



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:53 PM UTC

PDB ID : 6FHT / pdb_00006fht
Title : Crystal structure of an artificial phytochrome regulated adenylate/guanylate cyclase in its dark adapted Pr form
Authors : Etzl, S.; Winkler, A.
Deposited on : 2018-01-15
Resolution : 2.35 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

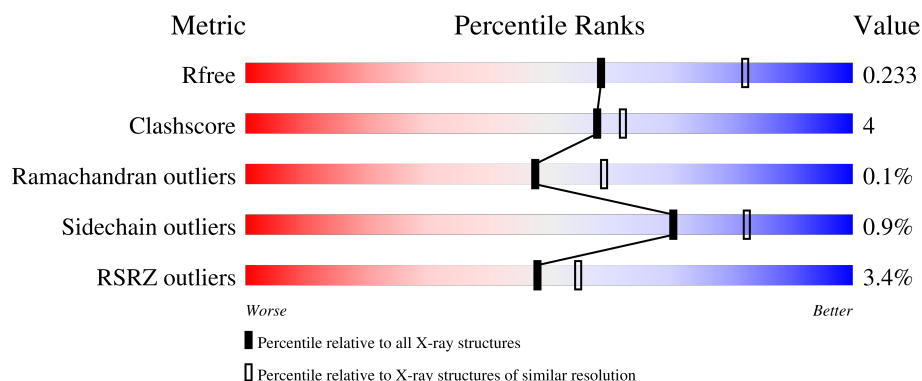
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	741	4% 88% 10%
1	B	741	3% 89% 10%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 11429 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.

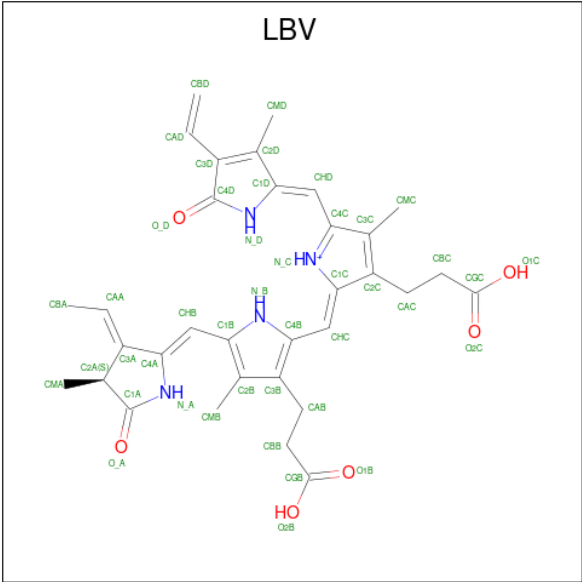
- Molecule 1 is a protein called Bacteriophytochrome,Adenylate cyclase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	730	Total	C	N	O	S	0	0	0
			5542	3508	978	1038	18			
1	B	735	Total	C	N	O	S	0	0	0
			5572	3526	983	1045	18			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	GLY	-	expression tag	UNP Q9RZA4
A	0	ALA	-	expression tag	UNP Q9RZA4
A	1	MET	-	expression tag	UNP Q9RZA4
A	2	ALA	-	expression tag	UNP Q9RZA4
A	589	LYS	GLU	engineered mutation	UNP P72951
B	-1	GLY	-	expression tag	UNP Q9RZA4
B	0	ALA	-	expression tag	UNP Q9RZA4
B	1	MET	-	expression tag	UNP Q9RZA4
B	2	ALA	-	expression tag	UNP Q9RZA4
B	589	LYS	GLU	engineered mutation	UNP P72951

- 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-methyl-5-oxidanylidene-pyrrolidin-2-ylidene]methyl]-4-methyl-1H-pyrrol-3-yl]propanoic acid (CCD ID: LBV) (formula: C₃₃H₃₇N₄O₆).



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		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



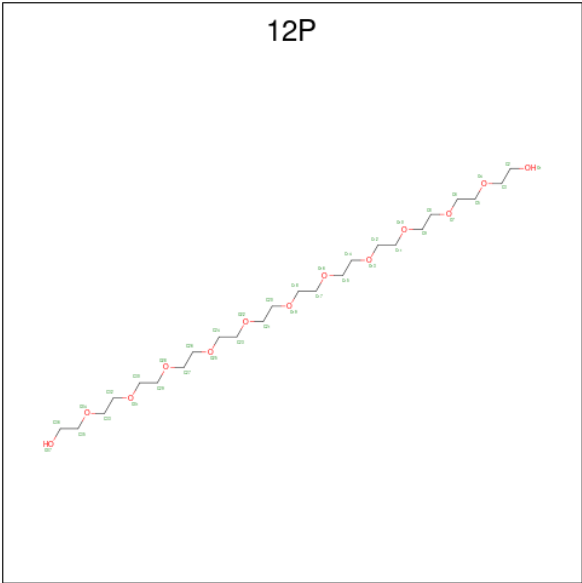
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is DODECAETHYLENE GLYCOL (CCD ID: 12P) (formula: C₂₄H₅₀O₁₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			37	24	13		

