



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:33 PM GMT

PDB ID : 1BBP  
Title : MOLECULAR STRUCTURE OF THE BILIN BINDING PROTEIN (BBP) FROM PIERIS BRASSICAE AFTER REFINEMENT AT 2.0 ANGSTROMS RESOLUTION.  
Authors : Huber, R.; Schneider, M.; Mayr, I.; Mueller, R.; Deutzmann, R.; Suter, F.; Zuber, H.; Falk, H.; Kayser, H.  
Deposited on : 1990-09-19  
Resolution : 2.00 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](mailto: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 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

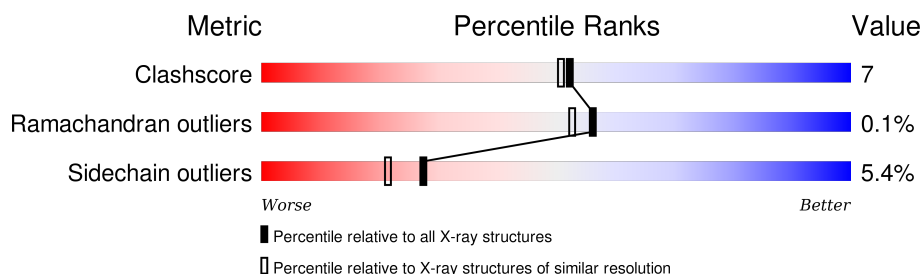
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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  $\geq 3$ , 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	173	
1	B	173	
1	C	173	
1	D	173	

2 Entry composition ⓘ

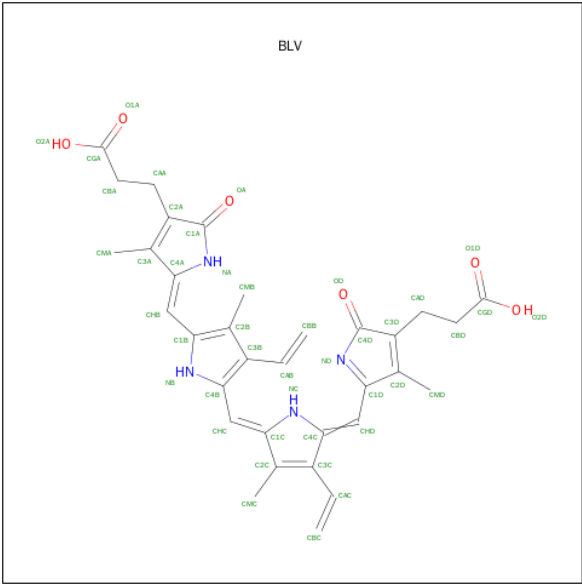
There are 3 unique types of molecules in this entry. The entry contains 6169 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 BILIN BINDING PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	173	Total	C	N	O	S	0	0	0
			1394	898	226	266	4			
1	B	173	Total	C	N	O	S	0	0	0
			1392	897	226	265	4			
1	C	173	Total	C	N	O	S	0	0	0
			1394	898	226	266	4			
1	D	173	Total	C	N	O	S	0	0	0
			1393	898	226	265	4			

- Molecule 2 is BILIVERDIN IX GAMMA CHROMOPHORE (three-letter code: BLV) (formula: C<sub>33</sub>H<sub>34</sub>N<sub>4</sub>O<sub>6</sub>).



*Continued from previous page...*

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

- Molecule 3 is water.

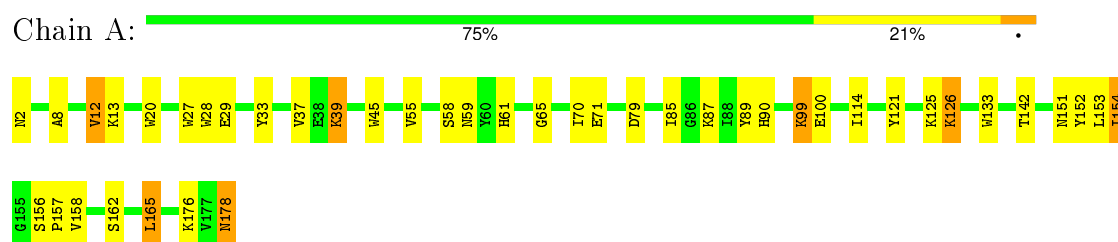
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	97	Total	O	0	0
			97	97		
3	B	100	Total	O	0	0
			100	100		
3	C	112	Total	O	0	0
			112	112		
3	D	115	Total	O	0	0
			115	115		

