



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 22, 2016 – 06:53 PM EDT

PDB ID : 5C5K
Title : Structure of the Pfr form of a canonical phytochrome
Authors : Burgie, E.S.; Vierstra, R.D.
Deposited on : 2015-06-20
Resolution : 3.31 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
<http://wwpdb.org/validation/2016/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
Mogul : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027107
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0122
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : rb-20027107

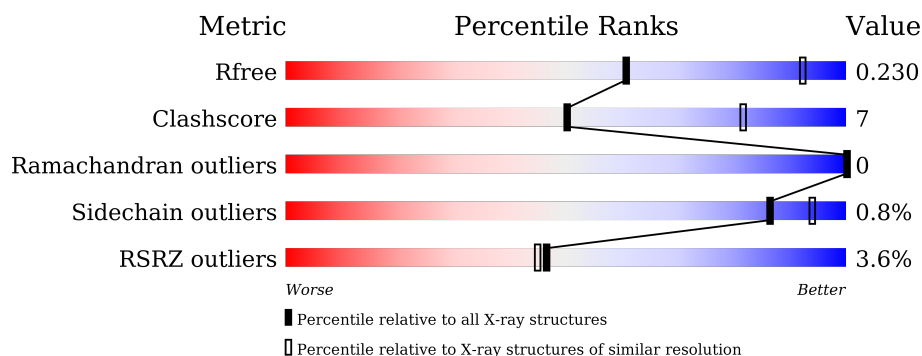
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.31 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1198 (3.40-3.24)
Clashscore	102246	1280 (3.40-3.24)
Ramachandran outliers	100387	1260 (3.40-3.24)
Sidechain outliers	100360	1259 (3.40-3.24)
RSRZ outliers	91569	1203 (3.40-3.24)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	524	<div> <div>2%</div> <div> <div></div> <div>76%</div> <div>17%</div> <div>6%</div> </div> </div>
1	B	524	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>14%</div> <div>7%</div> </div> </div>
1	C	524	<div> <div>3%</div> <div> <div></div> <div>75%</div> <div>15%</div> <div>9%</div> </div> </div>
1	D	524	<div> <div>5%</div> <div> <div></div> <div>76%</div> <div>14%</div> <div>10%</div> </div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	EDO	A	602	-	-	-	X
5	BGC	C	602	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 14949 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	490	Total	C	N	O	S	0	0	0
			3738	2379	660	688	11			
1	B	486	Total	C	N	O	S	0	0	0
			3716	2366	656	683	11			
1	C	476	Total	C	N	O	S	0	1	0
			3649	2324	644	670	11			
1	D	474	Total	C	N	O	S	0	1	0
			3636	2318	641	666	11			

There are 92 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	MET	-	initiating methionine	UNP Q9RZA4
A	-13	ALA	-	expression tag	UNP Q9RZA4
A	-12	LEU	-	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	469	TRP	PHE	engineered mutation	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

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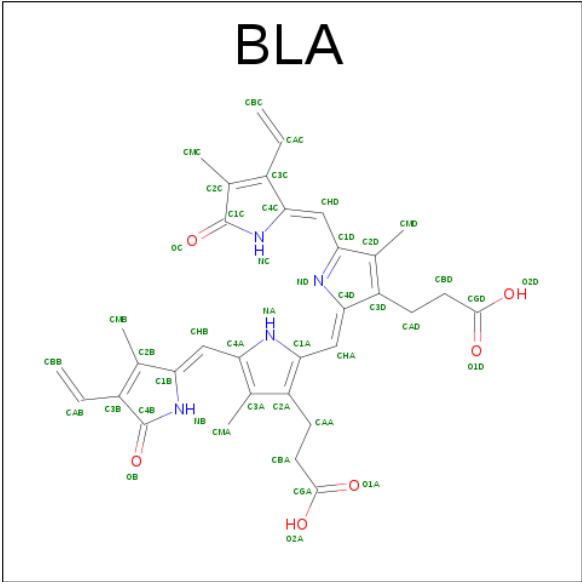
Chain	Residue	Modelled	Actual	Comment	Reference
A	508	HIS	-	expression tag	UNP Q9RZA4
A	509	HIS	-	expression tag	UNP Q9RZA4
B	-14	MET	-	initiating methionine	UNP Q9RZA4
B	-13	ALA	-	expression tag	UNP Q9RZA4
B	-12	LEU	-	expression tag	UNP Q9RZA4
B	-11	SER	-	expression tag	UNP Q9RZA4
B	-10	MET	-	expression tag	UNP Q9RZA4
B	-9	THR	-	expression tag	UNP Q9RZA4
B	-8	GLY	-	expression tag	UNP Q9RZA4
B	-7	GLY	-	expression tag	UNP Q9RZA4
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	469	TRP	PHE	engineered mutation	UNP Q9RZA4
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	-14	MET	-	initiating methionine	UNP Q9RZA4
C	-13	ALA	-	expression tag	UNP Q9RZA4
C	-12	LEU	-	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	469	TRP	PHE	engineered mutation	UNP Q9RZA4
C	503	GLU	-	expression tag	UNP Q9RZA4

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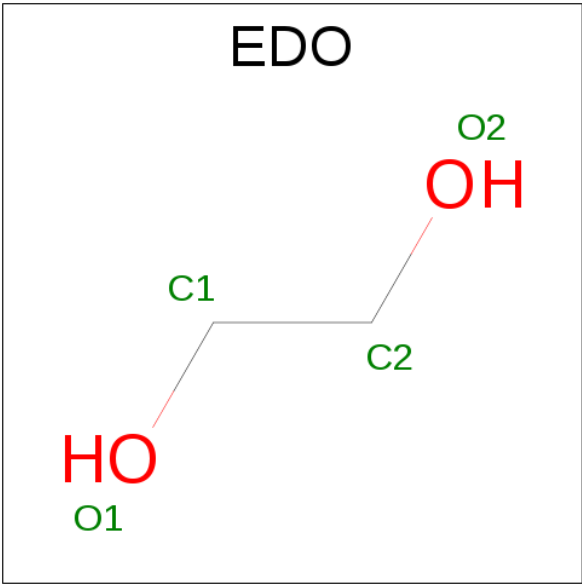
Chain	Residue	Modelled	Actual	Comment	Reference
C	504	HIS	-	expression tag	UNP Q9RZA4
C	505	HIS	-	expression tag	UNP Q9RZA4
C	506	HIS	-	expression tag	UNP Q9RZA4
C	507	HIS	-	expression tag	UNP Q9RZA4
C	508	HIS	-	expression tag	UNP Q9RZA4
C	509	HIS	-	expression tag	UNP Q9RZA4
D	-14	MET	-	initiating methionine	UNP Q9RZA4
D	-13	ALA	-	expression tag	UNP Q9RZA4
D	-12	LEU	-	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	469	TRP	PHE	engineered mutation	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 BILIVERDINE IX ALPHA (three-letter code: BLA) (formula: C₃₃H₃₄N₄O₆).