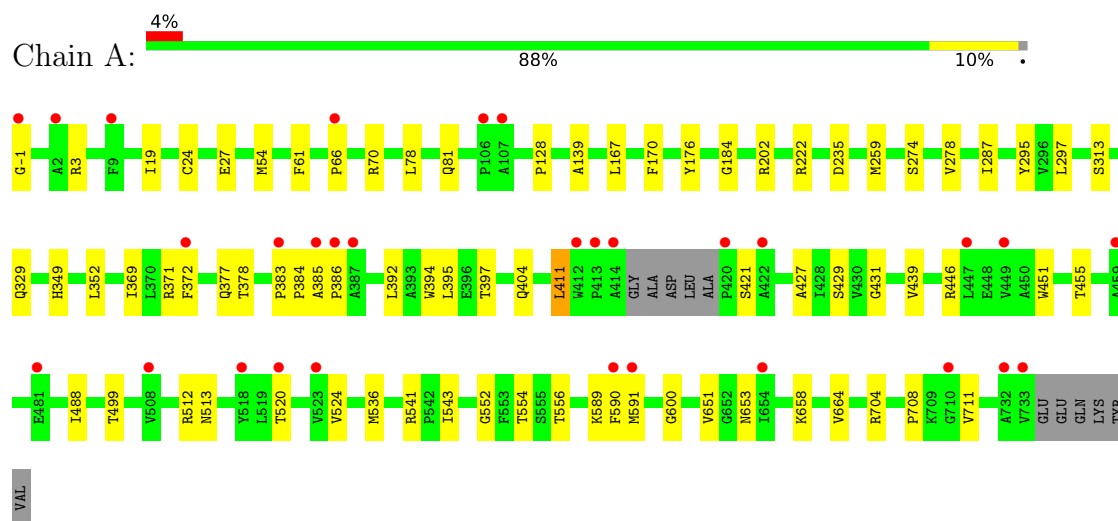
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	85	Total 85	O 85	0	0
5	B	57	Total 57	O 57	0	0

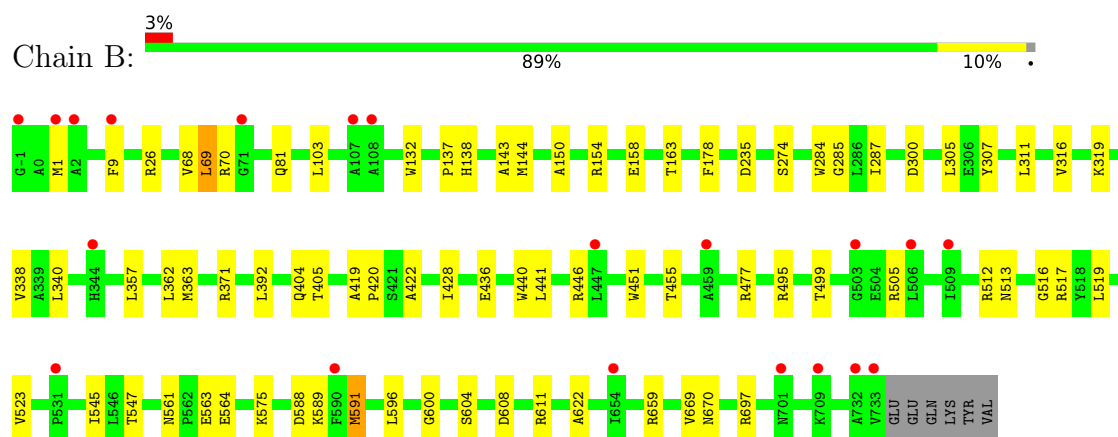
3 Residue-property plots

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,Adenylate cyclase



- Molecule 1: Bacteriophytochrome,Adenylate cyclase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	286.19Å 108.82Å 67.99Å 90.00° 100.84° 90.00°	Depositor
Resolution (Å)	46.85 – 2.35 46.85 – 2.35	Depositor EDS
% Data completeness (in resolution range)	96.2 (46.85-2.35) 96.2 (46.85-2.35)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.61 (at 2.34Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.180 , 0.227 0.185 , 0.233	Depositor DCC
R_{free} test set	4102 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	55.1	Xtriage
Anisotropy	0.322	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 46.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.074 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11429	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.21% 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: LBV, SO4, 12P

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	0.33	0/5669	0.50	0/7740
1	B	0.32	0/5700	0.51	0/7785
All	All	0.33	0/11369	0.51	0/15525

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	5542	0	5524	54	0
1	B	5572	0	5552	53	0
2	A	43	0	33	1	0
2	B	43	0	33	0	0
3	A	30	0	0	1	0
3	B	20	0	0	0	0
4	A	37	0	50	1	0
5	A	85	0	0	2	0
5	B	57	0	0	0	0
All	All	11429	0	11192	100	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including

hydrogen atoms). The all-atom clashscore for this structure is 4.

