



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 02:11 PM UTC

PDB ID : 3AJG / pdb_00003ajg
Title : Crystal structure of PcyA V225D-biliverdin IX alpha complex
Authors : Wada, K.; Hagiwara, Y.; Fukuyama, K.
Deposited on : 2010-06-05
Resolution : 1.90 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

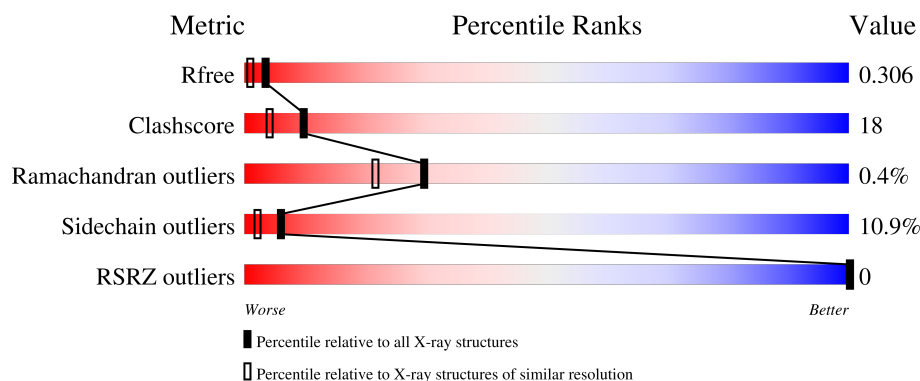
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	248	 65% 25% 6% .
1	B	248	 67% 24% 6% . .

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4139 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.

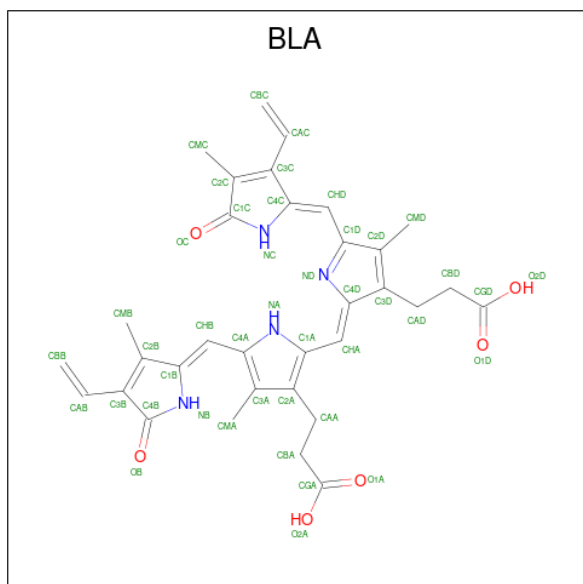
- Molecule 1 is a protein called Phycocyanobilin:ferredoxin oxidoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	238	Total	C	N	O	S	0	6	0
			1916	1221	326	356	13			
1	B	241	Total	C	N	O	S	0	3	0
			1935	1231	332	360	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	225	ASP	VAL	engineered mutation	UNP Q55891
B	225	ASP	VAL	engineered mutation	UNP Q55891

- Molecule 2 is BILIVERDINE IX ALPHA (CCD ID: BLA) (formula: $C_{33}H_{34}N_4O_6$).



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

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

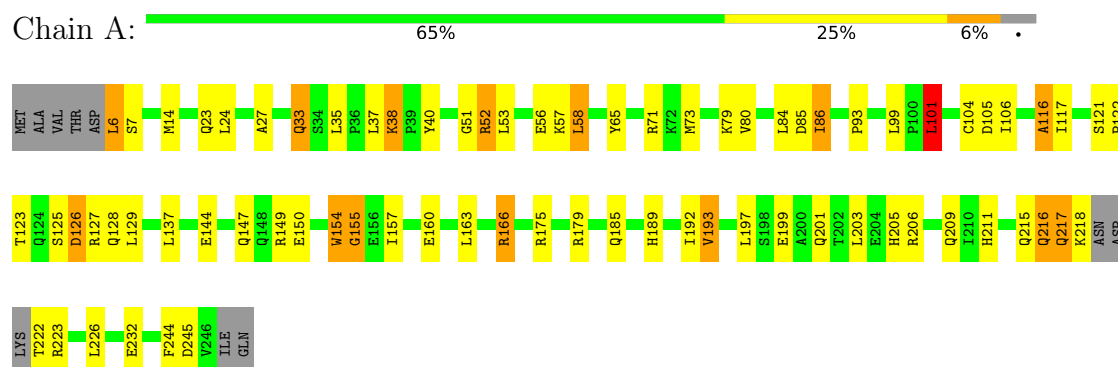
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	104	Total	O	0	0
			104	104		
3	B	98	Total	O	0	0
			98	98		

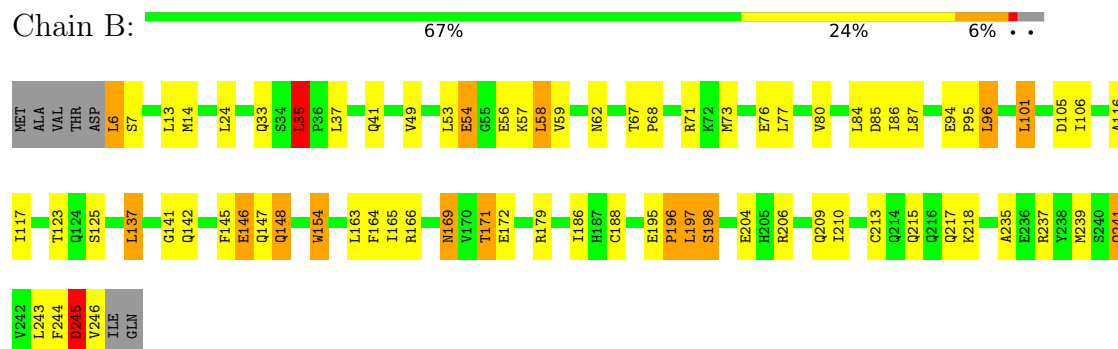
3 Residue-property plots [i](#)