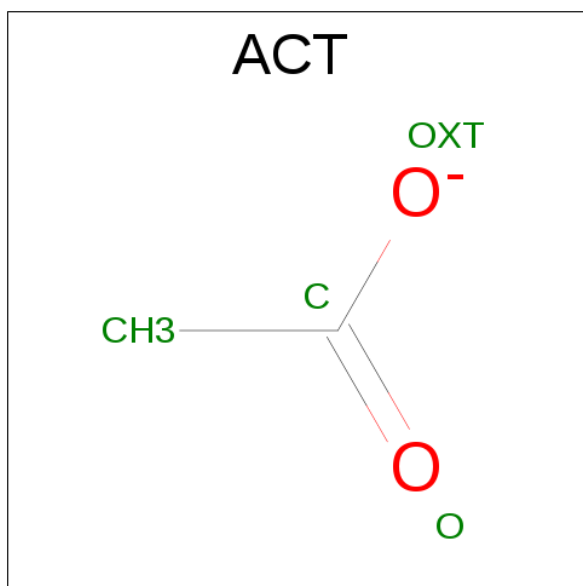
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			43	33	4	6		
2	B	1	Total	C	N	O	0	0
			43	33	4	6		
2	C	1	Total	C	N	O	0	0
			43	33	4	6		
2	D	1	Total	C	N	O	0	0
			43	33	4	6		

- Molecule 3 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



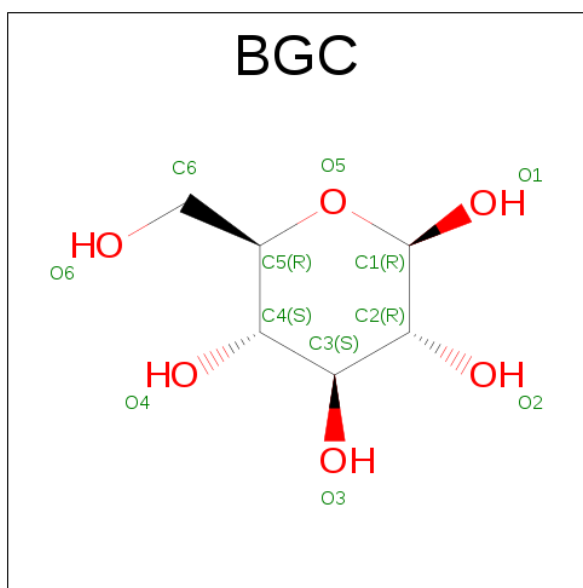
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 5 is BETA-D-GLUCOSE (three-letter code: BGC) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	C	O	0	0
			12	6	6		

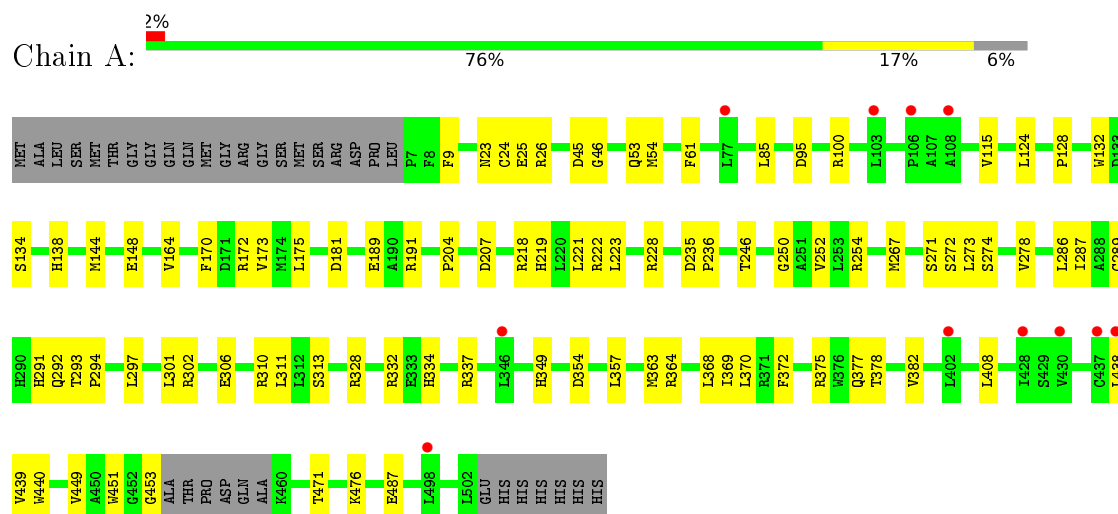
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	10	Total	O	0	0
			10	10		
6	B	6	Total	O	0	0
			6	6		
6	C	1	Total	O	0	0
			1	1		
6	D	1	Total	O	0	0
			1	1		