All (100) 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:591:MET:HE2	1:B:591:MET:HG2	1.71	0.73
1:B:589:LYS:HD3	1:B:591:MET:HE3	1.70	0.72
1:B:371:ARG:HE	1:B:392:LEU:HD13	1.54	0.71
1:A:70:ARG:NH2	1:A:235:ASP:O	2.24	0.71
1:B:512:ARG:NH1	1:B:519:LEU:O	2.25	0.70
1:A:512:ARG:NH2	3:A:804:SO4:O2	2.27	0.67
1:A:404:GLN:HG3	1:A:488:ILE:HD11	1.76	0.66
1:A:536:MET:HE3	1:A:653:ASN:HB2	1.79	0.64
1:A:371:ARG:HH21	1:A:392:LEU:HB3	1.64	0.61
1:B:446:ARG:HG2	1:B:477:ARG:HH21	1.68	0.58
1:A:383:PRO:HD3	1:A:421:SER:HB3	1.85	0.58
1:B:143:ALA:CB	1:B:163:THR:HG21	2.34	0.57
1:A:54:MET:HE1	1:A:66:PRO:HB3	1.86	0.57
1:A:541:ARG:NH2	1:A:600:GLY:O	2.38	0.57
1:A:384:PRO:HB2	1:A:386:PRO:HD2	1.86	0.57
1:A:139:ALA:HB1	1:A:167:LEU:HD21	1.88	0.56
1:B:143:ALA:HB2	1:B:163:THR:HG21	1.87	0.55
1:B:519:LEU:HD22	1:B:523:VAL:HG11	1.88	0.54
1:B:589:LYS:HB3	1:B:591:MET:HG3	1.89	0.54
1:A:19:ILE:HG13	1:A:27:GLU:HG3	1.90	0.53
1:A:589:LYS:HD2	1:B:589:LYS:HG2	1.90	0.53
1:A:-1:GLY:N	1:A:3:ARG:HH21	2.07	0.53
1:B:150:ALA:O	1:B:319:LYS:HE3	2.10	0.52
1:A:589:LYS:HG3	1:A:591:MET:HE3	1.93	0.51
1:B:371:ARG:NH2	1:B:392:LEU:HB3	2.26	0.51
1:A:352:LEU:HD12	1:A:372:PHE:HE2	1.75	0.51
1:A:554:THR:HG21	1:B:670:ASN:ND2	2.26	0.51
1:B:600:GLY:HA2	1:B:604:SER:HA	1.92	0.50
1:A:513:ASN:ND2	1:B:513:ASN:OD1	2.45	0.50
1:B:419:ALA:N	1:B:420:PRO:HD2	2.26	0.50
1:B:451:TRP:CD1	1:B:455:THR:HG22	2.46	0.50
1:A:295:TYR:CZ	4:A:808:12P:H182	2.47	0.50
1:A:499:THR:HG22	1:B:499:THR:HG22	1.94	0.49
1:B:575:LYS:HG3	1:B:622:ALA:HB1	1.94	0.49
1:A:451:TRP:CD1	1:A:455:THR:HG22	2.48	0.49
1:A:176:TYR:CZ	1:A:184:GLY:HA3	2.48	0.48
1:B:9:PHE:CE2	1:B:26:ARG:HD2	2.48	0.48
1:A:349:HIS:CE1	1:A:377:GLN:HG2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:589:LYS:HB3	1:A:591:MET:HG3	1.96	0.47
1:B:144:MET:HG3	1:B:311:LEU:HD21	1.97	0.47
1:A:352:LEU:HD12	1:A:372:PHE:CE2	2.50	0.47
1:A:395:LEU:HD21	1:A:427:ALA:HB1	1.96	0.47
1:A:651:VAL:HG22	1:A:664:VAL:HG12	1.96	0.47
1:A:552:GLY:O	1:A:556:THR:HG23	2.14	0.46
1:A:658:LYS:HE2	1:A:658:LYS:H	1.80	0.46
1:B:608:ASP:OD1	1:B:611:ARG:NH2	2.48	0.46
1:B:9:PHE:CZ	1:B:26:ARG:HD2	2.50	0.46
1:B:338:VAL:HG22	1:B:357:LEU:HG	1.97	0.46
1:B:697:ARG:HA	1:B:697:ARG:HE	1.81	0.46
1:A:520:THR:O	1:A:524:VAL:HG23	2.14	0.46
1:B:274:SER:HA	1:B:287:ILE:O	2.16	0.46
1:B:561:ASN:HB2	1:B:564:GLU:HG3	1.97	0.46
1:A:371:ARG:NH2	1:A:392:LEU:HB3	2.31	0.46
1:B:363:MET:HG3	1:B:440:TRP:CE2	2.51	0.46
1:A:19:ILE:HG22	1:A:259:MET:HE3	1.98	0.45
1:A:-1:GLY:H2	1:A:3:ARG:HH21	1.65	0.45
1:B:69:LEU:HD23	1:B:69:LEU:HA	1.88	0.44
1:A:429:SER:OG	1:A:431:GLY:O	2.26	0.44
1:A:513:ASN:O	1:B:517:ARG:NH1	2.28	0.44
1:A:78:LEU:HB3	1:A:81:GLN:HB2	1.99	0.44
1:B:516:GLY:HA3	1:B:519:LEU:HG	1.99	0.44
1:B:422:ALA:HB2	1:B:441:LEU:HD13	2.00	0.44
1:B:404:GLN:HG2	1:B:405:THR:N	2.33	0.44
1:B:436:GLU:OE2	1:B:505:ARG:NH2	2.44	0.44
1:B:436:GLU:CD	1:B:505:ARG:HH22	2.26	0.43
1:B:137:PRO:HG2	1:B:138:HIS:CE1	2.54	0.43
1:B:545:ILE:HD13	1:B:669:VAL:HA	2.00	0.43
1:A:274:SER:HA	1:A:287:ILE:O	2.18	0.43
1:B:178:PHE:CD2	1:B:285:GLY:HA2	2.54	0.43
1:B:307:TYR:CE2	1:B:311:LEU:HD22	2.53	0.43
1:B:307:TYR:HE2	1:B:311:LEU:HD22	1.84	0.43
1:A:394:TRP:O	1:A:397:THR:HB	2.19	0.42
1:A:170:PHE:CZ	1:A:297:LEU:HD11	2.54	0.42
1:A:369:ILE:HG12	1:A:378:THR:HG22	2.01	0.42
1:B:154:ARG:O	1:B:158:GLU:HG2	2.20	0.42
1:B:588:ASP:O	1:B:589:LYS:HG3	2.19	0.42
1:A:394:TRP:HZ2	1:A:411:LEU:HD22	1.84	0.42
1:A:704:ARG:HD2	5:A:935:HOH:O	2.19	0.42
1:B:81:GLN:HG3	1:B:103:LEU:HD21	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:371:ARG:HH21	1:B:392:LEU:HB3	1.84	0.42
1:B:547:THR:HG22	1:B:596:LEU:HD12	2.00	0.42
1:A:24:CYS:HB2	2:A:801:LBV:HAA1	1.95	0.42
1:B:697:ARG:HA	1:B:697:ARG:NE	2.35	0.42
1:B:371:ARG:NE	1:B:392:LEU:HD13	2.30	0.41
1:B:70:ARG:NH2	1:B:235:ASP:O	2.53	0.41
1:A:446:ARG:HA	1:A:446:ARG:HD3	1.92	0.41
1:A:349:HIS:CD2	1:A:372:PHE:HB2	2.55	0.41
1:A:278:VAL:O	1:A:313:SER:HB3	2.19	0.41
1:A:543:ILE:HD13	1:A:664:VAL:HG11	2.02	0.41
1:B:1:MET:HE3	1:B:1:MET:HB3	1.97	0.41
1:B:284:TRP:CE2	1:B:316:VAL:HG11	2.55	0.41
1:A:512:ARG:NH2	1:B:659:ARG:HE	2.19	0.41
1:A:61:PHE:HE2	1:A:128:PRO:HD3	1.85	0.41
1:A:385:ALA:N	1:A:386:PRO:HD2	2.36	0.41
1:B:362:LEU:HD23	1:B:363:MET:HE2	2.01	0.41
1:A:202:ARG:HD2	5:A:976:HOH:O	2.21	0.40
1:A:369:ILE:HB	1:A:439:VAL:HB	2.03	0.40
1:A:708:PRO:HG2	1:A:711:VAL:HG21	2.03	0.40
1:A:404:GLN:HG3	1:A:488:ILE:CD1	2.49	0.40
1:B:428:ILE:HD12	1:B:495:ARG:HB2	2.02	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	726/741 (98%)	712 (98%)	14 (2%)	0	100	100
1	B	733/741 (99%)	721 (98%)	11 (2%)	1 (0%)	48	59
All	All	1459/1482 (98%)	1433 (98%)	25 (2%)	1 (0%)	48	59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	132	TRP