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: Phycocyanobilin:ferredoxin oxidoreductase



- Molecule 1: Phycocyanobilin:ferredoxin oxidoreductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.05Å 74.88Å 84.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.88 – 1.90 32.88 – 1.91	Depositor EDS
% Data completeness (in resolution range)	98.3 (32.88-1.90) 96.1 (32.88-1.91)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.19 (at 1.91Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.231 , 0.295 0.245 , 0.306	Depositor DCC
R_{free} test set	1839 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	31.3	Xtriage
Anisotropy	0.289	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.149 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4139	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.04% 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: BLA

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	1.17	4/1986 (0.2%)	1.19	13/2694 (0.5%)
1	B	1.12	5/1994 (0.3%)	1.23	12/2704 (0.4%)
All	All	1.15	9/3980 (0.2%)	1.21	25/5398 (0.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2
1	B	0	1
All	All	0	3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	49	VAL	C-O	-7.69	1.15	1.24
1	A	117	ILE	C-N	-6.81	1.25	1.33
1	B	77	LEU	C-O	-6.13	1.16	1.24
1	A	86	ILE	C-O	-5.95	1.18	1.24
1	A	116	ALA	C-N	-5.83	1.25	1.33
1	B	14	MET	C-O	-5.64	1.19	1.24
1	B	171	THR	CA-CB	5.35	1.61	1.53
1	A	27	ALA	CA-CB	-5.17	1.45	1.53
1	B	105	ASP	C-N	-5.08	1.26	1.33

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	245	ASP	N-CA-C	8.50	122.55	110.23
1	A	157	ILE	N-CA-C	-8.02	102.57	113.00
1	B	85	ASP	CA-C-N	6.97	132.68	122.99
1	B	85	ASP	C-N-CA	6.97	132.68	122.99
1	B	14	MET	CA-C-N	-6.93	112.75	120.45
1	B	14	MET	C-N-CA	-6.93	112.75	120.45
1	A	121	SER	CA-C-N	6.07	126.39	119.83
1	A	121	SER	C-N-CA	6.07	126.39	119.83
1	B	13	LEU	N-CA-C	5.79	119.53	112.23
1	A	155	GLY	N-CA-C	-5.76	101.56	112.27
1	A	14	MET	CA-C-N	-5.62	113.89	119.56
1	A	14	MET	C-N-CA	-5.62	113.89	119.56
1	B	142	GLN	CA-C-N	5.61	126.07	120.52
1	B	142	GLN	C-N-CA	5.61	126.07	120.52
1	A	125	SER	CA-C-N	5.49	128.40	120.38
1	A	125	SER	C-N-CA	5.49	128.40	120.38
1	A	99	LEU	CA-C-N	5.44	125.42	120.03
1	A	99	LEU	C-N-CA	5.44	125.42	120.03
1	A	232	GLU	N-CA-C	5.44	118.13	111.82
1	B	244	PHE	CA-C-N	5.36	128.32	120.71
1	B	244	PHE	C-N-CA	5.36	128.32	120.71
1	A	40	TYR	N-CA-C	-5.09	100.00	108.34
1	A	101	LEU	CA-CB-CG	5.04	133.93	116.30
1	B	35	LEU	CA-C-N	-5.01	113.58	119.84
1	B	35	LEU	C-N-CA	-5.01	113.58	119.84

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	154	TRP	Peptide
1	A	155	GLY	Peptide
1	B	154	TRP	Peptide

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	1916	0	1878	68	0
1	B	1935	0	1904	66	0
2	A	43	0	32	13	0
2	B	43	0	32	0	0
3	A	104	0	0	5	0
3	B	98	0	0	2	0
All	All	4139	0	3846	137	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 18.