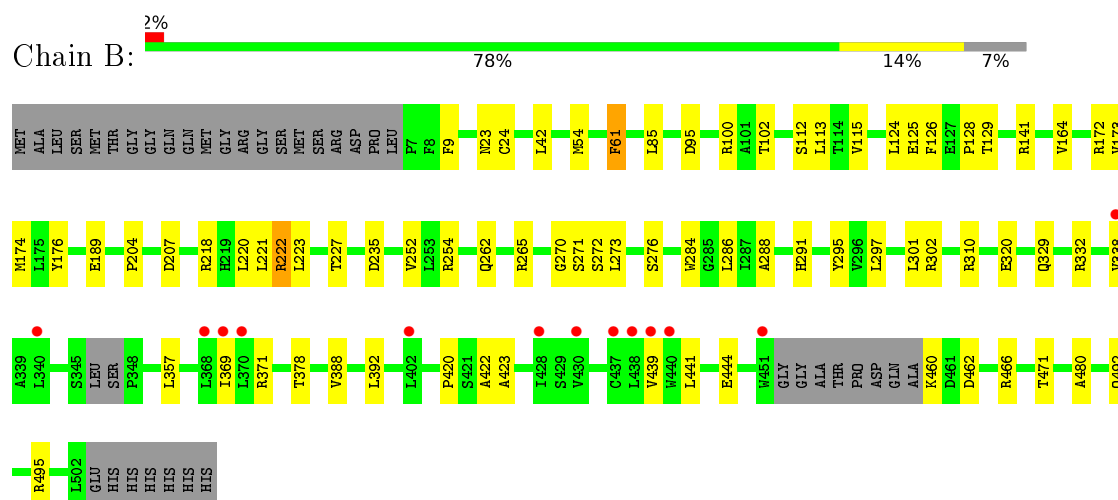
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

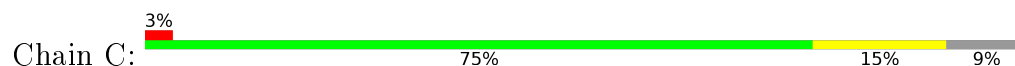
• Molecule 1: Bacteriophytochrome

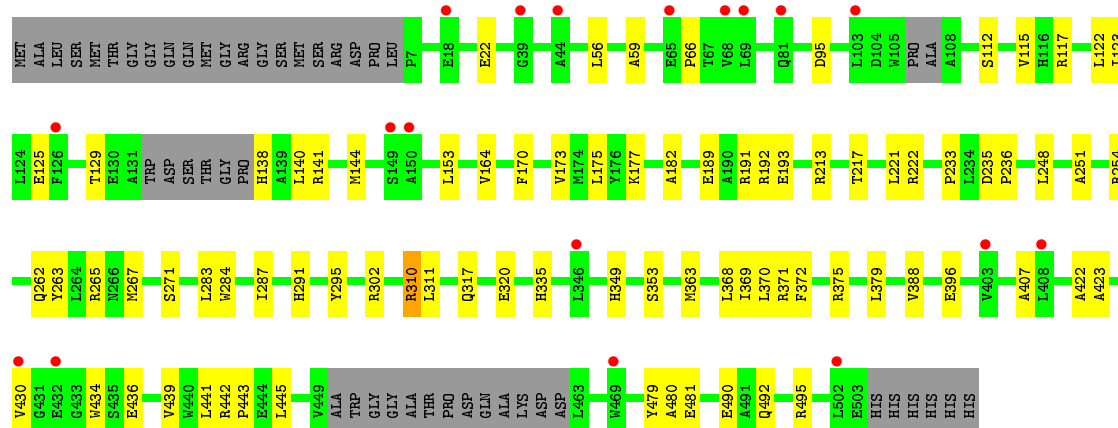


• Molecule 1: Bacteriophytochrome

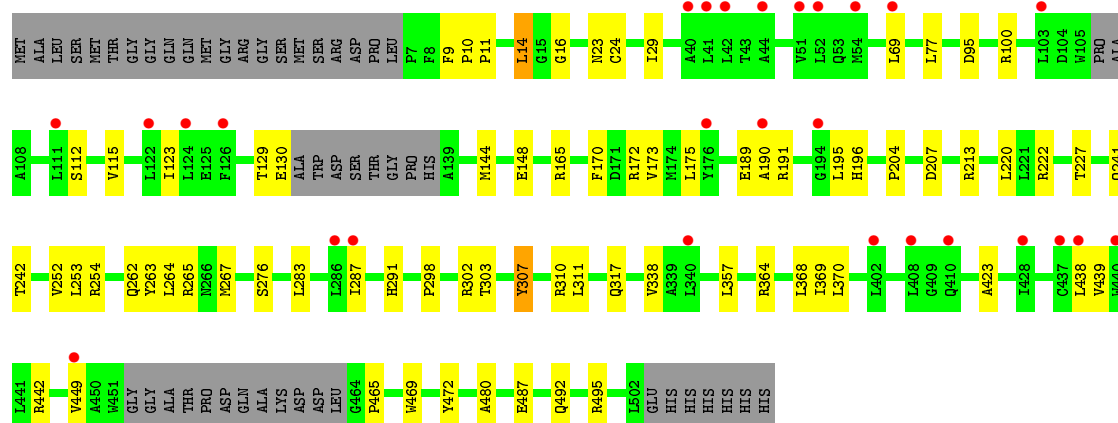
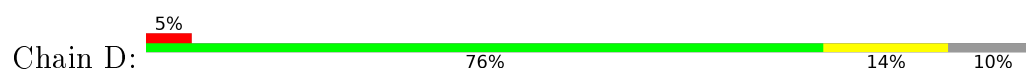


• Molecule 1: Bacteriophytochrome





● Molecule 1: Bacteriophytochrome



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	89.40Å 192.75Å 225.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.80 – 3.31 48.80 – 3.31	Depositor EDS
% Data completeness (in resolution range)	87.9 (48.80-3.31) 87.9 (48.80-3.31)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.11 (at 3.33Å)	Xtriage
Refinement program	PHENIX (1.10.1 _2155: ???)	Depositor
R, R_{free}	0.196 , 0.232 0.194 , 0.230	Depositor DCC
R_{free} test set	1882 reflections (3.64%)	DCC
Wilson B-factor (Å ²)	123.7	Xtriage
Anisotropy	0.012	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 89.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 51724 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14949	wwPDB-VP
Average B, all atoms (Å ²)	153.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.32% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: BGC, BLA, EDO, ACT