### 3 Residue-property plots

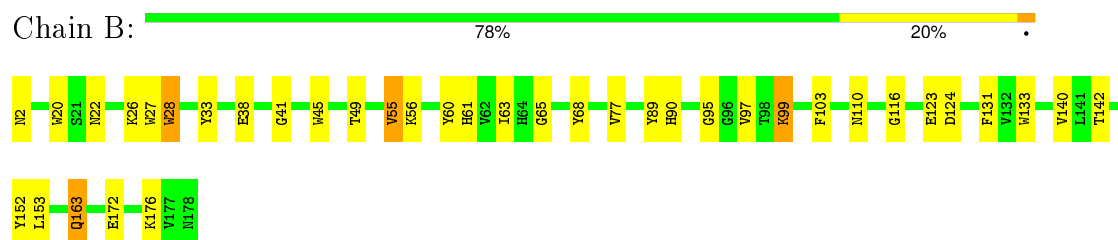
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

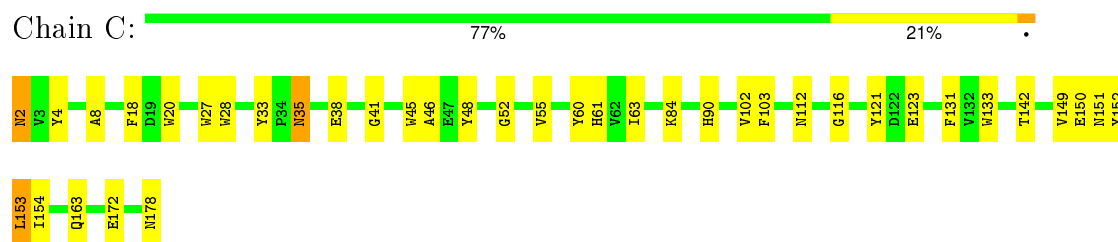
#### • Molecule 1: BILIN BINDING PROTEIN



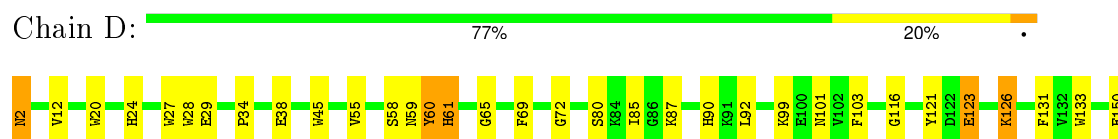
#### • Molecule 1: BILIN BINDING PROTEIN



#### • Molecule 1: BILIN BINDING PROTEIN



#### • Molecule 1: BILIN BINDING PROTEIN



H151	Y152	S162	Y166	K176	H178
L153	Q163			V177	
I154					

## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	132.10 Å   121.90 Å   63.80 Å 90.00°   90.00°   90.00°	Depositor
Resolution (Å)	(Not available) – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.00)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	EREF	Depositor
R, $R_{free}$	0.200 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	6169	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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

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	1.08	10/1436 (0.7%)	1.24	10/1947 (0.5%)
1	B	1.07	9/1434 (0.6%)	1.20	8/1944 (0.4%)
1	C	1.08	10/1436 (0.7%)	1.17	5/1947 (0.3%)
1	D	1.08	10/1435 (0.7%)	1.23	8/1946 (0.4%)
All	All	1.08	39/5741 (0.7%)	1.21	31/7784 (0.4%)

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	2
1	C	0	2
1	D	0	2
All	All	0	8

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	90	HIS	CB-CG	-15.19	1.22	1.50
1	C	90	HIS	CB-CG	-14.42	1.24	1.50
1	A	126	LYS	CD-CE	14.01	1.86	1.51
1	B	90	HIS	CB-CG	-13.86	1.25	1.50
1	B	99	LYS	CG-CD	10.85	1.89	1.52
1	A	99	LYS	CG-CD	9.08	1.83	1.52
1	D	90	HIS	CG-CD2	-8.93	1.20	1.35
1	B	90	HIS	CG-CD2	-8.28	1.21	1.35
1	C	90	HIS	CG-CD2	-8.02	1.22	1.35
1	D	27	TRP	NE1-CE2	-7.98	1.27	1.37