All (137) 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)
2:A:1131:BLA:HHA	2:A:1131:BLA:HBD2	1.31	1.09
1:B:117:ILE:HD11	1:B:164:PHE:CE2	1.87	1.08
1:B:117:ILE:CG1	1:B:164:PHE:CD2	2.43	1.02
1:A:244:PHE:HZ	2:A:1131:BLA:CMC	1.76	0.96
1:B:117:ILE:HG12	1:B:164:PHE:CD2	2.04	0.92
1:B:117:ILE:HD11	1:B:164:PHE:CD2	2.08	0.88
1:B:71:ARG:HH12	1:B:206:ARG:NH1	1.74	0.84
1:A:33:GLN:HE21	1:A:33:GLN:HA	1.43	0.84
1:B:33:GLN:HG3	3:B:274:HOH:O	1.76	0.83
2:A:1131:BLA:HHA	2:A:1131:BLA:CBD	2.10	0.82
1:A:6:LEU:HG	1:A:65:TYR:CZ	2.15	0.81
1:B:117:ILE:CD1	1:B:164:PHE:CD2	2.64	0.79
2:A:1131:BLA:HBD2	2:A:1131:BLA:CHA	2.12	0.79
1:B:96:LEU:C	1:B:96:LEU:HD23	2.07	0.79
1:B:117:ILE:CG1	1:B:164:PHE:HD2	1.90	0.79
1:A:244:PHE:HZ	2:A:1131:BLA:HMC1	1.49	0.77
1:A:244:PHE:CZ	2:A:1131:BLA:CMC	2.67	0.75
1:A:71:ARG:NH1	1:A:209:GLN:OE1	2.20	0.74
1:A:105:ASP:OD2	2:A:1131:BLA:HAC	1.88	0.73
1:A:24:LEU:HD21	1:A:104[A]:CYS:SG	2.29	0.73
1:A:244:PHE:HZ	2:A:1131:BLA:HMC2	1.53	0.71
1:B:96:LEU:HD21	1:B:196:PRO:HB3	1.73	0.71
1:B:84:LEU:HD11	1:B:86:ILE:HD11	1.73	0.71
1:A:127:ARG:HG2	3:A:283:HOH:O	1.90	0.70
1:A:244:PHE:CZ	2:A:1131:BLA:HMC2	2.27	0.69
1:B:117:ILE:HG13	1:B:164:PHE:HD2	1.61	0.65
1:B:117:ILE:CD1	1:B:164:PHE:CE2	2.74	0.64
1:A:33:GLN:HA	1:A:33:GLN:NE2	2.10	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:VAL:HB	1:B:84:LEU:HD23	1.79	0.63
1:A:106:ILE:HD13	1:A:116:ALA:HA	1.81	0.63
1:A:244:PHE:CZ	2:A:1131:BLA:HMC1	2.34	0.63
1:A:6:LEU:HB3	1:A:33:GLN:HE22	1.64	0.62
1:A:79:LYS:HD2	1:A:85:ASP:OD1	2.00	0.62
1:A:123:THR:HG23	1:A:205:HIS:ND1	2.14	0.61
1:B:106:ILE:HD13	1:B:116:ALA:HA	1.81	0.61
1:A:6:LEU:HD23	1:A:7:SER:N	2.16	0.61
1:A:216:GLN:HG2	1:A:244:PHE:CZ	2.37	0.60
1:B:117:ILE:HA	1:B:163:LEU:O	2.01	0.59
1:B:237:ARG:O	1:B:241:GLN:HB2	2.01	0.59
1:A:144:GLU:O	1:A:175:ARG:NH2	2.34	0.59
1:B:169:ASN:ND2	1:B:171:THR:HB	2.19	0.58
1:A:33:GLN:HE21	1:A:33:GLN:CA	2.13	0.57
1:A:53:LEU:O	1:A:56:GLU:HG2	2.04	0.57
1:A:201:GLN:O	1:A:205:HIS:HD2	1.87	0.57
1:B:71:ARG:NH1	1:B:206:ARG:NH1	2.50	0.57
1:B:117:ILE:HG13	1:B:164:PHE:CD2	2.34	0.57
1:B:24:LEU:C	1:B:24:LEU:HD13	2.31	0.55
1:B:24:LEU:HD13	1:B:24:LEU:O	2.07	0.54
1:B:154:TRP:H	1:B:154:TRP:CD1	2.22	0.53
1:B:145:PHE:CD2	1:B:165:ILE:CD1	2.92	0.53
1:B:101:LEU:HD11	1:B:209:GLN:HG2	1.90	0.52
1:B:245:ASP:O	1:B:246:VAL:HB	2.09	0.52
1:B:148:GLN:HG3	3:B:297:HOH:O	2.11	0.50
1:B:169:ASN:ND2	1:B:172:GLU:H	2.08	0.50
1:A:6:LEU:HD12	1:A:37:LEU:HD11	1.94	0.49
1:B:101:LEU:HD12	1:B:123:THR:CG2	2.42	0.49
1:B:197:LEU:HD12	1:B:198:SER:HA	1.94	0.49
1:A:6:LEU:CD2	1:A:6:LEU:C	2.86	0.49
1:B:213:CYS:O	1:B:217:GLN:HG2	2.13	0.49
1:A:6:LEU:HD23	1:A:6:LEU:C	2.38	0.49
1:A:123:THR:HG23	1:A:205:HIS:CE1	2.48	0.49
1:A:154:TRP:CB	1:A:218:LYS:HD2	2.43	0.48
1:B:169:ASN:HD22	1:B:172:GLU:H	1.60	0.48
1:A:217:GLN:CA	1:A:217:GLN:HE21	2.26	0.48
1:B:163:LEU:HD11	1:B:179:ARG:HG2	1.95	0.48
1:A:154:TRP:HB3	1:A:218:LYS:HD2	1.94	0.48
1:A:126:ASP:HB3	1:A:128:GLN:H	1.79	0.48
1:B:6:LEU:C	1:B:6:LEU:HD12	2.39	0.47
1:B:96:LEU:HD21	1:B:196:PRO:CB	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:THR:CG2	1:A:205:HIS:ND1	2.77	0.47