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	580/588 (99%)	576 (99%)	4 (1%)	76	86
1	B	582/588 (99%)	575 (99%)	7 (1%)	63	77
All	All	1162/1176 (99%)	1151 (99%)	11 (1%)	70	82

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	222	ARG
1	A	329	GLN
1	A	411	LEU
1	A	590	PHE
1	B	68	VAL
1	B	69	LEU
1	B	300	ASP
1	B	305	LEU
1	B	340	LEU
1	B	563	GLU
1	B	591	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	377	GLN
1	A	398	GLN
1	A	629	GLN
1	A	653	ASN
1	B	72	GLN

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Mol	Chain	Res	Type
1	B	196	HIS
1	B	467	HIS
1	B	728	ASN

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 ⓘ

13 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	$\# Z > 2$
4	12P	A	808	-	36,36,36	0.54	0	35,35,35	0.41	0
3	SO4	A	805	-	4,4,4	0.25	0	6,6,6	0.08	0
3	SO4	B	803	-	4,4,4	0.23	0	6,6,6	0.22	0
3	SO4	B	804	-	4,4,4	0.26	0	6,6,6	0.17	0
3	SO4	B	802	-	4,4,4	0.25	0	6,6,6	0.15	0
3	SO4	A	804	-	4,4,4	0.22	0	6,6,6	0.21	0
3	SO4	A	806	-	4,4,4	0.23	0	6,6,6	0.20	0
3	SO4	A	807	-	4,4,4	0.28	0	6,6,6	0.21	0
2	LBV	A	801	1	46,46,46	2.47	14 (30%)	58,67,67	1.27	6 (10%)
3	SO4	B	805	-	4,4,4	0.25	0	6,6,6	0.06	0
3	SO4	A	803	-	4,4,4	0.25	0	6,6,6	0.10	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	A	802	-	4,4,4	0.25	0	6,6,6	0.15	0
2	LBV	B	801	1	46,46,46	2.33	12 (26%)	58,67,67	1.22	8 (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
4	12P	A	808	-	-	20/34/34/34	-
2	LBV	A	801	1	-	8/26/74/74	0/4/4/4
2	LBV	B	801	1	-	9/26/74/74	0/4/4/4

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	LBV	C2A-C1A	-8.65	1.42	1.52
2	B	801	LBV	C2A-C1A	-7.79	1.43	1.52
2	B	801	LBV	C2A-C3A	-6.80	1.42	1.51
2	A	801	LBV	C2A-C3A	-5.95	1.44	1.51
2	A	801	LBV	CMC-C3C	-4.94	1.40	1.50
2	B	801	LBV	CMD-C2D	-4.72	1.41	1.50
2	A	801	LBV	CMB-C2B	-4.67	1.41	1.50
2	A	801	LBV	CMD-C2D	-4.65	1.41	1.50
2	B	801	LBV	CMB-C2B	-4.45	1.41	1.50
2	A	801	LBV	CMA-C2A	-3.99	1.41	1.53
2	B	801	LBV	CMA-C2A	-3.66	1.42	1.53
2	B	801	LBV	CMC-C3C	-3.28	1.44	1.50
2	B	801	LBV	CBB-CGB	-3.02	1.43	1.50
2	A	801	LBV	CAA-C3A	2.87	1.40	1.33
2	A	801	LBV	CAB-C3B	-2.75	1.44	1.51
2	B	801	LBV	CAB-C3B	-2.65	1.44	1.51
2	A	801	LBV	CBB-CGB	-2.65	1.44	1.50
2	A	801	LBV	O2B-CGB	-2.59	1.22	1.30
2	A	801	LBV	CHC-C1C	2.26	1.43	1.38
2	B	801	LBV	CAA-C3A	2.20	1.39	1.33
2	A	801	LBV	CBD-CAD	2.19	1.40	1.30
2	A	801	LBV	CAC-C2C	-2.16	1.45	1.51
2	B	801	LBV	C4B-N_B	2.14	1.41	1.37
2	B	801	LBV	CAC-C2C	-2.10	1.45	1.51
2	A	801	LBV	C1D-N_D	-2.06	1.34	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	801	LBV	CHC-C1C	2.04	1.42	1.38

