Interest" by depositor).



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

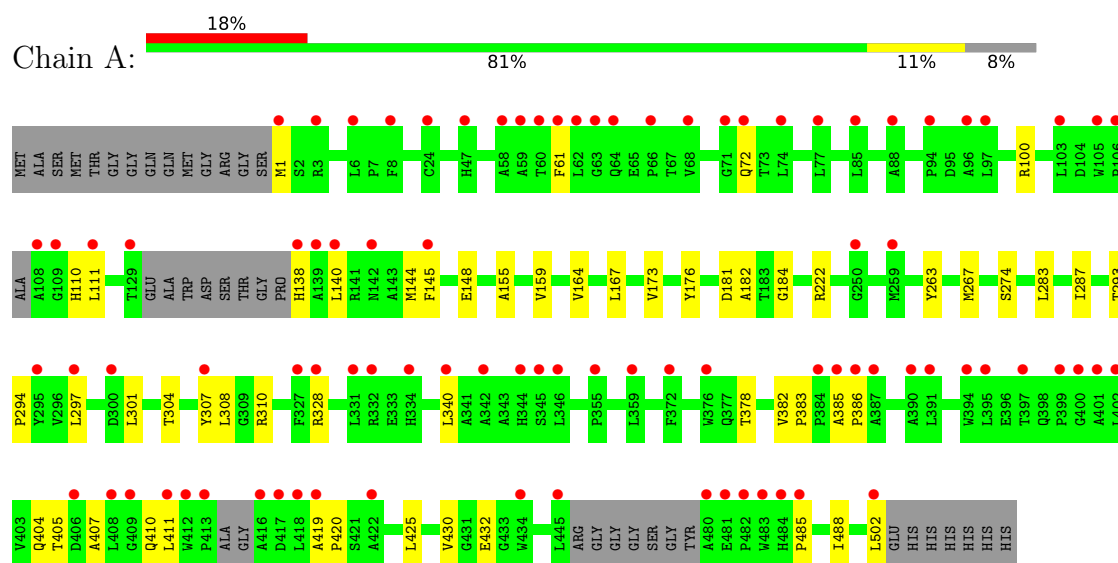
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	100	Total	O	0	0
			100	100		
3	B	94	Total	O	0	0
			94	94		
3	C	104	Total	O	0	0
			104	104		
3	D	128	Total	O	0	0
			128	128		

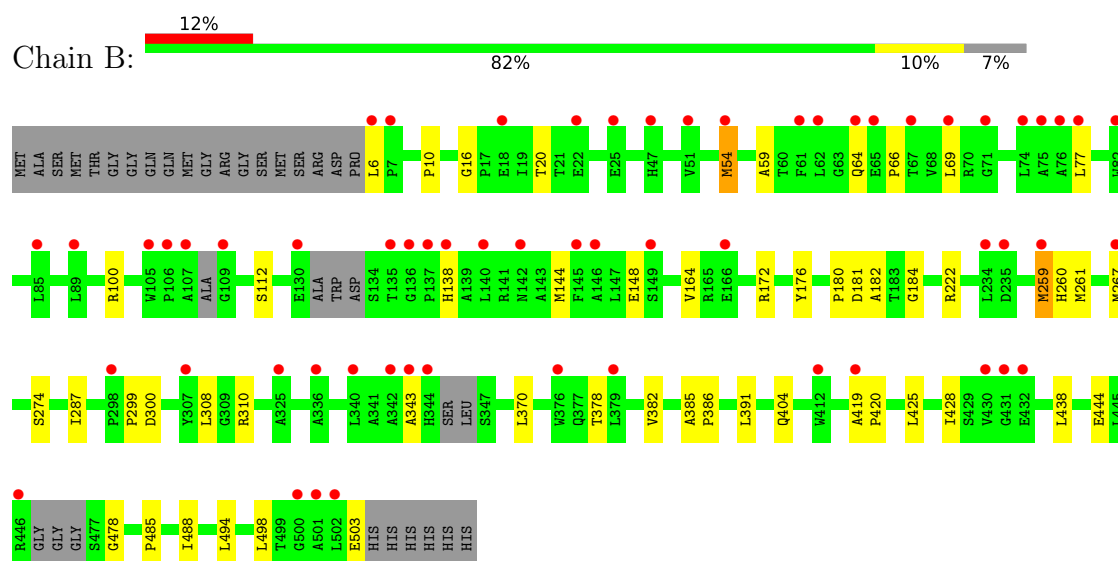
3 Residue-property plots [i](#)

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.

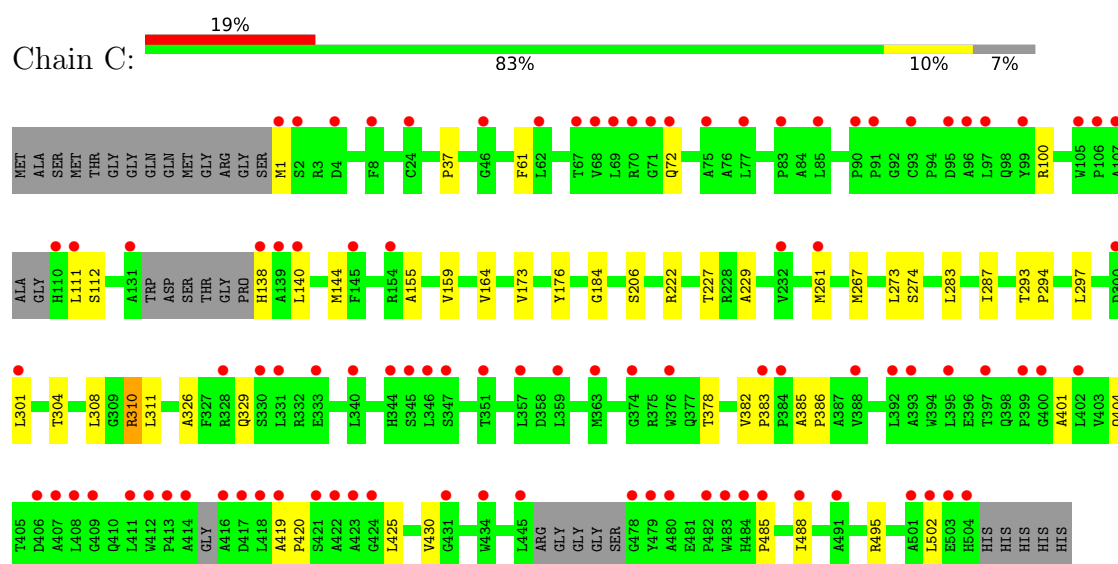
• Molecule 1: Bacteriophytochrome



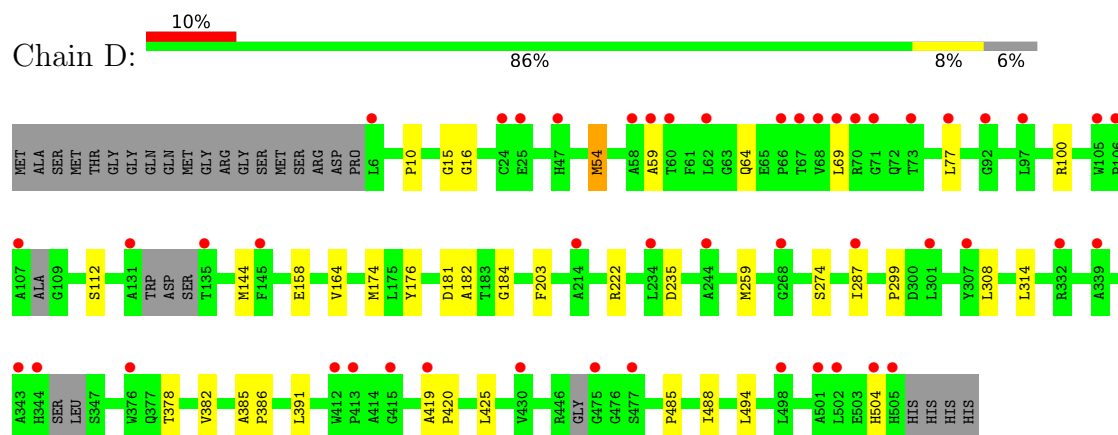
• Molecule 1: Bacteriophytochrome



• Molecule 1: Bacteriophytochrome



• Molecule 1: Bacteriophytochrome



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	119.63Å 64.59Å 131.04Å 90.00° 91.55° 90.00°	Depositor
Resolution (Å)	45.99 – 2.30 45.99 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.8 (45.99-2.30) 98.8 (45.99-2.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0352	Depositor
R, R_{free}	0.235 , 0.268 (Not available) , 0.263	Depositor DCC
R_{free} test set	4418 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	45.1	Xtriage
Anisotropy	0.312	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 32.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.046 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	14714	wwPDB-VP
Average B, all atoms (Å ²)	53.0	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 29.26 % 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 1.5991e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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](#)