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	0.25	0/3839	0.42	0/5257
1	B	0.25	0/3816	0.44	0/5224
1	C	0.24	0/3743	0.42	0/5120
1	D	0.25	0/3731	0.43	0/5105
All	All	0.25	0/15129	0.43	0/20706

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	3738	0	3695	58	1
1	B	3716	0	3673	51	1
1	C	3649	0	3616	54	0
1	D	3636	0	3602	51	0
2	A	43	0	31	4	0
2	B	43	0	31	6	0
2	C	43	0	31	2	0
2	D	43	0	31	4	0
3	A	4	0	6	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	4	0	3	0	0
5	C	12	0	12	0	0
6	A	10	0	0	1	0
6	B	6	0	0	2	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
All	All	14949	0	14731	208	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (208) 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:95:ASP:OD1	1:B:218:ARG:NH1	1.92	1.03
1:C:112:SER:HB2	1:C:129:THR:HB	1.67	0.75
1:B:273:LEU:HD13	1:B:297:LEU:HD12	1.71	0.72
1:A:207:ASP:HB3	2:A:601:BLA:HHB	1.73	0.70
1:C:144:MET:HG3	1:C:311:LEU:HD21	1.73	0.68
1:C:353:SER:HA	1:C:379:LEU:HD11	1.74	0.68
1:C:222:ARG:HE	1:C:254:ARG:NH1	1.92	0.67
1:D:112:SER:HB2	1:D:129:THR:HB	1.75	0.67
1:B:271:SER:HB2	1:B:295:TYR:H	1.59	0.67
1:B:460:LYS:HE2	1:B:462:ASP:HB3	1.78	0.66
1:D:29:ILE:HD12	2:D:601:BLA:HMD1	1.78	0.66
1:A:144:MET:HG3	1:A:311:LEU:HD11	1.78	0.65
1:B:371:ARG:HH11	1:B:392:LEU:HD13	1.60	0.65
1:D:220:LEU:HD12	1:D:310:ARG:NH1	2.11	0.65
1:A:9:PHE:HE2	1:A:23:ASN:HB2	1.62	0.64
1:A:274:SER:HB2	1:A:286:LEU:HD21	1.79	0.64
1:A:271:SER:HB2	1:A:294:PRO:HA	1.80	0.64
1:A:100:ARG:O	1:B:141:ARG:NH1	2.30	0.64
1:D:252:VAL:HG11	1:D:302:ARG:HH21	1.61	0.63
1:B:220:LEU:HD12	1:B:310:ARG:NH1	2.15	0.62
1:A:297:LEU:HD13	1:A:302:ARG:HB2	1.80	0.62
1:B:423:ALA:HB3	1:B:480:ALA:HB2	1.80	0.62
1:A:95:ASP:OD1	1:A:218:ARG:NH2	2.29	0.62
1:B:9:PHE:HE2	1:B:23:ASN:HB2	1.65	0.61
1:B:284:TRP:NE1	1:B:320:GLU:OE2	2.35	0.60
1:C:317:GLN:OE1	1:D:317:GLN:NE2	2.32	0.60
1:A:85:LEU:HD11	1:A:115:VAL:HG11	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:222:ARG:NH2	1:D:254:ARG:O	2.32	0.59
1:A:334:HIS:CD2	1:A:337:ARG:HH22	2.21	0.59
1:C:217:THR:HG22	1:C:283:LEU:HB2	1.85	0.58
1:A:292:GLN:HG3	1:A:293:THR:HG23	1.85	0.58
1:C:310:ARG:NH1	1:D:144:MET:SD	2.77	0.58
1:D:369:ILE:HB	1:D:439:VAL:HB	1.86	0.58
1:A:369:ILE:HB	1:A:439:VAL:HB	1.85	0.58
1:B:272:SER:OG	2:B:601:BLA:O1A	2.19	0.57
1:B:276:SER:HA	1:B:286:LEU:HD12	1.85	0.57
1:D:252:VAL:HG23	1:D:253:LEU:HD12	1.84	0.57
1:B:125:GLU:OE2	1:B:252:VAL:N	2.37	0.57
1:B:297:LEU:HD22	1:B:301:LEU:HD22	1.87	0.57
1:A:310:ARG:NH1	3:A:602:EDO:H12	2.19	0.57
1:A:170:PHE:CD1	1:A:291:HIS:HB2	2.40	0.57
1:C:222:ARG:HE	1:C:254:ARG:HH11	1.51	0.56
1:B:112:SER:HB2	1:B:129:THR:HB	1.86	0.56
1:C:153:LEU:HD11	1:C:320:GLU:HG2	1.87	0.56
1:A:354:ASP:HB3	1:A:357:LEU:HD13	1.88	0.56
1:C:423:ALA:HB3	1:C:480:ALA:HB2	1.87	0.55
1:C:140:LEU:HD21	1:D:307[A]:TYR:HE2	1.71	0.55
1:A:25:GLU:HG3	1:A:26:ARG:HG3	1.88	0.55
1:D:172:ARG:NH2	1:D:196:HIS:O	2.39	0.55
1:C:271:SER:HB2	1:C:295:TYR:H	1.71	0.55
1:B:338:VAL:HA	1:B:357:LEU:HD11	1.88	0.55
1:D:175:LEU:HD13	1:D:287:ILE:HG12	1.89	0.55
1:C:125:GLU:OE2	1:C:251:ALA:HA	2.07	0.54
1:A:451:TRP:HD1	1:A:453:GLY:H	1.55	0.54
1:A:278:VAL:O	1:A:313:SER:OG	2.13	0.54
1:A:115:VAL:HG12	1:A:124:LEU:HA	1.89	0.54
1:C:22:GLU:N	1:C:22:GLU:OE2	2.41	0.54
1:A:221:LEU:HD13	1:A:306:GLU:HG2	1.91	0.53
1:D:492:GLN:HG3	1:D:495:ARG:HH21	1.73	0.53