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	LBV	C4B-CHC-C1C	3.41	135.66	128.22
2	A	801	LBV	C4B-CHC-C1C	3.11	135.02	128.22
2	A	801	LBV	C2C-C1C-N_C	-2.82	106.05	110.04
2	B	801	LBV	C2C-C1C-N_C	-2.78	106.11	110.04
2	B	801	LBV	C1C-N_C-C4C	2.45	111.02	106.52
2	A	801	LBV	O_D-C4D-N_D	-2.43	119.39	125.08
2	B	801	LBV	O_A-C1A-C2A	2.34	129.22	126.74
2	A	801	LBV	CHD-C4C-C3C	2.32	130.83	124.87
2	A	801	LBV	O_A-C1A-N_A	-2.26	122.26	124.93
2	B	801	LBV	O_D-C4D-N_D	-2.21	119.90	125.08
2	B	801	LBV	O_A-C1A-N_A	-2.14	122.41	124.93
2	A	801	LBV	C1C-N_C-C4C	2.11	110.39	106.52
2	B	801	LBV	O2C-CGC-CBC	-2.11	116.41	123.09
2	B	801	LBV	O1B-CGB-CBB	-2.03	116.64	123.09

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	801	LBV	N_C-C4C-CHD-C1D
2	B	801	LBV	C3C-C4C-CHD-C1D
4	A	808	12P	O10-C11-C12-O13
4	A	808	12P	O25-C26-C27-O28
4	A	808	12P	O13-C14-C15-O16
4	A	808	12P	C11-C12-O13-C14
2	A	801	LBV	N_D-C1D-CHD-C4C
2	B	801	LBV	N_D-C1D-CHD-C4C
4	A	808	12P	O22-C23-C24-O25
4	A	808	12P	O19-C20-C21-O22
2	A	801	LBV	N_C-C4C-CHD-C1D
2	A	801	LBV	C3C-C4C-CHD-C1D
2	A	801	LBV	C2D-C1D-CHD-C4C
2	B	801	LBV	C2D-C1D-CHD-C4C
4	A	808	12P	C27-C26-O25-C24
4	A	808	12P	O16-C17-C18-O19
4	A	808	12P	O34-C35-C36-O37
4	A	808	12P	O28-C29-C30-O31

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Mol	Chain	Res	Type	Atoms
4	A	808	12P	C9-C8-O7-C6
4	A	808	12P	C2-C3-O4-C5
4	A	808	12P	C29-C30-O31-C32
4	A	808	12P	O7-C8-C9-O10
2	B	801	LBV	C2C-CAC-CBC-CGC
4	A	808	12P	C20-C21-O22-C23
4	A	808	12P	O1-C2-C3-O4
4	A	808	12P	C21-C20-O19-C18
4	A	808	12P	C36-C35-O34-C33
4	A	808	12P	C14-C15-O16-C17
2	A	801	LBV	CAB-CBB-CGB-O1B
2	B	801	LBV	CAB-CBB-CGB-O2B
2	A	801	LBV	CAB-CBB-CGB-O2B
2	B	801	LBV	CAB-CBB-CGB-O1B
2	B	801	LBV	CAC-CBC-CGC-O2C
2	A	801	LBV	CAC-CBC-CGC-O1C
2	A	801	LBV	CAC-CBC-CGC-O2C
2	B	801	LBV	CAC-CBC-CGC-O1C
4	A	808	12P	C18-C17-O16-C15