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

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	0.52	0/3606	0.86	0/4934
1	B	0.52	0/3614	0.87	0/4946
1	C	0.52	0/3629	0.87	0/4967
1	D	0.51	0/3632	0.86	0/4970
All	All	0.52	0/14481	0.87	0/19817

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 [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	3513	0	3497	37	0
1	B	3524	0	3496	42	0
1	C	3538	0	3511	33	0
1	D	3541	0	3510	30	0
2	A	43	0	33	0	0
2	B	43	0	33	1	0
2	C	43	0	34	2	0
2	D	43	0	33	3	0
3	A	100	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	94	0	0	1	0
3	C	104	0	0	1	0
3	D	128	0	0	2	0
All	All	14714	0	14147	132	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 5.

All (132) 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:326:ALA:O	1:C:329:GLN:HG3	1.79	0.83
1:B:54:MET:HE2	1:B:66:PRO:HA	1.64	0.79
1:D:54:MET:N	1:D:54:MET:SD	2.59	0.76
1:B:370:LEU:HD23	1:B:438[B]:LEU:CD2	2.16	0.75
1:B:100:ARG:HH22	1:B:299:PRO:HB3	1.54	0.72
1:D:158:GLU:HG2	3:D:1015:HOH:O	1.88	0.71
1:D:100:ARG:HH22	1:D:299:PRO:HB3	1.56	0.70
1:B:54:MET:HE1	1:B:69:LEU:HD12	1.74	0.69
1:A:138:HIS:CE1	1:B:100:ARG:HG3	2.28	0.69
1:D:100:ARG:HH22	1:D:299:PRO:CB	2.07	0.67
1:B:100:ARG:HH22	1:B:299:PRO:CB	2.07	0.67
1:B:378:THR:HB	1:B:382:VAL:HG11	1.78	0.66
1:B:370:LEU:HD23	1:B:438[B]:LEU:HD22	1.77	0.66
1:D:59:ALA:HA	1:D:64:GLN:O	1.96	0.65
1:C:297:LEU:HD11	1:C:301:LEU:HD23	1.79	0.65
1:A:1:MET:HE3	3:A:978:HOH:O	1.95	0.65
1:A:297:LEU:HD11	1:A:301:LEU:HD23	1.79	0.65
1:B:59:ALA:HA	1:B:64:GLN:O	1.97	0.65
1:B:261:MET:HE3	3:B:979:HOH:O	1.98	0.63
1:D:259:MET:HE2	2:D:800:LBV:HMA2	1.79	0.63
1:B:428:ILE:HD11	1:B:438[B]:LEU:HD12	1.81	0.62
1:B:54:MET:CE	1:B:66:PRO:HA	2.31	0.60
1:C:37:PRO:HG3	1:C:229:ALA:HB1	1.82	0.60
1:C:227:THR:HA	1:C:261:MET:HE3	1.84	0.60
1:C:310:ARG:NH1	3:C:902:HOH:O	2.34	0.59
1:B:54:MET:HE1	1:B:69:LEU:HB2	1.84	0.59
1:B:54:MET:HE1	1:B:69:LEU:CD1	2.33	0.58
1:C:293:THR:HG23	1:C:294:PRO:HD2	1.85	0.58
1:B:172:ARG:NH1	1:D:15:GLY:O	2.36	0.57
1:A:100:ARG:HB2	1:B:138:HIS:CE1	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:378:THR:HB	1:D:382:VAL:HG11	1.87	0.55
1:A:485:PRO:HA	1:A:488:ILE:HD12	1.89	0.55
1:A:293:THR:HG23	1:A:294:PRO:HD2	1.88	0.55
1:A:100:ARG:HD3	1:B:138:HIS:CD2	2.42	0.55
1:B:485:PRO:HA	1:B:488:ILE:HD12	1.89	0.55
1:D:485:PRO:HA	1:D:488:ILE:HD12	1.89	0.54
1:B:274:SER:HA	1:B:287:ILE:O	2.07	0.54
1:C:485:PRO:HA	1:C:488:ILE:HD12	1.89	0.54
1:D:259:MET:CE	2:D:800:LBV:HMA2	2.36	0.54
1:C:401:ALA:HB1	1:C:495:ARG:NE	2.23	0.54
1:D:274:SER:HA	1:D:287:ILE:O	2.08	0.54
1:B:69:LEU:CD2	1:B:77:LEU:HD21	2.38	0.53
1:D:176:TYR:CZ	1:D:184:GLY:HA3	2.44	0.53
1:C:140:LEU:HD22	1:C:304:THR:HG23	1.91	0.53
1:A:140:LEU:HD22	1:A:304:THR:HG23	1.91	0.53
1:A:176:TYR:CZ	1:A:184:GLY:HA3	2.44	0.52
1:B:176:TYR:CZ	1:B:184:GLY:HA3	2.44	0.52
1:B:10:PRO:HG2	1:B:16:GLY:HA2	1.91	0.52
1:C:274:SER:HA	1:C:287:ILE:O	2.09	0.52
1:C:176:TYR:CZ	1:C:184:GLY:HA3	2.45	0.52
1:C:138:HIS:CE1	1:D:100:ARG:HG3	2.44	0.52
1:A:274:SER:HA	1:A:287:ILE:O	2.10	0.52
1:A:145:PHE:CE1	1:B:310:ARG:NH1	2.78	0.51
1:D:10:PRO:HG2	1:D:16:GLY:HA2	1.91	0.51
1:C:206:SER:O	2:C:800:LBV:HBA3	2.11	0.50
1:C:297:LEU:CD1	1:C:301:LEU:HD23	2.41	0.50
1:A:138:HIS:CD2	1:B:100:ARG:HD3	2.47	0.50
1:A:382:VAL:HG13	1:A:383:PRO:HD2	1.93	0.50
1:A:297:LEU:CD1	1:A:301:LEU:HD23	2.42	0.50
1:C:382:VAL:HG13	1:C:383:PRO:HD2	1.92	0.49
1:D:382:VAL:HG22	3:D:947:HOH:O	2.12	0.49
1:B:54:MET:HE1	1:B:69:LEU:CB	2.43	0.49
1:B:260:HIS:NE2	2:B:800:LBV:HBC1	2.28	0.48
1:B:20:THR:HA	1:B:259:MET:HE2	1.94	0.48
1:C:267:MET:HE1	2:C:800:LBV:HBD1	1.95	0.48
1:A:140:LEU:HD21	1:A:307:TYR:CD2	2.48	0.48
1:A:378:THR:HB	1:A:382:VAL:HG21	1.96	0.48
1:A:407:ALA:O	1:A:410:GLN:HG2	2.14	0.47
1:D:391:LEU:C	1:D:391:LEU:HD23	2.39	0.47
1:C:430:VAL:HG11	1:C:502:LEU:HD12	1.95	0.47
1:C:100:ARG:NH1	1:C:112:SER:OG	2.46	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:382:VAL:CG1	1:A:383:PRO:HD2	2.45	0.47
1:B:100:ARG:NH2	1:B:112:SER:OG	2.48	0.47
1:C:164:VAL:HG22	1:C:308:LEU:HD13	1.96	0.47
1:D:100:ARG:NH2	1:D:112:SER:OG	2.48	0.47
1:C:378:THR:HB	1:C:382:VAL:HG21	1.97	0.46
1:A:61[B]:PHE:CE1	1:A:111:LEU:HD11	2.49	0.46
1:B:391:LEU:HD23	1:B:391:LEU:C	2.40	0.46
1:A:61[A]:PHE:CE1	1:A:111:LEU:HD11	2.49	0.46
1:D:419:ALA:N	1:D:420:PRO:CD	2.78	0.46
1:A:61[C]:PHE:CE1	1:A:111:LEU:HD11	2.50	0.46
1:B:444:GLU:OE1	1:B:478:GLY:N	2.44	0.46
1:C:37:PRO:CG	1:C:229:ALA:HB1	2.45	0.46
1:B:164:VAL:HG22	1:B:308:LEU:HD13	1.98	0.46
1:C:382:VAL:CG1	1:C:383:PRO:HD2	2.45	0.46
1:D:203:PHE:CZ	2:D:800:LBV:HAD1	2.51	0.46
1:D:164:VAL:HG22	1:D:308:LEU:HD13	1.97	0.45
1:B:308:LEU:O	1:B:308:LEU:HD23	2.17	0.45
1:C:326:ALA:C	1:C:329:GLN:HG3	2.42	0.45
1:D:69:LEU:CD2	1:D:77:LEU:HD21	2.46	0.45
1:B:419:ALA:N	1:B:420:PRO:CD	2.80	0.45
1:A:419:ALA:N	1:A:420:PRO:CD	2.80	0.44
1:C:61:PHE:CE1	1:C:111:LEU:HD11	2.52	0.44
1:C:385:ALA:HB3	1:C:386:PRO:HD3	2.00	0.44
1:C:419:ALA:N	1:C:420:PRO:CD	2.80	0.44
1:A:164:VAL:HG22	1:A:308:LEU:HD13	2.00	0.44
1:D:308:LEU:HD23	1:D:308:LEU:O	2.16	0.44
1:B:267:MET:O	1:D:10:PRO:HG3	2.18	0.44
1:B:385:ALA:HB3	1:B:386:PRO:HD3	1.99	0.44
1:C:311:LEU:HD22	1:D:314:LEU:HD12	1.99	0.44
1:D:100:ARG:NH2	1:D:299:PRO:HB3	2.28	0.44
1:A:308:LEU:O	1:A:308:LEU:HD23	2.18	0.43
1:A:410:GLN:HG3	1:A:411:LEU:N	2.33	0.43
1:B:180:PRO:HG2	1:B:404:GLN:CD	2.42	0.43
1:C:308:LEU:HD23	1:C:308:LEU:O	2.18	0.43
1:D:181:ASP:O	1:D:182:ALA:HB3	2.19	0.43
1:A:1:MET:HE2	1:A:283:LEU:O	2.19	0.43
1:A:385:ALA:HB3	1:A:386:PRO:HD3	1.99	0.43
1:A:405:THR:HB	1:A:411:LEU:HD12	2.01	0.43
1:B:20:THR:CA	1:B:259:MET:HE2	2.48	0.43
1:B:425:LEU:C	1:B:425:LEU:HD23	2.44	0.43
1:C:1:MET:HE2	1:C:283:LEU:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:273:LEU:HD13	1:C:297:LEU:HD22	2.01	0.42
1:B:100:ARG:NH2	1:B:299:PRO:HB3	2.27	0.42
1:D:385:ALA:HB3	1:D:386:PRO:HD3	2.01	0.42
1:D:425:LEU:C	1:D:425:LEU:HD23	2.45	0.42
1:A:263:TYR:CZ	1:A:267:MET:HE3	2.55	0.42
1:A:430:VAL:HG11	1:A:502:LEU:HD12	2.02	0.42
1:C:425:LEU:C	1:C:425:LEU:HD23	2.45	0.42
1:A:181:ASP:O	1:A:182:ALA:HB3	2.19	0.41
1:A:425:LEU:C	1:A:425:LEU:HD23	2.45	0.41
1:D:54:MET:SD	1:D:235:ASP:O	2.78	0.41
1:A:164:VAL:HB	1:A:173:VAL:HG11	2.02	0.41
1:B:181:ASP:O	1:B:182:ALA:HB3	2.20	0.41
1:C:155:ALA:O	1:C:159:VAL:HG23	2.20	0.41
1:A:140:LEU:HD21	1:A:307:TYR:HD2	1.85	0.41
1:A:155:ALA:O	1:A:159:VAL:HG23	2.21	0.41
1:A:328:ARG:HG3	1:A:328:ARG:HH11	1.85	0.41
1:B:54:MET:CE	1:B:69:LEU:HB2	2.49	0.41
1:D:174:MET:HE3	1:D:174:MET:HB2	1.84	0.41
1:A:340:LEU:HD23	1:B:343:ALA:O	2.22	0.40
1:C:164:VAL:HB	1:C:173:VAL:HG11	2.03	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	451/496 (91%)	445 (99%)	6 (1%)	0	100	100
1	B	454/496 (92%)	447 (98%)	7 (2%)	0	100	100
1	C	454/496 (92%)	445 (98%)	9 (2%)	0	100	100
1	D	457/496 (92%)	449 (98%)	8 (2%)	0	100	100
All	All	1816/1984 (92%)	1786 (98%)	30 (2%)	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	365/386 (95%)	356 (98%)	9 (2%)	42	60
1	B	364/386 (94%)	354 (97%)	10 (3%)	39	58
1	C	366/386 (95%)	361 (99%)	5 (1%)	59	76
1	D	364/386 (94%)	359 (99%)	5 (1%)	59	76
All	All	1459/1544 (94%)	1430 (98%)	29 (2%)	48	67