1:B:420:PRO:HA	1:B:444:GLU:HB2	1.91	0.53
1:A:164:VAL:HB	1:A:173:VAL:HG11	1.92	0.52
1:A:189:GLU:OE2	1:A:191:ARG:HD3	2.09	0.52
1:D:227:THR:HG21	1:D:264:LEU:HB2	1.91	0.52
1:A:375:ARG:NH2	1:A:377:GLN:OE1	2.42	0.52
1:B:174:MET:HG2	1:B:189:GLU:HG3	1.92	0.52
1:A:53:GLN:OE1	1:A:246:THR:OG1	2.23	0.52
2:B:601:BLA:NC	6:B:701:HOH:O	2.33	0.52
1:A:252:VAL:HG11	1:A:302:ARG:HH12	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:207:ASP:HB3	2:B:601:BLA:HHB	1.92	0.51
1:B:95:ASP:OD1	1:B:95:ASP:N	2.44	0.51
1:D:267:MET:SD	2:D:601:BLA:HAB	2.50	0.51
1:C:221:LEU:HD21	1:C:302:ARG:HD2	1.91	0.51
1:C:492:GLN:HA	1:C:495:ARG:HG2	1.92	0.51
1:D:241:GLN:HG3	1:D:242:THR:HG23	1.91	0.51
2:B:601:BLA:HMB3	2:B:601:BLA:CMA	2.40	0.51
1:B:221:LEU:HD21	1:B:302:ARG:HD2	1.93	0.50
1:B:222:ARG:HG3	1:B:254:ARG:HH12	1.77	0.50
1:D:172:ARG:HD2	1:D:195:LEU:HD13	1.93	0.50
1:D:165:ARG:HB2	1:D:173:VAL:HG11	1.93	0.50
1:D:222:ARG:HE	1:D:254:ARG:NH1	2.08	0.50
1:C:189:GLU:OE2	1:C:191:ARG:HD3	2.11	0.50
1:C:170:PHE:CD1	1:C:291:HIS:HB2	2.47	0.50
1:C:164:VAL:HB	1:C:173:VAL:HG11	1.94	0.49
1:B:85:LEU:HD11	1:B:115:VAL:HG21	1.95	0.49
1:B:460:LYS:N	1:B:466:ARG:HH11	2.11	0.49
1:C:141:ARG:NH1	1:D:100:ARG:HD3	2.28	0.49
1:A:222:ARG:HB3	1:A:274:SER:OG	2.13	0.49
1:B:254:ARG:NH1	6:B:702:HOH:O	2.45	0.48
2:B:601:BLA:HMB3	2:B:601:BLA:HMA1	1.95	0.48
1:A:223:LEU:HD13	1:A:273:LEU:HD13	1.95	0.48
1:B:262:GLN:OE1	1:B:265:ARG:NH1	2.39	0.48
1:D:10:PRO:HG2	1:D:16:GLY:HA2	1.96	0.48
1:C:445:LEU:HD13	1:C:481:GLU:HG2	1.96	0.48
1:D:267:MET:HG2	1:D:465:PRO:HG3	1.95	0.48
1:A:370:LEU:HD23	1:A:438:LEU:HD22	1.96	0.48
1:C:117:ARG:HH11	1:C:122:LEU:HD21	1.78	0.48
1:B:460:LYS:HG2	1:B:462:ASP:H	1.79	0.47
1:B:338:VAL:HG22	1:B:357:LEU:HG	1.95	0.47
1:B:371:ARG:NH1	1:B:392:LEU:HD13	2.28	0.47
1:B:113:LEU:HD11	1:B:124:LEU:HD23	1.97	0.47
1:D:364:ARG:NE	1:D:487:GLU:OE2	2.46	0.47
1:A:222:ARG:NE	1:A:254:ARG:NH1	2.62	0.47
2:A:601:BLA:O2A	2:A:601:BLA:HMA2	2.15	0.47
1:D:207:ASP:HB3	2:D:601:BLA:HHB	1.96	0.47
1:B:329:GLN:HA	1:B:332:ARG:HE	1.79	0.47
1:B:369:ILE:HD13	1:B:388:VAL:HG13	1.97	0.46
1:C:175:LEU:HD12	1:C:287:ILE:HG12	1.97	0.46
1:C:369:ILE:HB	1:C:439:VAL:HB	1.96	0.46
1:C:59:ALA:HB2	1:C:66:PRO:HB3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:95:ASP:OD1	1:D:95:ASP:N	2.48	0.46
1:D:204:PRO:HD3	1:D:472:TYR:HB2	1.97	0.46
1:A:172:ARG:NH1	6:A:701:HOH:O	2.38	0.46
1:C:422:ALA:HB1	1:C:441:LEU:HD22	1.98	0.46
1:C:371:ARG:O	1:C:436:GLU:HA	2.15	0.46
1:C:372:PHE:HA	1:C:436:GLU:HG2	1.97	0.46
1:A:138:HIS:NE2	1:B:102:THR:OG1	2.46	0.46
1:C:95:ASP:OD1	1:C:95:ASP:N	2.48	0.46
1:A:252:VAL:HG11	1:A:302:ARG:HH22	1.81	0.46
1:C:349:HIS:HE1	1:C:375:ARG:HG3	1.80	0.46
1:A:132:TRP:CZ2	1:A:134:SER:HB3	2.51	0.45
1:D:170:PHE:CD1	1:D:291:HIS:HB2	2.50	0.45
1:A:378:THR:HG21	1:A:382:VAL:HG11	1.97	0.45
1:C:335:HIS:HB2	1:C:490:GLU:OE2	2.16	0.45
1:A:181:ASP:OD2	1:A:476:LYS:NZ	2.42	0.45
1:D:213:ARG:HG3	1:D:283:LEU:HD23	1.97	0.45
1:B:54:MET:HE3	1:B:54:MET:HB2	1.67	0.45
1:C:369:ILE:HD13	1:C:388:VAL:HG13	1.98	0.45
1:C:396:GLU:HG2	1:C:434:TRP:HB2	1.98	0.45
1:D:222:ARG:NE	1:D:254:ARG:NH1	2.64	0.45
1:D:130:GLU:OE2	1:D:298:PRO:HB3	2.16	0.45
1:C:222:ARG:NE	1:C:254:ARG:HH11	2.13	0.45
1:C:368:LEU:HD21	1:C:370:LEU:HD21	1.99	0.45
1:A:61:PHE:HZ	1:A:128:PRO:HD3	1.82	0.45
1:A:95:ASP:OD1	1:A:95:ASP:N	2.48	0.45