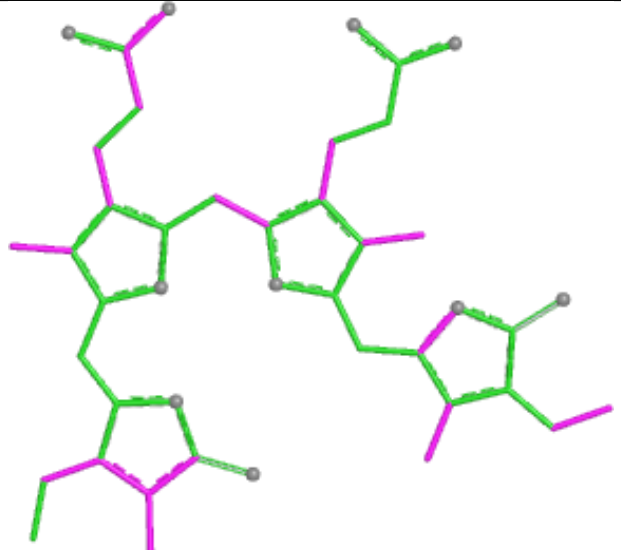
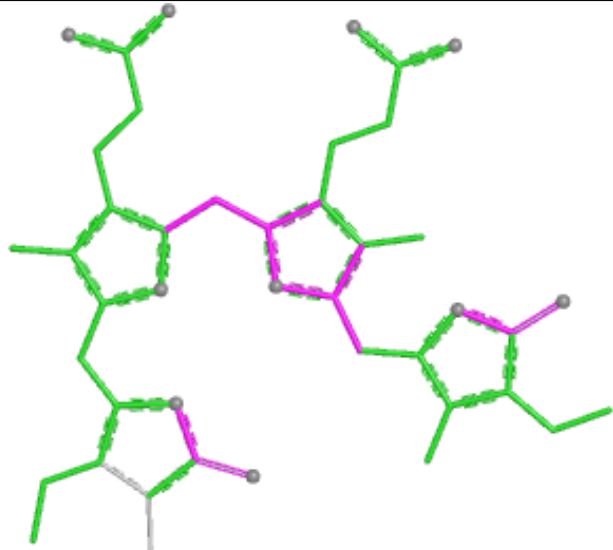
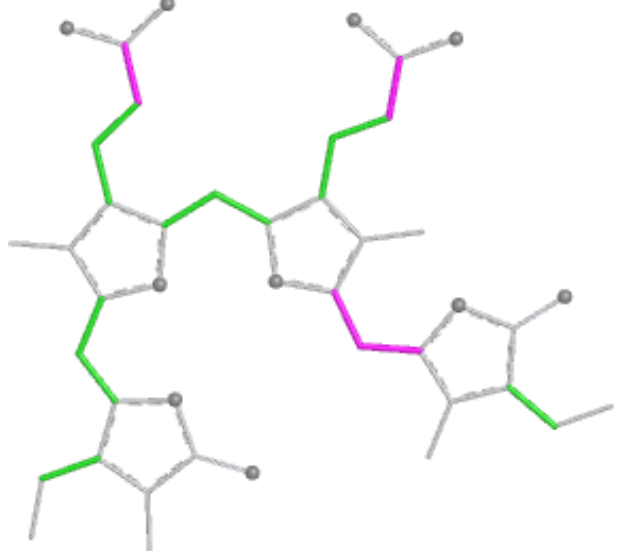
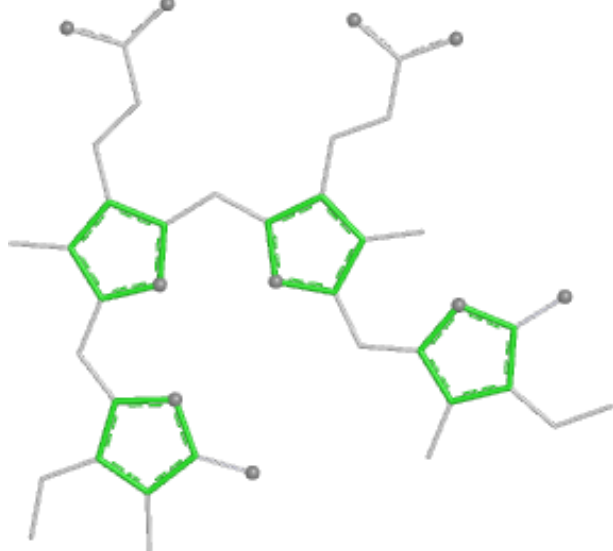
There are no ring outliers.

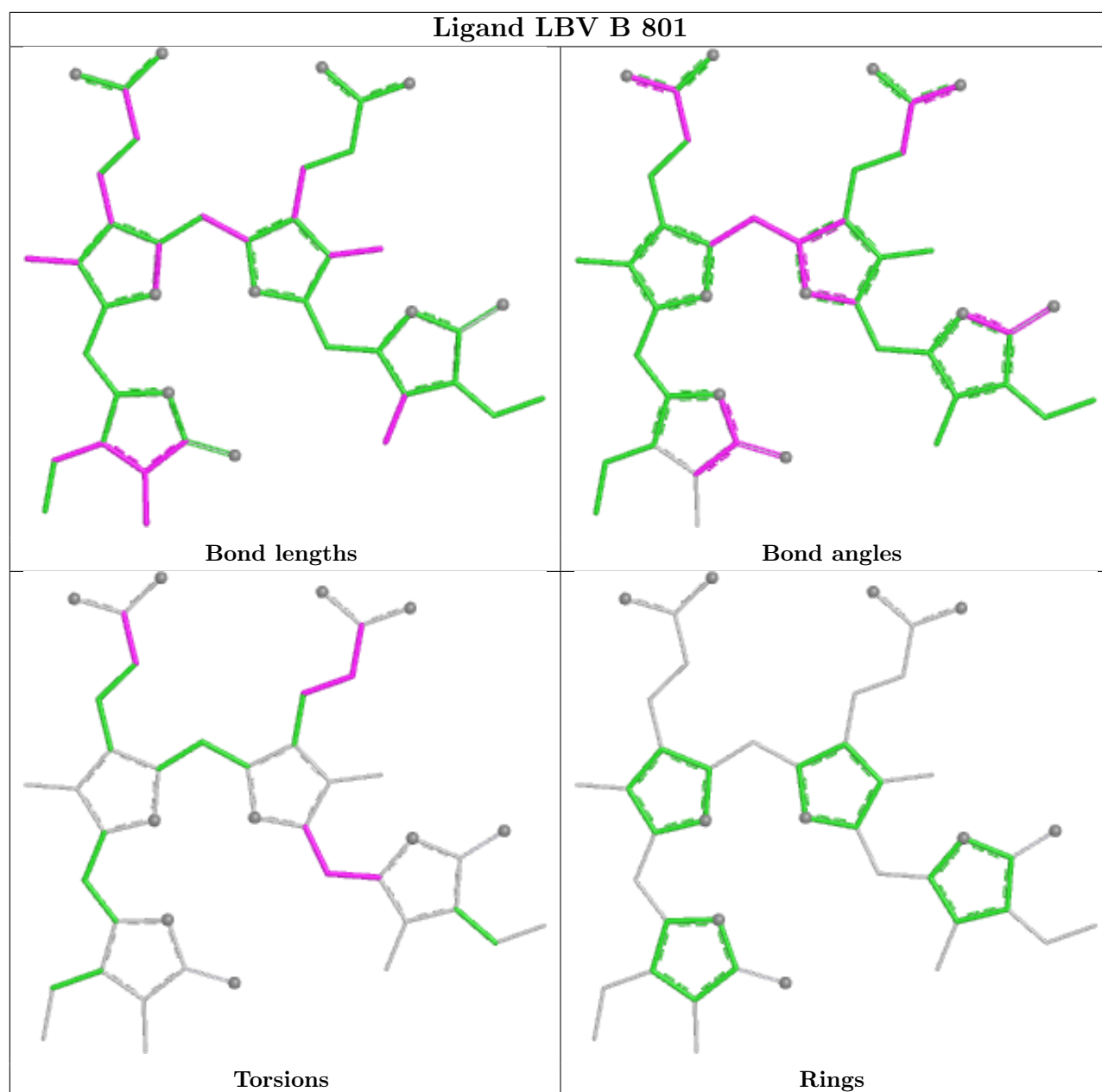
3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	808	12P	1	0
3	A	804	SO4	1	0
2	A	801	LBV	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.