1:A:106:ILE:CD1	1:A:116:ALA:CB	2.92	0.47
1:A:216:GLN:CG	2:A:1131:BLA:HMC2	2.44	0.47
1:B:217:GLN:HA	1:B:217:GLN:OE1	2.15	0.47
1:A:6:LEU:HD12	1:A:37:LEU:CD1	2.45	0.47
1:B:145:PHE:CD2	1:B:165:ILE:HD13	2.50	0.47
1:A:53:LEU:O	1:A:56:GLU:CG	2.62	0.47
1:A:71:ARG:HG2	1:A:93:PRO:O	2.14	0.47
1:A:105:ASP:O	1:A:106:ILE:HD13	2.15	0.47
1:B:58:LEU:HD13	1:B:80:VAL:HA	1.96	0.47
1:B:169:ASN:HD22	1:B:169:ASN:C	2.23	0.47
1:A:217:GLN:HA	1:A:217:GLN:NE2	2.30	0.46
1:B:146:GLU:HG3	1:B:147:GLN:HG3	1.97	0.46
1:B:245:ASP:OD1	1:B:246:VAL:HG23	2.16	0.46
1:A:217:GLN:HE21	1:A:217:GLN:HA	1.79	0.46
1:A:216:GLN:HG2	2:A:1131:BLA:HMC2	1.98	0.46
1:B:37:LEU:HD21	1:B:73:MET:HE1	1.98	0.46
1:A:71:ARG:NH2	1:A:245:ASP:OD2	2.48	0.46
1:A:106:ILE:CD1	1:A:116:ALA:HB1	2.45	0.46
1:B:6:LEU:O	1:B:6:LEU:CG	2.61	0.46
1:A:147:GLN:OE1	1:A:166[A]:ARG:HD3	2.15	0.45
1:B:117:ILE:HG12	1:B:164:PHE:CG	2.49	0.45
1:B:6:LEU:O	1:B:6:LEU:HG	2.15	0.45
1:B:87:LEU:C	1:B:87:LEU:HD23	2.42	0.45
1:B:148:GLN:HA	1:B:165:ILE:HG22	1.99	0.45
1:A:93:PRO:HD2	1:A:101:LEU:HD13	2.00	0.45
1:B:58:LEU:HD12	1:B:59:VAL:N	2.32	0.45
1:B:239:MET:SD	1:B:243:LEU:HD12	2.57	0.44
1:A:51:GLY:O	1:A:52:ARG:HD2	2.18	0.44
1:B:35:LEU:HD11	1:B:188:CYS:SG	2.57	0.44
1:A:222:THR:HG23	2:A:1131:BLA:C3B	2.47	0.44
1:B:94:GLU:OE2	1:B:95:PRO:HD2	2.18	0.44
1:B:53:LEU:O	1:B:54:GLU:C	2.61	0.44
1:B:197:LEU:HG	1:B:198:SER:OG	2.18	0.44
1:A:58:LEU:HD13	1:A:80:VAL:HA	2.00	0.43
1:A:57:LYS:HE2	1:A:57:LYS:HB3	1.82	0.43
1:A:122:PRO:HD3	1:A:129:LEU:HD11	2.00	0.43
1:B:145:PHE:CG	1:B:165:ILE:HD13	2.53	0.43
1:B:235:ALA:O	1:B:239:MET:HG2	2.19	0.43
1:B:67:THR:HB	1:B:68:PRO:HD2	1.99	0.43
1:A:6:LEU:HB3	1:A:33:GLN:NE2	2.31	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:137:LEU:HD13	1:B:186:ILE:HD13	2.01	0.43
1:B:206:ARG:O	1:B:210:ILE:HG13	2.19	0.42
1:A:80:VAL:HB	1:A:84:LEU:HD23	2.01	0.42
1:B:101:LEU:HD12	1:B:123:THR:HG22	2.01	0.42
1:B:245:ASP:C	1:B:246:VAL:HG23	2.44	0.42
1:B:154:TRP:HA	1:B:215:GLN:NE2	2.34	0.42
1:A:223:ARG:O	1:A:226:LEU:N	2.53	0.42
1:A:6:LEU:HG	1:A:65:TYR:CE2	2.55	0.41
1:B:147:GLN:HB2	1:B:166:ARG:O	2.20	0.41
1:A:123:THR:HG21	3:A:325:HOH:O	2.20	0.41
1:A:147:GLN:OE1	1:A:166[B]:ARG:CZ	2.68	0.41
1:A:201:GLN:NE2	3:A:304:HOH:O	2.53	0.41
1:A:211:HIS:HD2	3:A:296:HOH:O	2.03	0.41
1:A:201:GLN:O	1:A:205:HIS:CD2	2.72	0.41
1:B:246:VAL:O	1:B:246:VAL:HG12	2.20	0.41
1:A:38:LYS:HE3	1:A:38:LYS:HB2	1.92	0.41
1:B:62:ASN:HD21	1:B:76:GLU:HG3	1.85	0.41
1:B:67:THR:HB	1:B:68:PRO:CD	2.51	0.41
1:A:189:HIS:O	1:A:193:VAL:HG13	2.20	0.41
1:A:215:GLN:HA	1:A:215:GLN:NE2	2.36	0.41
1:A:217:GLN:CA	1:A:217:GLN:NE2	2.83	0.41
1:A:185:GLN:NE2	3:A:324:HOH:O	2.47	0.40
1:A:163:LEU:HD11	1:A:179:ARG:HG2	2.04	0.40
1:B:117:ILE:HG12	1:B:164:PHE:HD2	1.58	0.40
1:A:52:ARG:HD2	1:A:52:ARG:HA	1.95	0.40
1:A:154:TRP:HB3	1:A:218:LYS:CD	2.52	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	240/248 (97%)	233 (97%)	7 (3%)	0	100	100
1	B	242/248 (98%)	234 (97%)	6 (2%)	2 (1%)	16	8
All	All	482/496 (97%)	467 (97%)	13 (3%)	2 (0%)	30	22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	141	GLY
1	B	196	PRO