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	110	HIS
1	A	144	MET
1	A	148	GLU
1	A	167	LEU
1	A	222	ARG
1	A	310	ARG
1	A	404	GLN
1	A	432	GLU
1	B	6	LEU
1	B	54	MET
1	B	144	MET
1	B	148	GLU
1	B	222	ARG
1	B	259	MET
1	B	300	ASP
1	B	494	LEU
1	B	498	LEU
1	B	503	GLU
1	C	72	GLN
1	C	144	MET
1	C	222	ARG

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Mol	Chain	Res	Type
1	C	310	ARG
1	C	404	GLN
1	D	54	MET
1	D	144	MET
1	D	222	ARG
1	D	494	LEU
1	D	504	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	86	GLN
1	A	377	GLN
1	B	86	GLN
1	B	98	GLN
1	B	110	HIS
1	B	282	GLN
1	C	86	GLN
1	C	377	GLN
1	D	142	ASN
1	D	243	ASN
1	D	282	GLN
1	D	344	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

4 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	$\# Z > 2$
2	LBV	C	800	1	46,46,46	0.73	0	58,67,67	1.26	7 (12%)
2	LBV	B	800	1	46,46,46	0.84	0	58,67,67	1.31	6 (10%)
2	LBV	D	800	1	46,46,46	0.91	0	58,67,67	1.52	6 (10%)
2	LBV	A	800	1	46,46,46	0.81	0	58,67,67	1.42	6 (10%)

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
2	LBV	C	800	1	-	11/26/74/74	0/4/4/4
2	LBV	B	800	1	-	11/26/74/74	0/4/4/4
2	LBV	D	800	1	-	9/26/74/74	0/4/4/4
2	LBV	A	800	1	-	12/26/74/74	0/4/4/4

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	800	LBV	CAA-C3A-C4A	-6.03	119.08	126.36
2	A	800	LBV	CAA-C3A-C4A	-4.36	121.09	126.36
2	A	800	LBV	CHD-C4C-N_C	-4.35	115.56	124.95
2	A	800	LBV	CHD-C4C-C3C	4.13	135.47	124.87
2	B	800	LBV	CAA-C3A-C4A	-4.05	121.47	126.36
2	B	800	LBV	CHD-C4C-C3C	3.88	134.83	124.87
2	B	800	LBV	CHD-C4C-N_C	-3.80	116.75	124.95
2	D	800	LBV	CHD-C4C-N_C	-3.61	117.15	124.95
2	D	800	LBV	CHD-C4C-C3C	3.58	134.06	124.87
2	C	800	LBV	CHD-C4C-N_C	-3.43	117.54	124.95
2	C	800	LBV	CHD-C4C-C3C	3.39	133.57	124.87
2	D	800	LBV	C2C-C1C-N_C	-3.02	105.76	110.04

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	800	LBV	O1C-CGC-CBC	3.02	123.54	114.00
2	C	800	LBV	C2C-C1C-N_C	-2.84	106.03	110.04
2	C	800	LBV	CAA-C3A-C4A	-2.82	122.96	126.36
2	B	800	LBV	C2C-C1C-N_C	-2.77	106.12	110.04
2	A	800	LBV	C2C-C1C-N_C	-2.32	106.76	110.04
2	B	800	LBV	O1C-CGC-CBC	2.29	121.25	114.00
2	D	800	LBV	C1C-N_C-C4C	2.23	110.61	106.52
2	A	800	LBV	O_A-C1A-N_A	-2.22	122.32	124.93
2	A	800	LBV	O1C-CGC-CBC	2.22	121.00	114.00
2	C	800	LBV	C4B-CHC-C1C	2.17	132.96	128.22
2	C	800	LBV	CMC-C3C-C4C	2.10	128.36	125.10
2	B	800	LBV	C1C-N_C-C4C	2.04	110.27	106.52
2	C	800	LBV	CMD-C2D-C1D	-2.03	121.69	124.16

There are no chirality outliers.