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	45	TRP	NE1-CE2	-7.71	1.27	1.37
1	B	27	TRP	NE1-CE2	-7.71	1.27	1.37
1	A	28	TRP	NE1-CE2	-7.64	1.27	1.37
1	A	20	TRP	NE1-CE2	-7.51	1.27	1.37
1	D	45	TRP	NE1-CE2	-7.45	1.27	1.37
1	D	28	TRP	NE1-CE2	-7.44	1.27	1.37
1	C	45	TRP	NE1-CE2	-7.36	1.27	1.37
1	C	20	TRP	NE1-CE2	-7.36	1.27	1.37
1	C	27	TRP	NE1-CE2	-7.35	1.27	1.37
1	A	27	TRP	NE1-CE2	-7.34	1.28	1.37
1	D	20	TRP	NE1-CE2	-7.23	1.28	1.37
1	A	133	TRP	NE1-CE2	-7.17	1.28	1.37
1	A	45	TRP	NE1-CE2	-7.10	1.28	1.37
1	B	20	TRP	NE1-CE2	-7.07	1.28	1.37
1	C	28	TRP	NE1-CE2	-7.07	1.28	1.37
1	C	123	GLU	CG-CD	7.06	1.62	1.51
1	A	90	HIS	CB-CG	-7.02	1.37	1.50
1	C	133	TRP	NE1-CE2	-7.01	1.28	1.37
1	C	123	GLU	CD-OE2	7.00	1.33	1.25
1	B	28	TRP	NE1-CE2	-6.98	1.28	1.37
1	B	133	TRP	NE1-CE2	-6.92	1.28	1.37
1	D	133	TRP	NE1-CE2	-6.88	1.28	1.37
1	B	124	ASP	CA-CB	5.91	1.67	1.53
1	D	101	ASN	N-CA	5.46	1.57	1.46
1	D	126	LYS	CE-NZ	5.42	1.62	1.49
1	A	114	ILE	N-CA	5.32	1.56	1.46
1	D	131	PHE	CA-CB	-5.31	1.42	1.53
1	A	114	ILE	CB-CG1	5.15	1.68	1.54
1	C	131	PHE	CA-CB	-5.11	1.42	1.53

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	90	HIS	ND1-CG-CD2	10.42	123.39	108.80
1	C	90	HIS	ND1-CG-CD2	10.01	122.82	108.80
1	B	90	HIS	ND1-CG-CD2	9.41	121.97	108.80
1	A	126	LYS	CD-CE-NZ	-8.92	91.19	111.70
1	D	131	PHE	CB-CG-CD2	-8.02	115.18	120.80
1	A	114	ILE	CB-CA-C	7.86	127.31	111.60
1	C	131	PHE	CB-CG-CD2	-7.57	115.50	120.80
1	D	126	LYS	CD-CE-NZ	7.42	128.77	111.70
1	A	99	LYS	CD-CE-NZ	7.14	128.13	111.70

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	99	LYS	CD-CE-NZ	7.11	128.05	111.70
1	A	99	LYS	CG-CD-CE	-6.95	91.05	111.90
1	B	99	LYS	CG-CD-CE	-6.69	91.84	111.90
1	C	33	TYR	CB-CG-CD1	-6.61	117.03	121.00
1	A	33	TYR	CB-CG-CD1	-6.53	117.08	121.00
1	B	33	TYR	CB-CG-CD1	-6.42	117.15	121.00
1	C	123	GLU	OE1-CD-OE2	-6.08	116.01	123.30
1	A	114	ILE	CA-CB-CG1	6.06	122.51	111.00
1	B	131	PHE	CB-CG-CD2	-6.01	116.59	120.80
1	D	90	HIS	CG-CD2-NE2	-5.77	98.23	109.20
1	D	80	SER	CB-CA-C	5.68	120.89	110.10
1	C	90	HIS	CG-CD2-NE2	-5.62	98.52	109.20
1	B	99	LYS	CB-CG-CD	-5.56	97.14	111.60
1	B	55	VAL	CA-CB-CG1	5.54	119.22	110.90
1	A	29	GLU	N-CA-CB	5.48	120.46	110.60
1	A	99	LYS	CB-CG-CD	-5.47	97.39	111.60
1	D	85	ILE	CB-CA-C	5.46	122.53	111.60
1	B	90	HIS	CG-CD2-NE2	-5.42	98.90	109.20
1	A	39	LYS	N-CA-CB	-5.39	100.89	110.60
1	A	114	ILE	CG1-CB-CG2	-5.07	100.25	111.40
1	D	99	LYS	CB-CA-C	-5.06	100.27	110.40
1	D	87	LYS	N-CA-CB	5.00	119.61	110.60