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	211/216 (98%)	187 (89%)	24 (11%)	5	2
1	B	212/216 (98%)	190 (90%)	22 (10%)	7	2
All	All	423/432 (98%)	377 (89%)	46 (11%)	6	2

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LEU
1	A	23	GLN
1	A	33	GLN
1	A	35	LEU
1	A	38	LYS
1	A	52	ARG
1	A	58	LEU
1	A	86	ILE
1	A	101	LEU
1	A	126	ASP
1	A	137	LEU
1	A	149	ARG
1	A	150	GLU
1	A	160	GLU

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Mol	Chain	Res	Type
1	A	166[A]	ARG
1	A	166[B]	ARG
1	A	192	ILE
1	A	193	VAL
1	A	197	LEU
1	A	199	GLU
1	A	203	LEU
1	A	206	ARG
1	A	216	GLN
1	A	217	GLN
1	B	6	LEU
1	B	7	SER
1	B	35	LEU
1	B	41	GLN
1	B	54	GLU
1	B	56	GLU
1	B	57	LYS
1	B	58	LEU
1	B	96	LEU
1	B	101	LEU
1	B	125	SER
1	B	137	LEU
1	B	146	GLU
1	B	148	GLN
1	B	169	ASN
1	B	195	GLU
1	B	197	LEU
1	B	198	SER
1	B	204	GLU
1	B	218	LYS
1	B	241	GLN
1	B	245	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	18	ASN
1	A	23	GLN
1	A	33	GLN
1	A	62	ASN
1	A	124	GLN
1	A	134	GLN

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Mol	Chain	Res	Type
1	A	211	HIS
1	A	215	GLN
1	A	216	GLN
1	A	217	GLN
1	B	33	GLN
1	B	62	ASN
1	B	124	GLN
1	B	134	GLN
1	B	169	ASN
1	B	214	GLN

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 ⓘ

2 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	1131	-	46,46,46	2.53	16 (34%)	66,67,67	2.69	28 (42%)
2	BLA	B	1130	-	46,46,46	2.23	16 (34%)	66,67,67	3.03	31 (46%)

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	1131	-	-	14/26/74/74	0/4/4/4
2	BLA	B	1130	-	-	4/26/74/74	0/4/4/4

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1130	BLA	C3C-C2C	5.59	1.48	1.37
2	A	1131	BLA	CHA-C4D	5.58	1.49	1.38
2	A	1131	BLA	C3B-C2B	5.47	1.48	1.37
2	B	1130	BLA	CHA-C4D	5.06	1.48	1.38
2	A	1131	BLA	C3D-C2D	5.05	1.47	1.36
2	A	1131	BLA	C3C-C2C	4.91	1.47	1.37
2	A	1131	BLA	CHB-C1B	4.73	1.49	1.37
2	A	1131	BLA	CHB-C4A	4.64	1.50	1.40
2	B	1130	BLA	C3B-C2B	4.62	1.46	1.37
2	B	1130	BLA	C3D-C2D	4.39	1.46	1.36
2	A	1131	BLA	CHD-C4C	4.21	1.47	1.37
2	A	1131	BLA	CHA-C1A	3.87	1.49	1.40
2	A	1131	BLA	C2A-C3A	3.78	1.48	1.38
2	B	1130	BLA	CHB-C1B	3.69	1.46	1.37
2	B	1130	BLA	CHA-C1A	3.62	1.48	1.40
2	A	1131	BLA	C4A-C3A	3.61	1.50	1.41
2	B	1130	BLA	CHD-C4C	3.60	1.46	1.37
2	B	1130	BLA	CHB-C4A	3.39	1.48	1.40
2	A	1131	BLA	OC-C1C	3.29	1.29	1.23
2	A	1131	BLA	CHD-C1D	3.21	1.48	1.40
2	A	1131	BLA	C1B-C2B	3.12	1.50	1.45
2	A	1131	BLA	C1A-C2A	3.10	1.49	1.42
2	B	1130	BLA	CHD-C1D	3.06	1.47	1.40
2	B	1130	BLA	OB-C4B	3.05	1.29	1.23
2	A	1131	BLA	C4D-C3D	3.00	1.50	1.45
2	A	1131	BLA	OB-C4B	2.82	1.28	1.23
2	B	1130	BLA	C4D-C3D	2.65	1.49	1.45
2	B	1130	BLA	C4C-NC	-2.60	1.33	1.37
2	B	1130	BLA	C2A-C3A	2.56	1.45	1.38
2	B	1130	BLA	C1A-C2A	2.43	1.48	1.42
2	B	1130	BLA	OC-C1C	2.31	1.27	1.23
2	B	1130	BLA	C4A-C3A	2.23	1.46	1.41