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	800	LBV	C2A-C3A-CAA-CBA
2	A	800	LBV	C4A-C3A-CAA-CBA
2	A	800	LBV	N_C-C4C-CHD-C1D
2	B	800	LBV	C2A-C3A-CAA-CBA
2	B	800	LBV	C4A-C3A-CAA-CBA
2	B	800	LBV	N_C-C4C-CHD-C1D
2	C	800	LBV	C4A-C3A-CAA-CBA
2	C	800	LBV	N_C-C4C-CHD-C1D
2	D	800	LBV	C2A-C3A-CAA-CBA
2	D	800	LBV	C4A-C3A-CAA-CBA
2	D	800	LBV	N_C-C4C-CHD-C1D
2	A	800	LBV	C3C-C4C-CHD-C1D
2	B	800	LBV	C3C-C4C-CHD-C1D
2	C	800	LBV	C3C-C4C-CHD-C1D
2	D	800	LBV	C3C-C4C-CHD-C1D
2	A	800	LBV	N_A-C4A-CHB-C1B
2	C	800	LBV	N_A-C4A-CHB-C1B
2	C	800	LBV	N_D-C1D-CHD-C4C
2	C	800	LBV	C2D-C1D-CHD-C4C
2	A	800	LBV	N_D-C1D-CHD-C4C
2	C	800	LBV	C2C-CAC-CBC-CGC
2	A	800	LBV	C2D-C1D-CHD-C4C
2	B	800	LBV	N_D-C1D-CHD-C4C
2	C	800	LBV	CAC-CBC-CGC-O2C

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Mol	Chain	Res	Type	Atoms
2	A	800	LBV	C3A-C4A-CHB-C1B
2	C	800	LBV	CAB-CBB-CGB-O1B
2	C	800	LBV	CAB-CBB-CGB-O2B
2	C	800	LBV	CAC-CBC-CGC-O1C
2	B	800	LBV	CAB-CBB-CGB-O2B
2	B	800	LBV	C2C-CAC-CBC-CGC
2	B	800	LBV	CAB-CBB-CGB-O1B
2	A	800	LBV	CAB-CBB-CGB-O2B
2	A	800	LBV	CAB-CBB-CGB-O1B
2	B	800	LBV	CAC-CBC-CGC-O1C
2	A	800	LBV	CAC-CBC-CGC-O1C
2	D	800	LBV	CAC-CBC-CGC-O1C
2	A	800	LBV	CAC-CBC-CGC-O2C
2	B	800	LBV	C2D-C1D-CHD-C4C
2	B	800	LBV	CAC-CBC-CGC-O2C
2	D	800	LBV	CAC-CBC-CGC-O2C
2	D	800	LBV	CAB-CBB-CGB-O1B
2	D	800	LBV	CAB-CBB-CGB-O2B
2	D	800	LBV	N_D-C1D-CHD-C4C

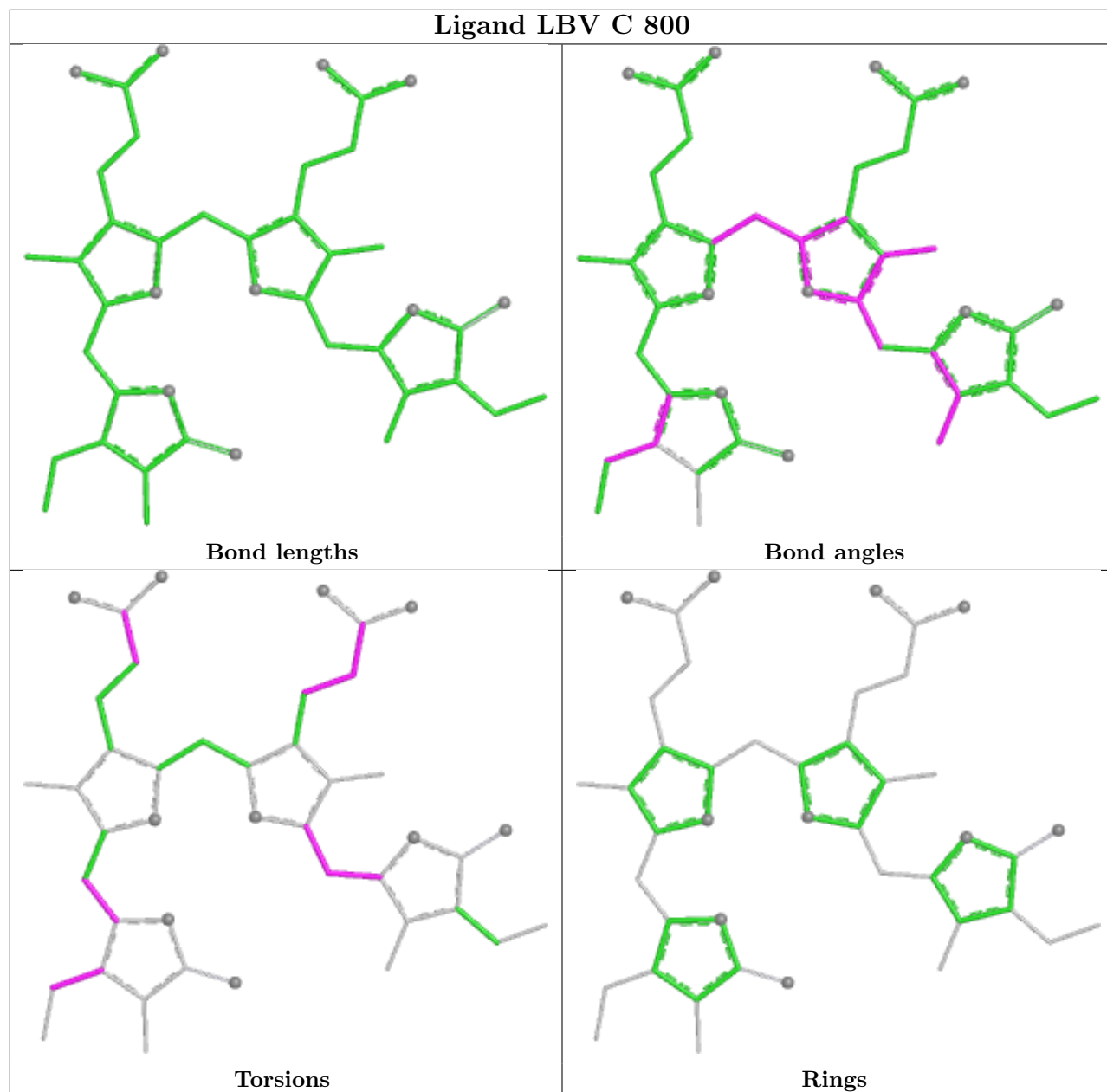
There are no ring outliers.