Ligand 12P A 808	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand LBV A 801	
 Bond lengths	 Bond angles
 Torsions	 Rings



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	730/741 (98%)	0.17	30 (4%)	41 47	32, 62, 113, 139	0
1	B	735/741 (99%)	0.12	20 (2%)	56 63	37, 63, 108, 138	0
All	All	1465/1482 (98%)	0.14	50 (3%)	48 55	32, 63, 111, 139	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	420	PRO	6.2
1	A	459	ALA	4.6
1	B	733	VAL	4.2
1	B	732	ALA	3.8
1	A	414	ALA	3.6
1	B	2	ALA	3.5
1	A	733	VAL	3.4
1	B	506	LEU	3.3
1	B	459	ALA	3.3
1	B	447	LEU	3.3
1	A	523	VAL	3.2
1	A	385	ALA	3.1
1	B	107	ALA	3.1
1	A	447	LEU	3.1
1	B	531	PRO	3.1
1	A	2	ALA	3.0
1	A	422	ALA	2.9
1	A	732	ALA	2.9
1	A	413	PRO	2.7
1	A	590	PHE	2.7
1	B	590	PHE	2.7
1	A	383	PRO	2.7
1	B	108	ALA	2.7
1	A	508	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	-1	GLY	2.5
1	A	710	GLY	2.5
1	A	449	VAL	2.4
1	B	654	ILE	2.4
1	A	518	TYR	2.4
1	A	106	PRO	2.4
1	A	386	PRO	2.4
1	B	1	MET	2.4
1	A	412	TRP	2.4
1	A	481	GLU	2.4
1	B	71	GLY	2.4
1	A	107	ALA	2.3
1	A	387	ALA	2.2
1	A	654	ILE	2.2
1	A	66	PRO	2.2
1	A	9	PHE	2.2
1	A	591	MET	2.2
1	B	9	PHE	2.2
1	A	520	THR	2.2
1	B	509	ILE	2.1
1	B	709	LYS	2.1
1	B	701	ASN	2.1
1	B	344	HIS	2.0
1	B	-1	GLY	2.0
1	B	503	GLY	2.0
1	A	372	PHE	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](#)