1:D:449:VAL:HG22	1:D:469:TRP:NE1	2.32	0.45
1:D:173:VAL:HG13	1:D:190:ALA:HB3	1.98	0.45
1:A:267:MET:SD	2:A:601:BLA:HAB	2.57	0.45
1:C:138:HIS:ND1	1:C:138:HIS:O	2.50	0.44
1:D:338:VAL:HA	1:D:357:LEU:HD22	1.99	0.44
1:D:69:LEU:HD11	1:D:77:LEU:HD21	1.99	0.44
1:B:369:ILE:HB	1:B:439:VAL:HB	1.99	0.44
1:B:24:CYS:HB2	2:B:601:BLA:HAC	1.76	0.44
1:A:54:MET:O	1:A:235:ASP:N	2.33	0.44
1:B:172:ARG:HD3	1:B:174:MET:SD	2.57	0.44
1:D:423:ALA:HB3	1:D:480:ALA:HB2	1.99	0.44
1:A:364:ARG:HD3	1:A:487:GLU:OE2	2.18	0.44
1:C:56:LEU:HD12	1:C:233:PRO:HD2	2.00	0.44
1:D:115:VAL:HA	1:D:123:ILE:O	2.18	0.44
1:D:423:ALA:N	1:D:442:ARG:O	2.49	0.44
1:C:140:LEU:HD21	1:D:307[A]:TYR:CE2	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:368:LEU:HD11	1:D:438:LEU:HD11	2.00	0.43
1:C:310:ARG:HG2	1:D:148:GLU:OE2	2.18	0.43
1:D:189:GLU:OE2	1:D:191:ARG:HD3	2.19	0.43
1:A:204:PRO:HG3	1:A:471:THR:HG23	1.99	0.43
1:D:144:MET:HG3	1:D:311:LEU:HD21	2.00	0.43
1:A:222:ARG:HE	1:A:254:ARG:NH1	2.17	0.43
1:C:267:MET:HE1	2:C:601:BLA:HAB	2.01	0.43
1:B:54:MET:O	1:B:235:ASP:N	2.37	0.42
1:C:213:ARG:O	1:C:217:THR:HG23	2.19	0.42
1:A:132:TRP:CZ2	1:A:301:LEU:HD13	2.54	0.42
1:A:223:LEU:HD12	1:A:272:SER:O	2.19	0.42
1:D:220:LEU:HD12	1:D:310:ARG:HH11	1.82	0.42
1:D:9:PHE:HE2	1:D:23:ASN:HB2	1.85	0.42
1:A:363:MET:HG3	1:A:440:TRP:CD2	2.55	0.42
1:B:270:GLY:N	1:B:291:HIS:O	2.44	0.42
1:A:349:HIS:NE2	1:A:372:PHE:HB2	2.34	0.42
1:A:219:HIS:NE2	1:A:250:GLY:HA2	2.35	0.42
1:D:220:LEU:HA	1:D:276:SER:HB2	2.02	0.42
1:B:42:LEU:HD13	1:B:126:PHE:CE1	2.55	0.42
1:D:11:PRO:HD2	1:D:14:LEU:HD21	2.02	0.42
1:A:45:ASP:OD1	1:A:46:GLY:N	2.53	0.41
1:B:61:PHE:HE2	1:B:128:PRO:HD3	1.85	0.41
1:B:204:PRO:HG3	1:B:471:THR:HG23	2.02	0.41
1:D:24:CYS:HB2	2:D:601:BLA:HAC	1.89	0.41
1:A:148:GLU:OE2	1:B:310:ARG:NE	2.50	0.41
1:B:176:TYR:CE2	1:B:288:ALA:HB2	2.54	0.41
1:C:363:MET:O	1:C:442:ARG:NH2	2.38	0.41
1:D:370:LEU:HD23	1:D:438:LEU:CD1	2.50	0.41
1:A:175:LEU:HD13	1:A:287:ILE:HG12	2.01	0.41
1:C:177:LYS:HD3	1:C:284:TRP:CZ2	2.55	0.41
1:C:235:ASP:HA	1:C:236:PRO:HA	1.93	0.41
1:A:170:PHE:CG	1:A:289:CYS:HB3	2.56	0.41
1:A:328:ARG:O	1:A:332:ARG:HG2	2.20	0.41
1:C:222:ARG:NH2	1:C:254:ARG:O	2.52	0.41
1:C:267:MET:HE1	2:C:601:BLA:HMB1	2.01	0.41
1:A:24:CYS:HB2	2:A:601:BLA:HAC	1.59	0.41
1:B:492:GLN:HG3	1:B:495:ARG:HH21	1.85	0.41
1:A:53:GLN:HA	1:A:236:PRO:HD2	2.02	0.41
1:B:222:ARG:HD3	1:B:223:LEU:N	2.36	0.41
1:C:248:LEU:O	1:C:254:ARG:HD3	2.21	0.41
1:C:407:ALA:HB2	1:C:479:TYR:HA	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:164:VAL:HB	1:B:173:VAL:HG11	2.02	0.41
1:B:329:GLN:HG3	1:B:332:ARG:HH21	1.86	0.41
1:C:115:VAL:HA	1:C:123:ILE:O	2.20	0.41
1:C:262:GLN:HG2	1:C:265:ARG:NH1	2.36	0.41
1:B:100:ARG:HA	1:B:113:LEU:O	2.21	0.41
1:B:422:ALA:HB2	1:B:441:LEU:HD23	2.02	0.41
1:C:442:ARG:HA	1:C:443:PRO:HD3	1.95	0.41
1:C:141:ARG:HD3	1:D:100:ARG:HE	1.83	0.41
1:D:262:GLN:HG2	1:D:265:ARG:NH1	2.35	0.41
1:A:297:LEU:HD12	1:A:297:LEU:O	2.21	0.40
1:A:368:LEU:HD21	1:A:370:LEU:HD21	2.04	0.40
1:C:192:ARG:HD2	1:C:193:GLU:O	2.21	0.40
1:D:303:THR:HG22	1:D:307[B]:TYR:CZ	2.55	0.40
1:A:363:MET:HG3	1:A:440:TRP:CE2	2.56	0.40
1:C:182:ALA:O	1:C:213:ARG:NH2	2.54	0.40
1:A:310:ARG:HH12	3:A:602:EDO:H12	1.83	0.40
1:B:460:LYS:HB3	1:B:460:LYS:NZ	2.37	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:228:ARG:NH1	1:B:128:PRO:O[4_445]	2.09	0.11