3 monomers are involved in 6 short contacts:

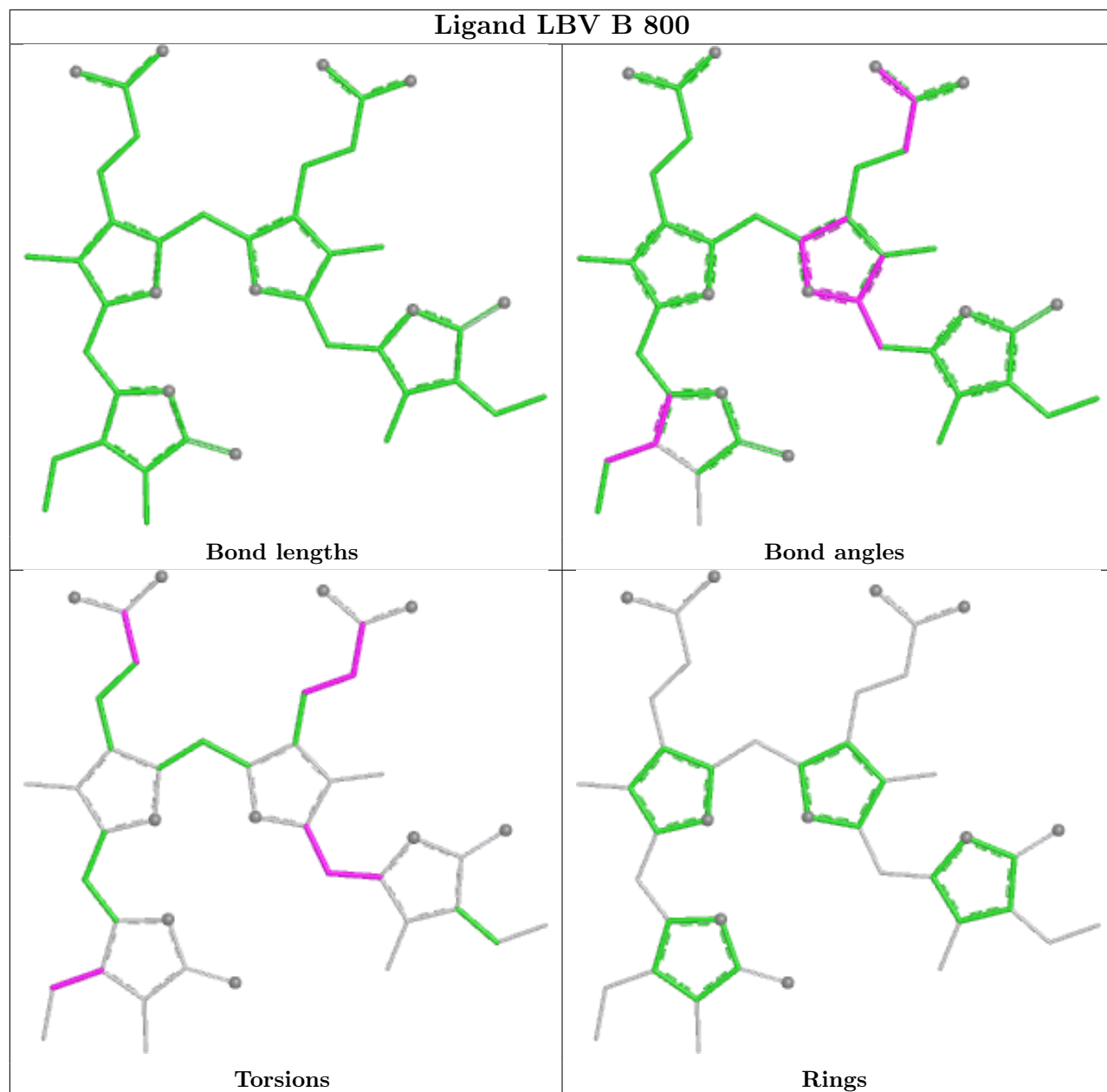
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	800	LBV	2	0
2	B	800	LBV	1	0
2	D	800	LBV	3	0

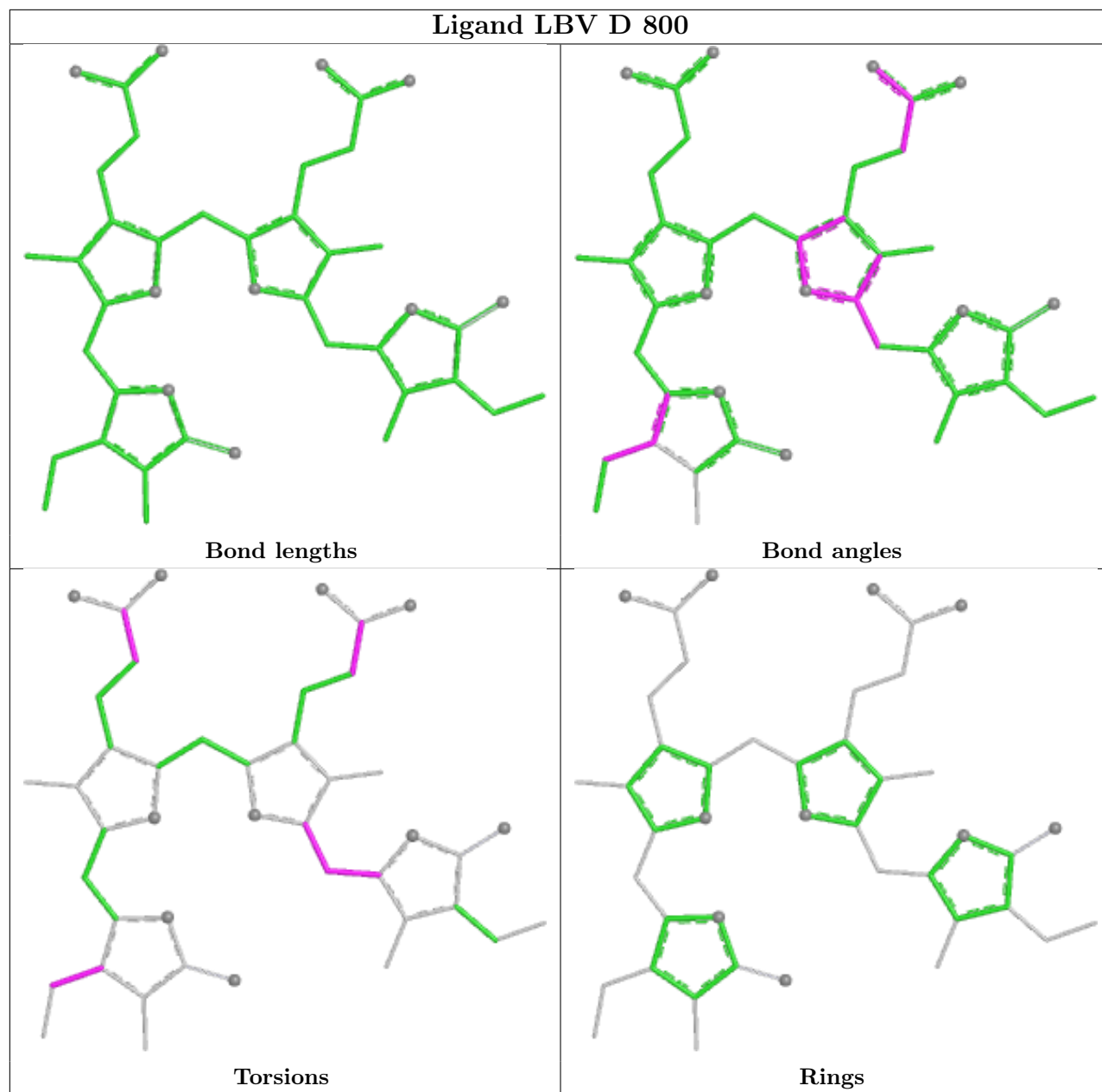
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