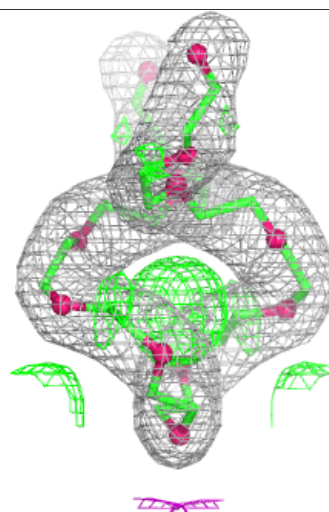
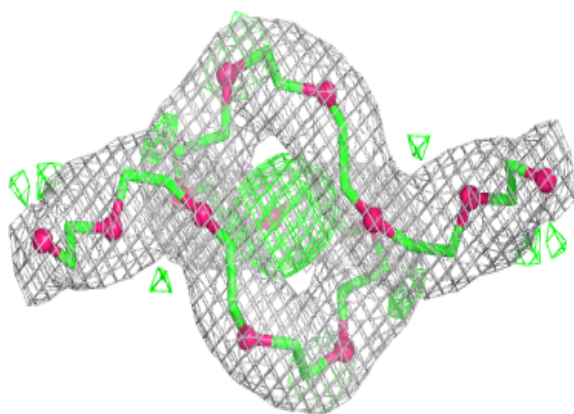
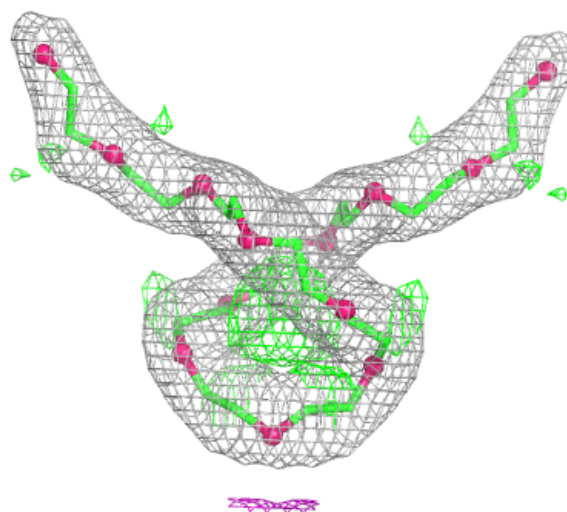
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	A	806	5/5	0.68	0.12	110,111,113,114	0
3	SO4	A	805	5/5	0.74	0.13	143,143,144,147	0
3	SO4	A	803	5/5	0.74	0.14	118,118,119,119	0
3	SO4	B	802	5/5	0.80	0.09	115,115,117,118	0
3	SO4	A	804	5/5	0.81	0.12	106,110,114,114	0
3	SO4	A	807	5/5	0.82	0.10	96,99,103,107	0
3	SO4	A	802	5/5	0.82	0.09	113,115,117,121	0
3	SO4	B	804	5/5	0.83	0.10	105,106,107,111	0
3	SO4	B	805	5/5	0.84	0.08	130,131,131,132	0
3	SO4	B	803	5/5	0.88	0.09	94,98,101,102	0
4	12P	A	808	37/37	0.92	0.12	53,60,76,77	37
2	LBV	A	801	43/43	0.96	0.07	28,43,55,59	0
2	LBV	B	801	43/43	0.97	0.06	35,47,56,62	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.

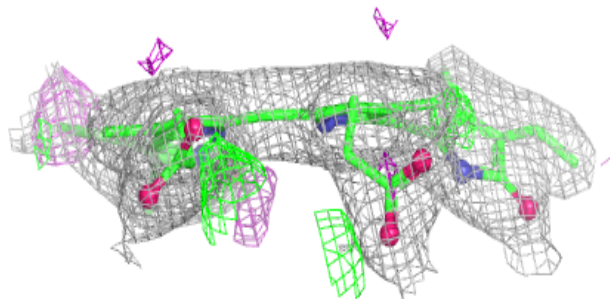
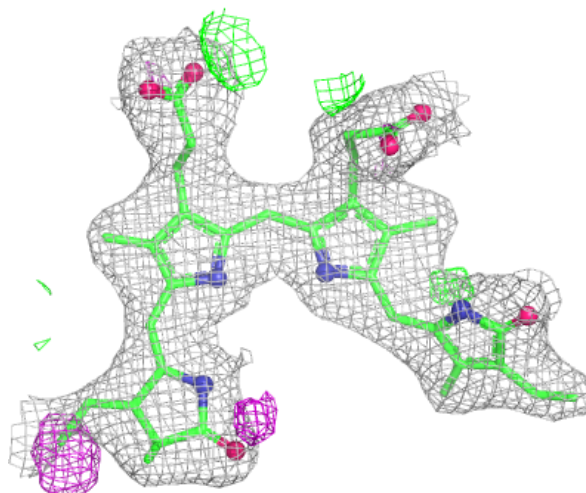
Electron density around 12P A 808:

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



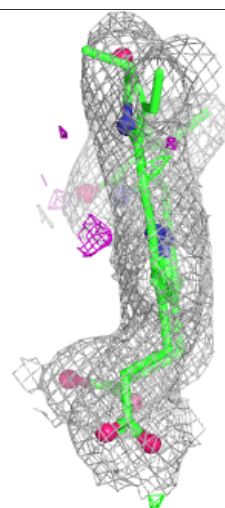
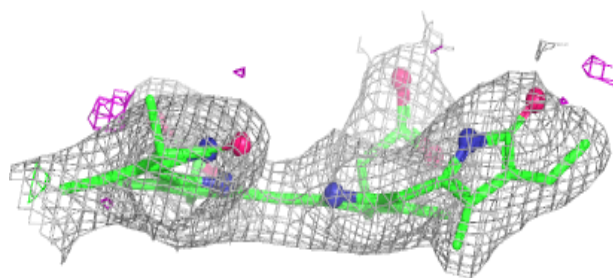
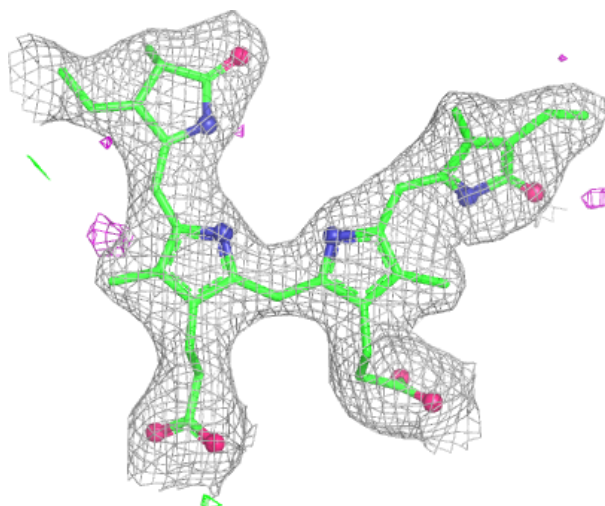
Electron density around LBV A 801:

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



Electron density around LBV B 801:

$2mF_o-DF_c$ (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.