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	486/524 (93%)	478 (98%)	8 (2%)	0	100	100
1	B	480/524 (92%)	473 (98%)	7 (2%)	0	100	100
1	C	469/524 (90%)	463 (99%)	6 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	467/524 (89%)	460 (98%)	7 (2%)	0	100	100
All	All	1902/2096 (91%)	1874 (98%)	28 (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	384/411 (93%)	382 (100%)	2 (0%)	92	96
1	B	382/411 (93%)	378 (99%)	4 (1%)	82	92
1	C	376/411 (92%)	373 (99%)	3 (1%)	86	94
1	D	374/411 (91%)	370 (99%)	4 (1%)	80	91
All	All	1516/1644 (92%)	1503 (99%)	13 (1%)	86	93

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	408	LEU
1	A	449	VAL
1	B	61	PHE
1	B	222	ARG
1	B	227	THR
1	B	378	THR
1	C	263	TYR
1	C	310	ARG
1	C	430	VAL
1	D	14	LEU
1	D	263	TYR
1	D	307[A]	TYR
1	D	307[B]	TYR

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

Mol	Chain	Res	Type
1	B	260	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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

7 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	BLA	A	601	1	36,46,46	3.72	11 (30%)	45,67,67	1.36	5 (11%)
3	EDO	A	602	-	3,3,3	0.45	0	2,2,2	0.40	0
4	ACT	A	603	-	0,3,3	0.00	-	0,3,3	0.00	-
2	BLA	B	601	1	36,46,46	3.87	10 (27%)	45,67,67	1.67	7 (15%)
2	BLA	C	601	1	36,46,46	3.69	11 (30%)	45,67,67	1.28	5 (11%)
5	BGC	C	602	-	12,12,12	0.54	0	17,17,17	0.67	0
2	BLA	D	601	1	36,46,46	3.72	11 (30%)	45,67,67	1.34	6 (13%)

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	BLA	A	601	1	-	1/22/74/74	0/4/4/4
3	EDO	A	602	-	-	0/1/1/1	0/0/0/0
4	ACT	A	603	-	-	0/0/0/0	0/0/0/0
2	BLA	B	601	1	-	1/22/74/74	0/4/4/4
2	BLA	C	601	1	-	1/22/74/74	0/4/4/4
5	BGC	C	602	-	-	0/2/22/22	0/1/1/1
2	BLA	D	601	1	-	1/22/74/74	0/4/4/4

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	601	BLA	C3C-C4C	-3.60	1.40	1.45
2	A	601	BLA	C3C-C4C	-3.33	1.40	1.45
2	A	601	BLA	C4D-C3D	-3.30	1.40	1.45
2	D	601	BLA	C3C-C4C	-3.24	1.40	1.45
2	D	601	BLA	C4D-C3D	-3.18	1.40	1.45
2	B	601	BLA	C3C-C4C	-3.18	1.40	1.45
2	C	601	BLA	C4D-C3D	-3.17	1.40	1.45
2	B	601	BLA	C4D-C3D	-3.08	1.40	1.45
2	A	601	BLA	C1C-C2C	-2.89	1.39	1.47
2	C	601	BLA	C1C-C2C	-2.89	1.39	1.47
2	D	601	BLA	C1C-C2C	-2.87	1.39	1.47
2	B	601	BLA	C1C-C2C	-2.79	1.39	1.47
2	B	601	BLA	C3B-C4B	-2.73	1.39	1.47
2	A	601	BLA	C3B-C4B	-2.71	1.39	1.47
2	C	601	BLA	C1B-C2B	-2.68	1.40	1.45
2	D	601	BLA	C1B-C2B	-2.67	1.40	1.45
2	C	601	BLA	C1D-C2D	-2.66	1.39	1.45
2	D	601	BLA	C3B-C4B	-2.65	1.39	1.47
2	D	601	BLA	C1D-C2D	-2.65	1.40	1.45
2	C	601	BLA	C3B-C4B	-2.65	1.39	1.47
2	B	601	BLA	C1D-C2D	-2.60	1.40	1.45
2	A	601	BLA	C1B-C2B	-2.58	1.40	1.45
2	A	601	BLA	C1D-C2D	-2.58	1.40	1.45
2	C	601	BLA	CHB-C1B	2.58	1.39	1.34
2	D	601	BLA	CHB-C1B	2.62	1.39	1.34
2	A	601	BLA	CHB-C1B	2.66	1.39	1.34
2	B	601	BLA	CHB-C1B	3.32	1.41	1.34
2	D	601	BLA	C4A-CHB	3.40	1.53	1.40
2	A	601	BLA	C4A-CHB	3.46	1.53	1.40
2	C	601	BLA	C4A-CHB	3.46	1.53	1.40
2	B	601	BLA	C4A-CHB	3.74	1.54	1.40
2	C	601	BLA	CBC-CAC	4.07	1.50	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	601	BLA	CBC-CAC	4.12	1.50	1.30
2	D	601	BLA	CBC-CAC	4.13	1.50	1.30
2	B	601	BLA	CBC-CAC	4.20	1.50	1.30
2	C	601	BLA	CHD-C1D	5.25	1.53	1.40
2	D	601	BLA	CHD-C1D	5.37	1.53	1.40
2	A	601	BLA	CHD-C1D	5.39	1.53	1.40
2	B	601	BLA	CHD-C1D	5.47	1.53	1.40
2	C	601	BLA	CHA-C4D	18.76	1.52	1.35
2	A	601	BLA	CHA-C4D	18.87	1.52	1.35
2	D	601	BLA	CHA-C4D	18.88	1.52	1.35
2	B	601	BLA	CHA-C4D	19.80	1.53	1.35