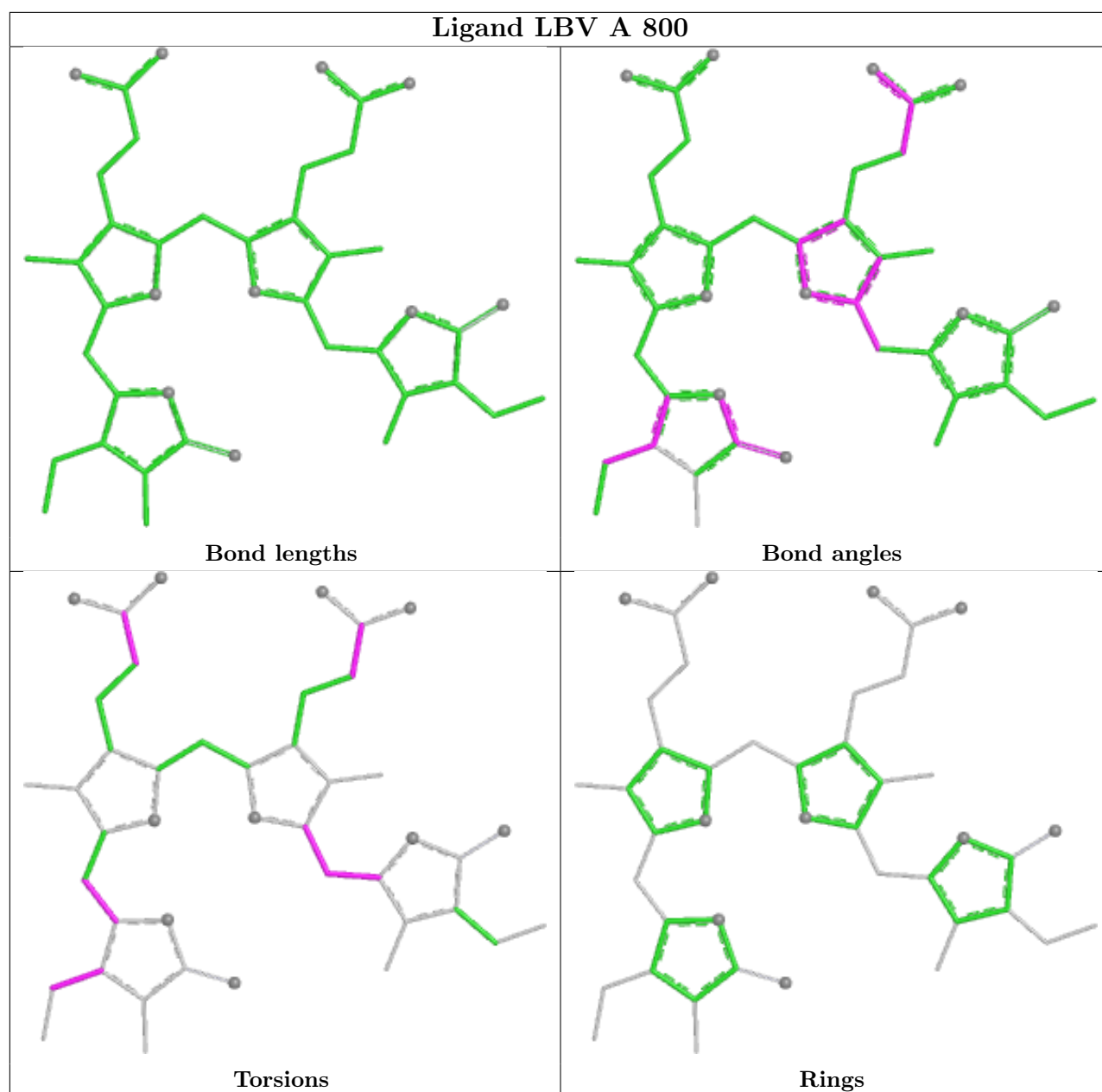
Ligand LBV C 800



Ligand LBV B 800







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	457/496 (92%)	1.15	89 (19%)	3 3	21, 53, 97, 134	3 (0%)
1	B	462/496 (93%)	0.92	60 (12%)	7 8	21, 46, 84, 113	2 (0%)
1	C	463/496 (93%)	1.18	96 (20%)	2 3	28, 53, 92, 128	1 (0%)
1	D	466/496 (93%)	0.88	48 (10%)	12 13	19, 45, 78, 118	1 (0%)
All	All	1848/1984 (93%)	1.03	293 (15%)	5 5	19, 48, 90, 134	7 (0%)