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1130	BLA	C2B-C1B-NB	6.86	116.97	106.97
2	A	1131	BLA	C3B-C4B-NB	6.63	114.46	106.13
2	B	1130	BLA	C3B-C2B-C1B	-6.46	100.72	107.92
2	B	1130	BLA	C3A-C4A-NA	6.26	115.97	107.43
2	B	1130	BLA	C3B-C4B-NB	6.25	113.98	106.13
2	A	1131	BLA	C2A-C1A-NA	5.73	114.75	107.57
2	B	1130	BLA	C1B-NB-C4B	-5.56	103.83	110.66
2	B	1130	BLA	C4C-NC-C1C	-5.30	104.15	110.66
2	A	1131	BLA	C4C-NC-C1C	-5.21	104.26	110.66
2	A	1131	BLA	C3B-C2B-C1B	-5.19	102.14	107.92
2	A	1131	BLA	CMB-C2B-C1B	5.17	130.45	124.16
2	B	1130	BLA	CMC-C2C-C1C	5.07	131.81	121.21
2	B	1130	BLA	CHB-C4A-C3A	-5.04	115.18	127.53
2	B	1130	BLA	C2A-C1A-NA	5.01	113.85	107.57
2	A	1131	BLA	CMC-C2C-C1C	5.01	131.69	121.21
2	A	1131	BLA	C2B-C1B-NB	4.99	114.24	106.97
2	B	1130	BLA	CHA-C1A-C2A	-4.97	116.54	127.22
2	B	1130	BLA	C2C-C1C-NC	4.80	117.30	106.55
2	A	1131	BLA	C1A-C2A-C3A	-4.47	102.42	107.62
2	A	1131	BLA	C3C-C4C-NC	4.43	113.86	106.64
2	A	1131	BLA	C3A-C4A-NA	4.37	113.40	107.43
2	B	1130	BLA	C3C-C4C-NC	4.36	113.75	106.64
2	A	1131	BLA	C1B-NB-C4B	-4.26	105.43	110.66
2	B	1130	BLA	CAD-C3D-C4D	4.25	132.48	125.02
2	A	1131	BLA	C1A-CHA-C4D	4.24	137.49	128.22
2	B	1130	BLA	C4A-NA-C1A	-4.08	102.62	109.78
2	B	1130	BLA	CMB-C2B-C1B	4.00	129.02	124.16
2	A	1131	BLA	C2C-C1C-NC	3.91	115.30	106.55
2	B	1130	BLA	C4A-C3A-C2A	-3.69	103.14	107.76
2	B	1130	BLA	C1D-C2D-C3D	-3.50	102.50	106.48
2	B	1130	BLA	C3D-C4D-ND	3.33	114.76	110.04
2	B	1130	BLA	CHB-C1B-NB	-3.33	118.96	126.06
2	A	1131	BLA	CAD-C3D-C4D	3.32	130.85	125.02
2	A	1131	BLA	CHA-C1A-C2A	-3.29	120.15	127.22
2	A	1131	BLA	C4B-C3B-C2B	-3.27	103.73	107.92
2	A	1131	BLA	C1D-C2D-C3D	-3.21	102.84	106.48
2	B	1130	BLA	C4D-C3D-C2D	-3.16	103.31	106.73
2	A	1131	BLA	C4A-C3A-C2A	-3.15	103.82	107.76
2	A	1131	BLA	OB-C4B-C3B	-3.14	122.14	129.71
2	A	1131	BLA	CHB-C1B-NB	-2.98	119.70	126.06
2	B	1130	BLA	C1A-C2A-C3A	-2.91	104.23	107.62
2	B	1130	BLA	C2D-C1D-ND	2.86	116.84	110.58

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1131	BLA	C4D-C3D-C2D	-2.70	103.81	106.73
2	A	1131	BLA	C3D-C4D-ND	2.68	113.83	110.04
2	B	1130	BLA	C4B-C3B-C2B	-2.67	104.50	107.92
2	B	1130	BLA	C1A-CHA-C4D	2.65	134.02	128.22
2	B	1130	BLA	OC-C1C-NC	-2.54	119.13	125.08
2	A	1131	BLA	CHD-C4C-C3C	-2.49	120.82	127.94
2	A	1131	BLA	CHD-C1D-C2D	-2.47	118.52	124.87
2	B	1130	BLA	O1A-CGA-CBA	-2.46	115.30	123.09
2	B	1130	BLA	C4C-CHD-C1D	-2.43	122.09	128.06
2	A	1131	BLA	C4A-NA-C1A	-2.38	105.61	109.78
2	B	1130	BLA	CAA-CBA-CGA	-2.31	107.53	113.67
2	A	1131	BLA	C2D-C1D-ND	2.27	115.55	110.58
2	A	1131	BLA	CAD-CBD-CGD	-2.20	107.83	113.67
2	B	1130	BLA	CMD-C2D-C1D	2.19	128.50	125.10
2	B	1130	BLA	C4D-ND-C1D	-2.15	102.57	106.52
2	B	1130	BLA	CHD-C1D-C2D	-2.14	119.39	124.87
2	A	1131	BLA	CHB-C4A-C3A	-2.10	122.39	127.53

There are no chirality outliers.

All (18) torsion outliers are listed below:

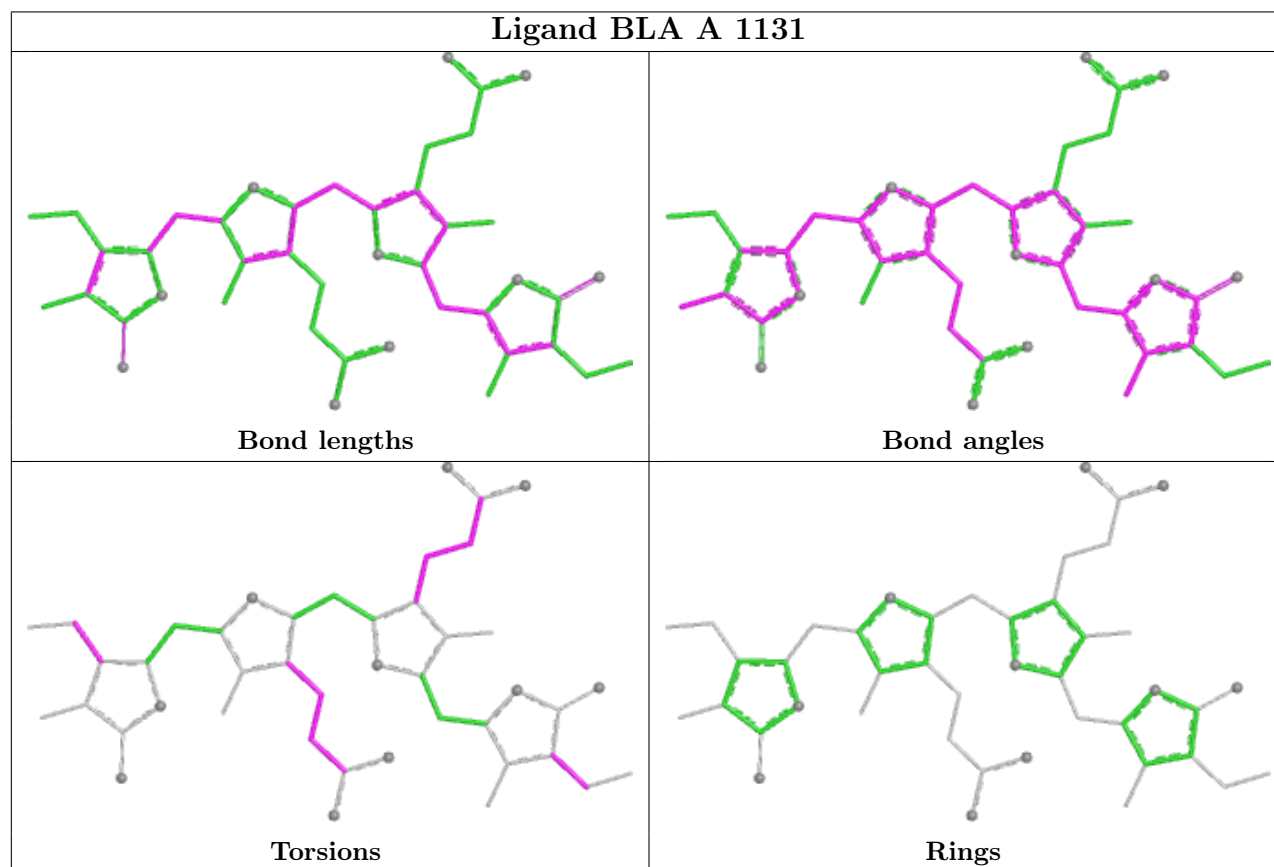
Mol	Chain	Res	Type	Atoms
2	A	1131	BLA	C1A-C2A-CAA-CBA
2	A	1131	BLA	C3A-C2A-CAA-CBA
2	A	1131	BLA	C2C-C3C-CAC-CBC
2	A	1131	BLA	C2D-C3D-CAD-CBD
2	A	1131	BLA	C4D-C3D-CAD-CBD
2	A	1131	BLA	C2A-CAA-CBA-CGA
2	A	1131	BLA	C3D-CAD-CBD-CGD
2	A	1131	BLA	C2B-C3B-CAB-CBB
2	A	1131	BLA	C4C-C3C-CAC-CBC
2	A	1131	BLA	C4B-C3B-CAB-CBB
2	A	1131	BLA	CAA-CBA-CGA-O1A
2	A	1131	BLA	CAD-CBD-CGD-O1D
2	B	1130	BLA	CAD-CBD-CGD-O2D
2	B	1130	BLA	CAD-CBD-CGD-O1D
2	A	1131	BLA	CAA-CBA-CGA-O2A
2	A	1131	BLA	CAD-CBD-CGD-O2D
2	B	1130	BLA	CAA-CBA-CGA-O2A
2	B	1130	BLA	CAA-CBA-CGA-O1A

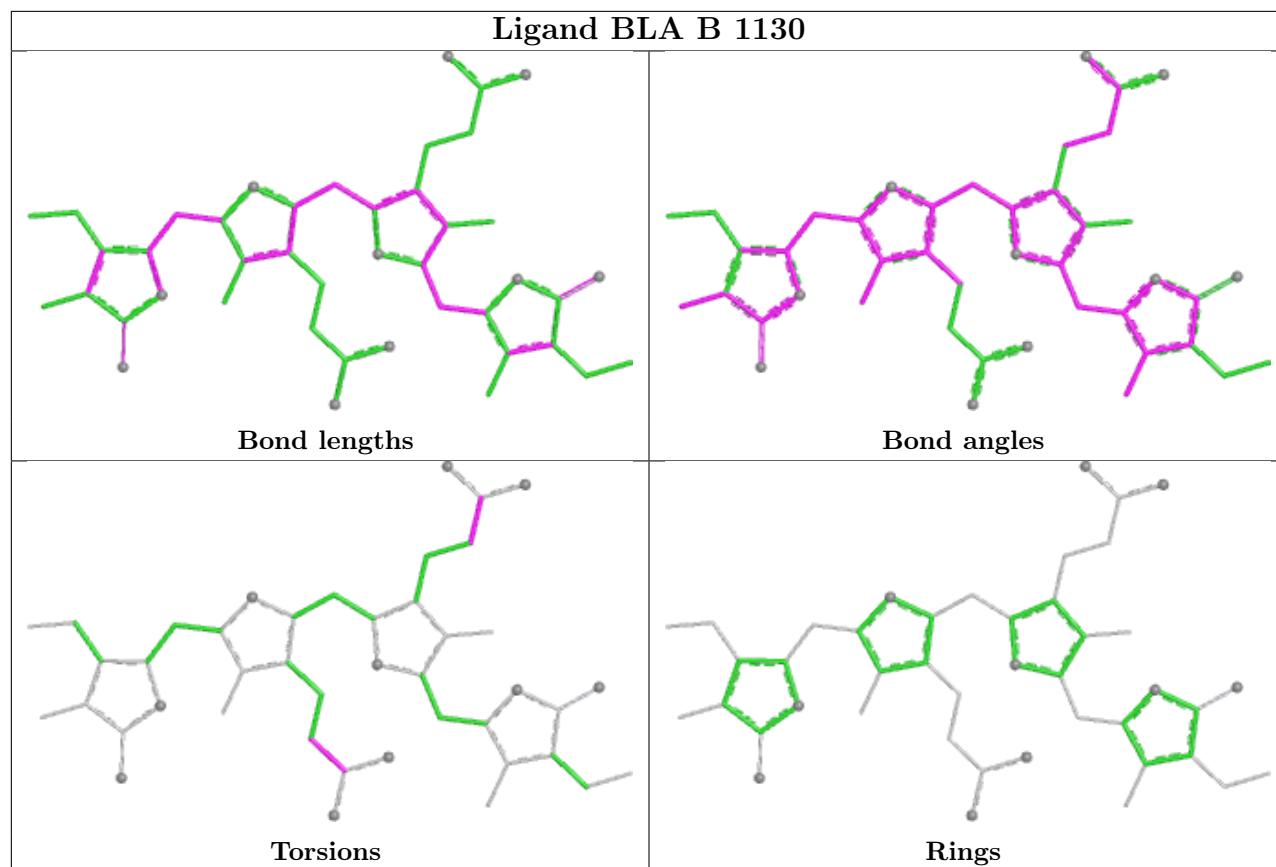
There are no ring outliers.