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	BLA	CHB-C1B-NB	-3.49	118.13	130.81
2	B	601	BLA	CMB-C2B-C3B	-2.60	122.35	128.58
2	B	601	BLA	C3A-C4A-NA	-2.37	106.40	110.29
2	D	601	BLA	CHB-C1B-NB	-2.25	122.62	130.81
2	A	601	BLA	CHB-C1B-NB	-2.21	122.79	130.81
2	D	601	BLA	C4C-CHD-C1D	-2.16	122.59	128.03
2	C	601	BLA	CHB-C1B-NB	-2.08	123.25	130.81
2	A	601	BLA	CAD-CBD-CGD	2.26	117.18	112.78
2	D	601	BLA	CAD-CBD-CGD	2.37	117.39	112.78
2	C	601	BLA	CAD-CBD-CGD	2.51	117.67	112.78
2	C	601	BLA	CAA-CBA-CGA	2.52	117.68	112.78
2	C	601	BLA	CMB-C2B-C1B	2.66	127.67	124.20
2	D	601	BLA	CAA-CBA-CGA	2.74	118.12	112.78
2	C	601	BLA	CBA-CAA-C2A	2.83	117.47	112.49
2	D	601	BLA	CMB-C2B-C1B	2.83	127.90	124.20
2	A	601	BLA	CAA-CBA-CGA	3.16	118.94	112.78
2	A	601	BLA	CMB-C2B-C1B	3.31	128.52	124.20
2	B	601	BLA	CAA-CBA-CGA	3.40	119.40	112.78
2	D	601	BLA	CBA-CAA-C2A	3.68	118.96	112.49
2	A	601	BLA	CBA-CAA-C2A	3.69	118.99	112.49
2	B	601	BLA	CBA-CAA-C2A	4.20	119.88	112.49
2	B	601	BLA	CHB-C1B-C2B	4.36	135.43	126.96
2	B	601	BLA	CMB-C2B-C1B	4.59	130.18	124.20

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	601	BLA	C4A-CHB-C1B-NB
2	A	601	BLA	C4A-CHB-C1B-NB
2	D	601	BLA	C4A-CHB-C1B-NB
2	B	601	BLA	C4A-CHB-C1B-NB

There are no ring outliers.

5 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	BLA	4	0
3	A	602	EDO	2	0
2	B	601	BLA	6	0
2	C	601	BLA	2	0
2	D	601	BLA	4	0

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	490/524 (93%)	0.16	11 (2%) 65 65	60, 121, 201, 257	0
1	B	486/524 (92%)	0.20	13 (2%) 58 57	66, 121, 213, 273	1 (0%)
1	C	476/524 (90%)	0.11	18 (3%) 44 42	102, 169, 231, 297	0
1	D	474/524 (90%)	0.34	27 (5%) 27 26	106, 182, 247, 313	0
All	All	1926/2096 (91%)	0.20	69 (3%) 46 45	60, 156, 228, 313	1 (0%)

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	44	ALA	6.4
1	B	430	VAL	4.3
1	B	402	LEU	4.3
1	D	408	LEU	3.7
1	B	437	CYS	3.6
1	A	106	PRO	3.6
1	B	369	ILE	3.5
1	D	103	LEU	3.4
1	D	54	MET	3.4
1	D	40	ALA	3.3
1	B	438	LEU	3.3
1	C	126	PHE	3.1
1	C	65	GLU	3.1
1	B	440	TRP	3.0
1	D	124	LEU	3.0
1	D	69	LEU	3.0
1	D	126	PHE	2.9
1	B	368	LEU	2.9
1	B	340	LEU	2.9
1	C	149	SER	2.9
1	A	438	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	108	ALA	2.9
1	A	430	VAL	2.8
1	A	437	CYS	2.8
1	D	51	VAL	2.7
1	C	502	LEU	2.7
1	A	77	LEU	2.7
1	A	346	LEU	2.6
1	D	440	TRP	2.6
1	D	42	LEU	2.6
1	D	428	ILE	2.6
1	D	194	GLY	2.6
1	D	437	CYS	2.6
1	C	44	ALA	2.5
1	D	176	TYR	2.5
1	C	39	GLY	2.5
1	D	286	LEU	2.5
1	C	150	ALA	2.5
1	D	190	ALA	2.5
1	D	111	LEU	2.4
1	C	69	LEU	2.4
1	D	122	LEU	2.4
1	D	402	LEU	2.4
1	A	428	ILE	2.4
1	C	432	GLU	2.4
1	C	430	VAL	2.4
1	B	451	TRP	2.3
1	C	18	GLU	2.3
1	D	410	GLN	2.3
1	D	41	LEU	2.3
1	D	340	LEU	2.3
1	D	438	LEU	2.3
1	B	428	ILE	2.3
1	A	498	LEU	2.3
1	D	287	ILE	2.2
1	C	346	LEU	2.2
1	B	370	LEU	2.1
1	C	103	LEU	2.1
1	C	469	TRP	2.1
1	A	402	LEU	2.1
1	D	449	VAL	2.1
1	C	408	LEU	2.1
1	B	338	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	403	VAL	2.1
1	C	81	GLN	2.1
1	A	103	LEU	2.0
1	B	439	VAL	2.0
1	D	52	LEU	2.0
1	C	68	VAL	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	LLDF	B-factors(\AA^2)	Q<0.9
3	EDO	A	602	4/4	0.87	0.50	14.70	101,108,130,143	0
5	BGC	C	602	12/12	0.86	0.31	3.51	139,208,225,229	0
2	BLA	B	601	43/43	0.89	0.34	1.29	72,109,164,187	0
2	BLA	C	601	43/43	0.92	0.35	1.18	112,151,220,232	0
2	BLA	A	601	43/43	0.93	0.31	1.06	54,87,121,156	0
4	ACT	A	603	4/4	0.95	0.31	0.87	94,120,123,135	0
2	BLA	D	601	43/43	0.87	0.34	0.63	128,168,232,252	0

6.5 Other polymers [i](#)

There are no such residues in this entry.