All (293) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	424	GLY	5.7
1	C	131	ALA	5.3
1	A	502	LEU	4.9
1	A	412	TRP	4.8
1	A	108	ALA	4.8
1	A	139	ALA	4.8
1	C	504	HIS	4.5
1	C	107	ALA	4.3
1	C	478	GLY	4.3
1	A	485	PRO	4.2
1	C	412	TRP	4.2
1	A	445	LEU	4.1
1	B	379	LEU	4.0
1	C	445	LEU	4.0
1	A	106	PRO	4.0
1	D	502	LEU	4.0
1	C	395	LEU	3.9
1	A	480	ALA	3.9
1	C	71	GLY	3.9
1	D	344	HIS	3.9
1	C	416	ALA	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	502	LEU	3.8
1	B	343	ALA	3.8
1	D	505	HIS	3.8
1	A	483	TRP	3.8
1	C	479[A]	TYR	3.8
1	A	418	LEU	3.8
1	A	109	GLY	3.7
1	C	105	TRP	3.7
1	B	77	LEU	3.7
1	D	60	THR	3.7
1	D	67	THR	3.7
1	B	419	ALA	3.7
1	A	395	LEU	3.6
1	A	386	PRO	3.6
1	B	235	ASP	3.6
1	C	138	HIS	3.6
1	C	345	SER	3.5
1	B	6	LEU	3.5
1	B	76	ALA	3.5
1	D	131	ALA	3.5
1	D	62	LEU	3.5
1	B	344	HIS	3.5
1	A	411	LEU	3.5
1	C	414	ALA	3.5
1	B	67	THR	3.4
1	A	1	MET	3.4
1	C	344	HIS	3.4
1	B	109	GLY	3.4
1	C	359	LEU	3.4
1	B	430	VAL	3.4
1	A	385	ALA	3.4
1	C	145	PHE	3.4
1	A	60	THR	3.4
1	A	77	LEU	3.3
1	A	331	LEU	3.3
1	C	346	LEU	3.3
1	C	411	LEU	3.3
1	B	47[A]	HIS	3.3
1	A	74	LEU	3.3
1	C	392	LEU	3.3
1	D	504	HIS	3.3
1	C	419	ALA	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	58	ALA	3.2
1	B	234	LEU	3.2
1	D	135	THR	3.2
1	B	107	ALA	3.2
1	C	96	ALA	3.2
1	D	105	TRP	3.2
1	A	413	PRO	3.2
1	D	332	ARG	3.1
1	A	400	GLY	3.1
1	C	331	LEU	3.1
1	C	431	GLY	3.1
1	A	66	PRO	3.1
1	D	106	PRO	3.1
1	A	391	LEU	3.1
1	C	485	PRO	3.1
1	B	25	GLU	3.1
1	B	105	TRP	3.1
1	C	422	ALA	3.1
1	C	69	LEU	3.1
1	C	408	LEU	3.1
1	A	105	TRP	3.0
1	C	139	ALA	3.0
1	C	480	ALA	3.0
1	A	138	HIS	3.0
1	A	484	HIS	3.0
1	C	417	ASP	3.0
1	A	399	PRO	3.0
1	D	343	ALA	3.0
1	D	419	ALA	3.0
1	A	346	LEU	2.9
1	D	24	CYS	2.9
1	A	142	ASN	2.9
1	A	406	ASP	2.9
1	B	501	ALA	2.9
1	D	107	ALA	2.9
1	B	22	GLU	2.9
1	C	393	ALA	2.9
1	D	59	ALA	2.9
1	D	430	VAL	2.9
1	C	67	THR	2.9
1	B	431	GLY	2.8
1	C	383	PRO	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	6	LEU	2.8
1	A	416	ALA	2.8
1	C	421	SER	2.8
1	B	137	PRO	2.8
1	D	66	PRO	2.8
1	D	234	LEU	2.8
1	D	498	LEU	2.8
1	D	214	ALA	2.8
1	B	412	TRP	2.8
1	A	71	GLY	2.8
1	D	475	GLY	2.8
1	A	94	PRO	2.8
1	A	307	TYR	2.8
1	A	297	LEU	2.7
1	C	1	MET	2.7
1	B	64	GLN	2.7
1	C	111	LEU	2.7
1	A	387	ALA	2.7
1	C	110	HIS	2.7
1	A	8	PHE	2.7
1	A	145	PHE	2.7
1	C	24	CYS	2.7
1	C	99	TYR	2.7
1	A	300	ASP	2.7
1	A	63	GLY	2.7
1	A	434	TRP	2.7
1	B	502	LEU	2.7
1	C	402	LEU	2.7
1	C	75	ALA	2.6
1	D	415	GLY	2.6
1	A	344	HIS	2.6
1	C	484	HIS	2.6
1	D	58	ALA	2.6
1	C	83	PRO	2.6
1	A	342	ALA	2.6
1	A	376	TRP	2.6
1	C	232	VAL	2.6
1	A	62	LEU	2.6
1	B	62	LEU	2.6
1	C	418	LEU	2.6
1	D	68	VAL	2.6
1	A	402	LEU	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	408	LEU	2.5
1	B	446	ARG	2.5
1	A	397	THR	2.5
1	B	106	PRO	2.5
1	C	68	VAL	2.5
1	A	295	TYR	2.5
1	B	130	GLU	2.5
1	A	140	LEU	2.5
1	B	89	LEU	2.5
1	C	62	LEU	2.5
1	C	77	LEU	2.5
1	D	97	LEU	2.5
1	A	482	PRO	2.5
1	B	136	GLY	2.5
1	C	400	GLY	2.5
1	A	327	PHE	2.5
1	D	477	SER	2.5
1	B	65	GLU	2.5
1	B	140	LEU	2.5
1	D	77	LEU	2.5
1	A	401	ALA	2.5
1	A	334	HIS	2.5
1	A	64[A]	GLN	2.5
1	A	61[A]	PHE	2.5
1	A	372	PHE	2.5
1	A	417	ASP	2.5
1	C	328	ARG	2.4
1	A	340	LEU	2.4
1	B	342	ALA	2.4
1	C	407	ALA	2.4
1	C	423	ALA	2.4
1	D	501	ALA	2.4
1	C	351	THR	2.4
1	C	483	TRP	2.4
1	C	154	ARG	2.4
1	A	355	PRO	2.4
1	C	347	SER	2.4
1	A	390	ALA	2.4
1	A	409	GLY	2.4
1	A	419	ALA	2.4
1	A	24	CYS	2.4
1	A	68	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	D	145	PHE	2.4
1	C	4	ASP	2.3
1	B	149	SER	2.3
1	C	8	PHE	2.3
1	B	85	LEU	2.3
1	A	481	GLU	2.3
1	C	406	ASP	2.3
1	D	412	TRP	2.3
1	A	59	ALA	2.3
1	A	88	ALA	2.3
1	B	54	MET	2.3
1	A	103	LEU	2.3
1	C	399	PRO	2.3
1	C	434	TRP	2.3
1	D	244	ALA	2.3
1	B	135	THR	2.3
1	A	3	ARG	2.3
1	A	345	SER	2.3
1	A	97	LEU	2.3
1	B	69	LEU	2.3
1	C	93	CYS	2.3
1	C	409	GLY	2.3
1	D	376	TRP	2.3
1	C	488	ILE	2.3
1	C	388	VAL	2.3
1	B	500	GLY	2.2
1	C	70	ARG	2.2
1	B	336	ALA	2.2
1	C	72	GLN	2.2
1	C	333	GLU	2.2
1	D	307	TYR	2.2
1	A	394	TRP	2.2
1	C	330	SER	2.2
1	A	259	MET	2.2
1	D	25	GLU	2.2
1	B	51	VAL	2.2
1	C	95	ASP	2.2
1	C	413	PRO	2.2
1	B	75	ALA	2.2
1	C	491	ALA	2.2
1	D	73	THR	2.2
1	C	376	TRP	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	359	LEU	2.2
1	C	301	LEU	2.2
1	B	18	GLU	2.2
1	C	46	GLY	2.2
1	C	374	GLY	2.2
1	D	47	HIS	2.2
1	D	71	GLY	2.2
1	A	96	ALA	2.2
1	D	339	ALA	2.2
1	A	85	LEU	2.1
1	A	111	LEU	2.1
1	B	259	MET	2.1
1	B	267	MET	2.1
1	B	376	TRP	2.1
1	C	97	LEU	2.1
1	D	301	LEU	2.1
1	C	300	ASP	2.1
1	B	145	PHE	2.1
1	C	106	PRO	2.1
1	C	503	GLU	2.1
1	A	72	GLN	2.1
1	A	332	ARG	2.1
1	B	340	LEU	2.1
1	B	82	TRP	2.1
1	D	268	GLY	2.1
1	A	129	THR	2.1
1	C	2	SER	2.1
1	B	74	LEU	2.1
1	B	432	GLU	2.1
1	C	85	LEU	2.1
1	C	357	LEU	2.1
1	D	69	LEU	2.1
1	B	142	ASN	2.1
1	A	47	HIS	2.1
1	B	138	HIS	2.1
1	B	7	PRO	2.1
1	B	298	PRO	2.1
1	D	413	PRO	2.1
1	B	61	PHE	2.1
1	B	71	GLY	2.1
1	B	146	ALA	2.1
1	B	325	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	363	MET	2.1
1	B	166	GLU	2.1
1	A	6	LEU	2.1
1	A	384	PRO	2.1
1	C	90	PRO	2.1
1	C	501	ALA	2.0
1	C	140	LEU	2.0
1	C	340	LEU	2.0
1	D	287	ILE	2.0
1	C	384	PRO	2.0
1	B	307	TYR	2.0
1	A	328	ARG	2.0
1	D	70	ARG	2.0
1	A	250	GLY	2.0
1	D	92	GLY	2.0
1	C	261	MET	2.0
1	C	397	THR	2.0
1	A	422	ALA	2.0
1	C	91	PRO	2.0
1	C	482	PRO	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
2	LBV	B	800	43/43	0.91	0.10	33,38,59,70	0
2	LBV	C	800	43/43	0.92	0.11	32,38,54,78	0
2	LBV	D	800	43/43	0.92	0.09	30,37,49,62	0

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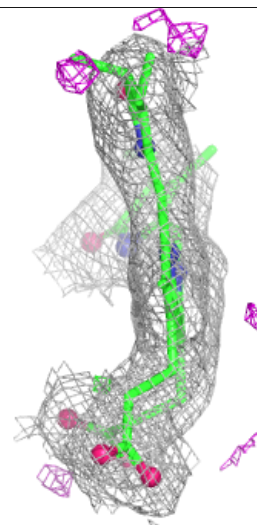
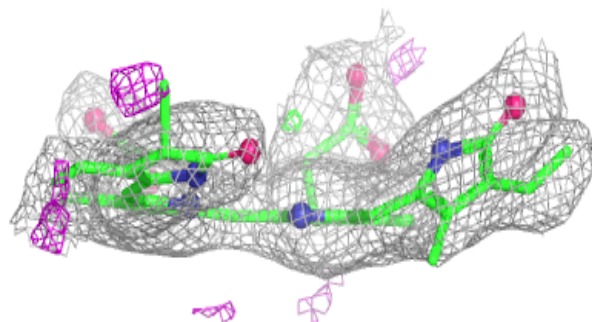
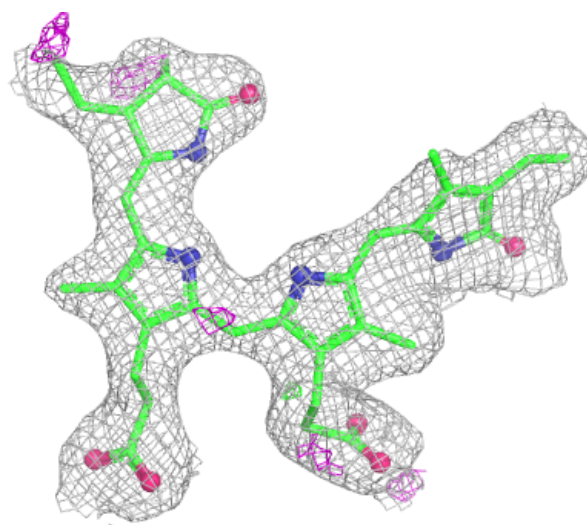
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	LBV	A	800	43/43	0.93	0.10	32,38,52,65	0

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.