1 monomer is involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1131	BLA	13	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	238/248 (95%)	-0.84	0 100 100	18, 34, 50, 58	7 (2%)
1	B	241/248 (97%)	-0.78	0 100 100	20, 35, 50, 58	3 (1%)
All	All	479/496 (96%)	-0.81	0 100 100	18, 35, 50, 58	10 (2%)

There are no RSRZ outliers to report.

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](#)

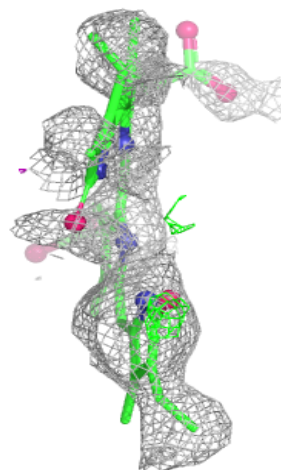
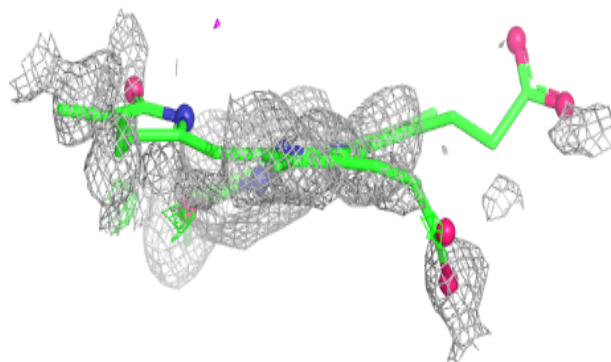
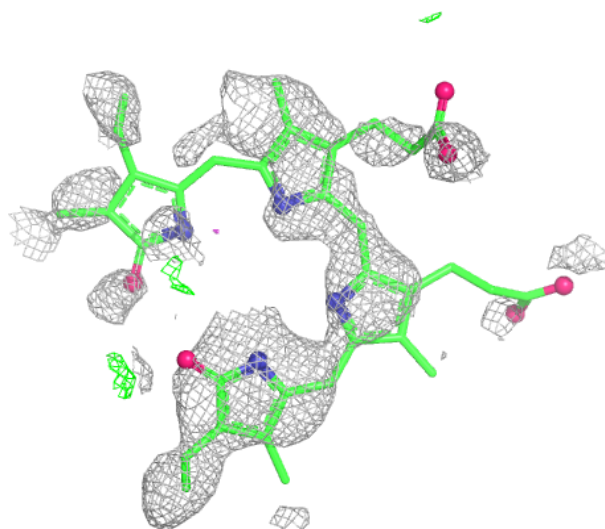
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists 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	BLA	A	1131	43/43	0.97	0.07	63,70,75,76	0
2	BLA	B	1130	43/43	0.98	0.05	33,43,48,50	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.

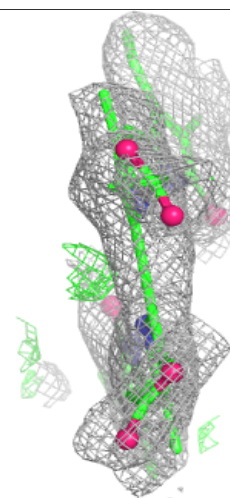
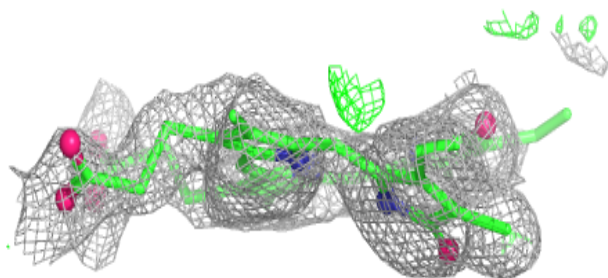
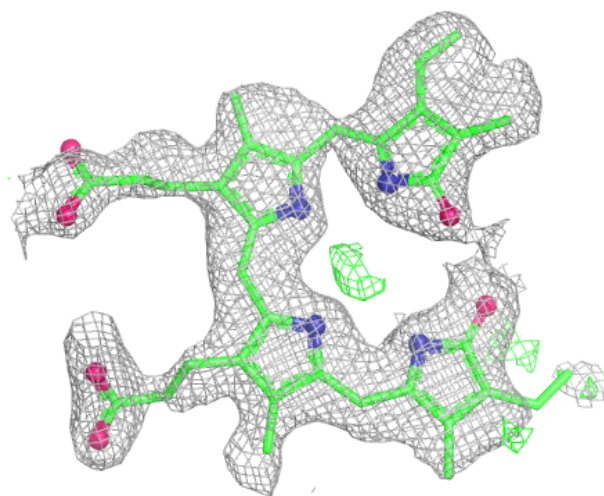
Electron density around BLA A 1131:

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



Electron density around BLA B 1130:

$2mF_o-DF_c$ (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.