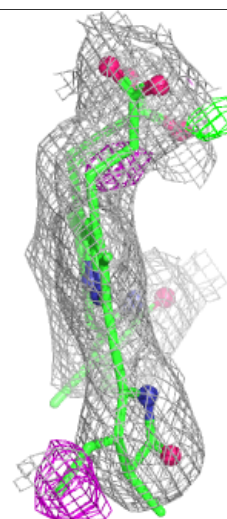
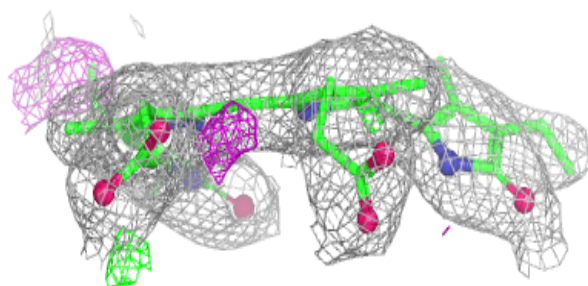
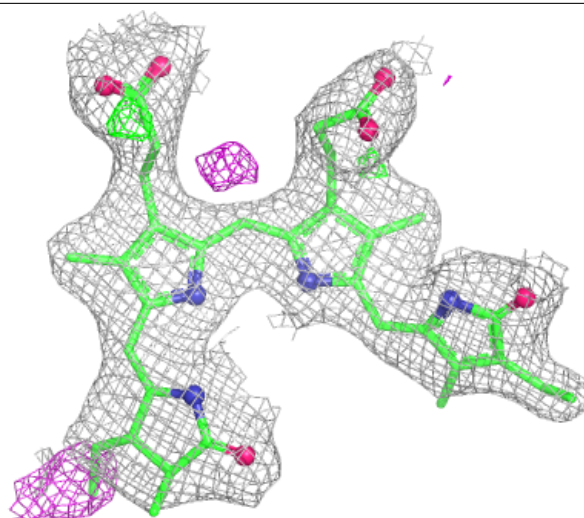
Electron density around LBV B 800:

2mF_o-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



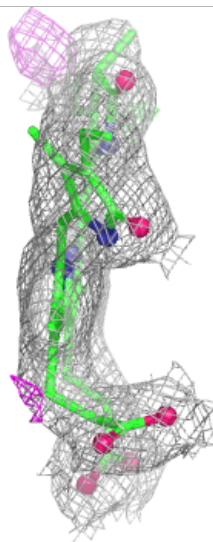
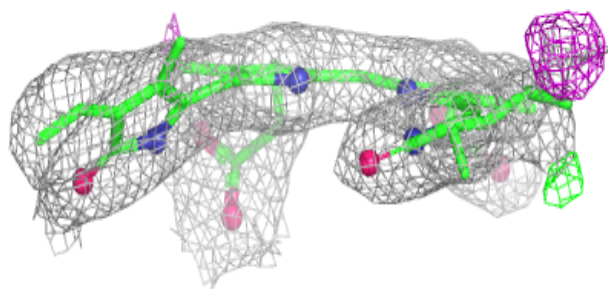
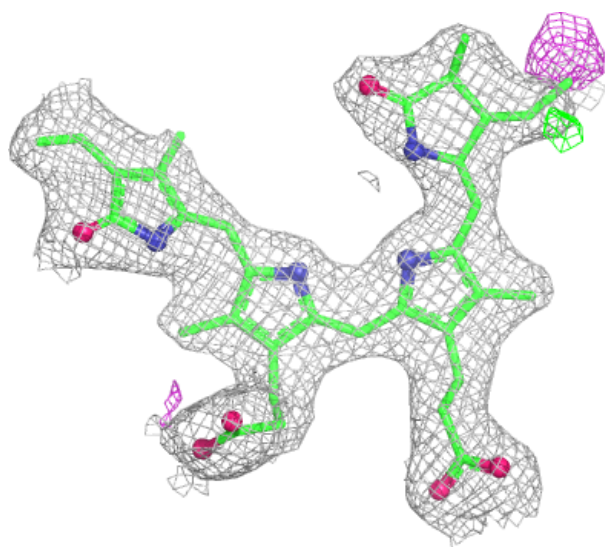
Electron density around LBV C 800:

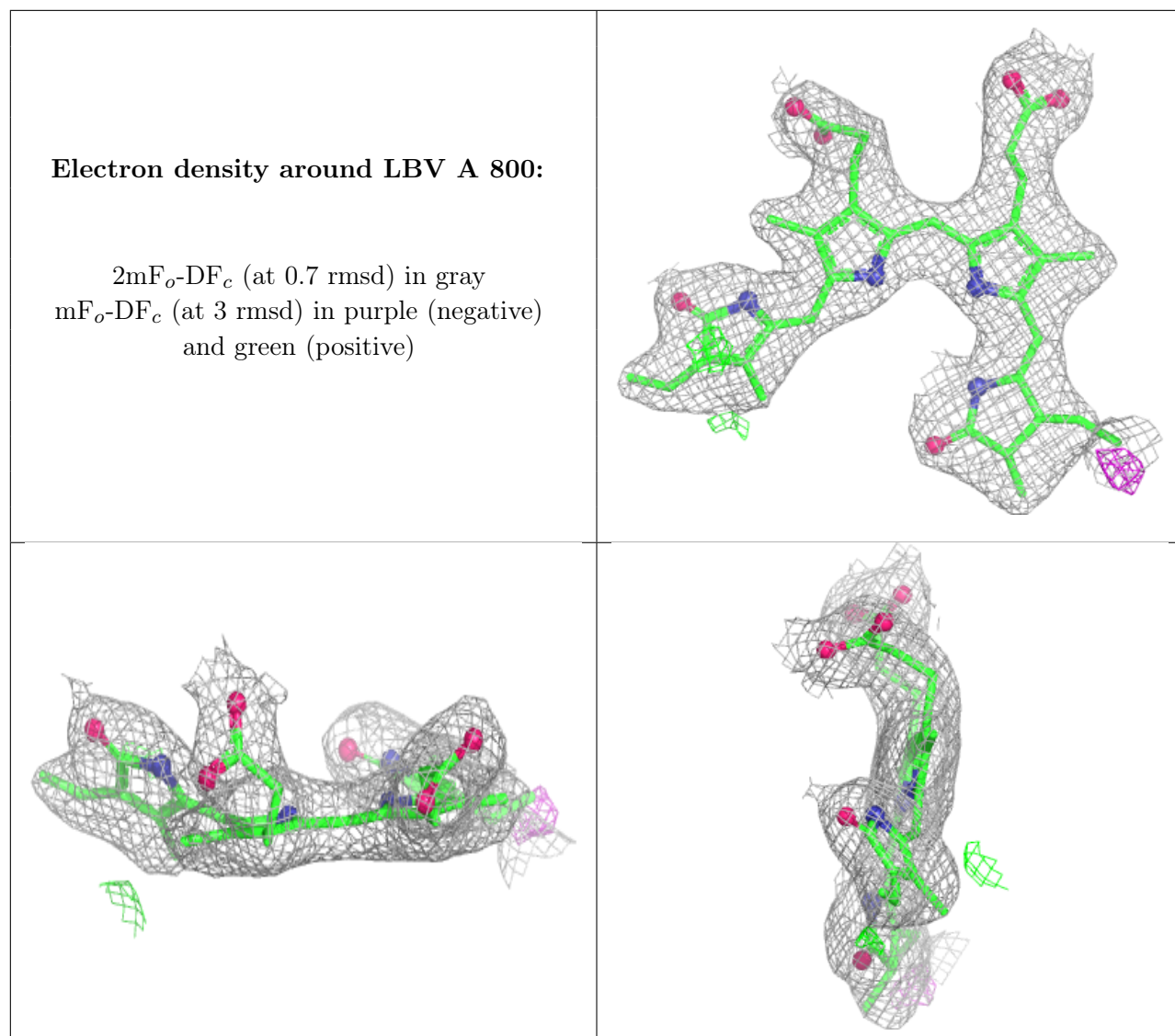
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around LBV D 800:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.