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	152	TYR	Sidechain
1	A	87	LYS	Mainchain
1	B	152	TYR	Sidechain
1	B	60	TYR	Mainchain
1	C	152	TYR	Sidechain
1	C	60	TYR	Mainchain
1	D	152	TYR	Sidechain
1	D	60	TYR	Mainchain

## 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	1394	0	1322	26	0
1	B	1392	0	1317	21	1
1	C	1394	0	1322	19	0
1	D	1393	0	1319	12	1
2	A	43	0	32	2	0
2	B	43	0	32	2	0
2	C	43	0	32	1	0
2	D	43	0	32	1	0
3	A	97	0	0	1	0
3	B	100	0	0	2	0
3	C	112	0	0	2	0
3	D	115	0	0	0	0
All	All	6169	0	5408	75	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 (75) 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:99:LYS:CG	1:A:99:LYS:CD	1.83	1.54
1:A:126:LYS:CD	1:A:126:LYS:CE	1.86	1.50
1:B:99:LYS:CD	1:B:99:LYS:CG	1.89	1.48
1:A:99:LYS:CG	1:A:99:LYS:CE	2.42	0.97
1:B:99:LYS:CE	1:B:99:LYS:CG	2.48	0.92
1:A:126:LYS:NZ	1:A:126:LYS:CD	2.40	0.85
1:B:99:LYS:CB	1:B:99:LYS:CD	2.61	0.79
1:A:99:LYS:CB	1:A:99:LYS:CD	2.61	0.78
1:A:126:LYS:HE2	3:A:589:HOH:O	1.89	0.72
1:A:126:LYS:CG	1:A:126:LYS:CE	2.65	0.72
1:A:99:LYS:CG	1:A:99:LYS:HE3	2.18	0.72
1:B:95:GLY:O	1:B:97:VAL:HG23	1.98	0.64
1:A:154:ILE:HG22	1:B:140:VAL:HG23	1.80	0.62
1:C:172:GLU:H	1:C:172:GLU:CD	2.04	0.61
1:B:103:PHE:CZ	1:B:116:GLY:HA3	2.37	0.59
1:D:2:ASN:HD22	1:D:2:ASN:N	2.00	0.59
1:B:41:GLY:HA2	1:B:63:ILE:HG12	1.84	0.59
1:C:154:ILE:HG22	3:C:506:HOH:O	2.04	0.57
2:A:500:BLV:HHD	2:A:500:BLV:HBC1	1.86	0.56
1:B:99:LYS:HE2	1:B:99:LYS:CG	2.35	0.56

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:89:TYR:CD2	1:A:100:GLU:HB3	2.41	0.56
1:B:172:GLU:CD	1:B:172:GLU:H	2.09	0.55
1:C:2:ASN:N	1:C:2:ASN:HD22	2.04	0.54
1:B:123:GLU:HG3	3:B:531:HOH:O	2.08	0.53
1:A:176:LYS:HE3	1:A:178:ASN:OD1	2.09	0.53
1:C:103:PHE:CZ	1:C:116:GLY:HA3	2.43	0.53
1:C:8:ALA:HA	1:C:121:TYR:CE1	2.44	0.52
1:A:154:ILE:HG22	1:B:140:VAL:CG2	2.40	0.51
1:C:112:ASN:HB3	3:C:589:HOH:O	2.10	0.51
1:A:2:ASN:N	1:A:2:ASN:HD22	2.08	0.51
1:B:2:ASN:N	1:B:2:ASN:HD22	2.07	0.51
1:C:150:GLU:O	1:C:154:ILE:HG23	2.09	0.51
1:D:34:PRO:HG3	1:D:166:VAL:HG21	1.92	0.51
1:A:162:SER:HA	1:A:165:LEU:HD22	1.94	0.50
1:B:163:GLN:HE22	1:C:163:GLN:HE21	1.58	0.50
1:D:60:TYR:CD1	1:D:69:PHE:HB3	2.47	0.49
1:D:59:ASN:HB3	2:D:500:BLV:C1B	2.43	0.49
1:D:65:GLY:HA2	1:D:176:LYS:O	2.14	0.48
1:B:77:VAL:HA	1:B:89:TYR:CE2	2.48	0.48
2:A:500:BLV:CHD	2:A:500:BLV:HBC1	2.43	0.48
1:C:48:TYR:HB3	1:C:55:VAL:HG11	1.96	0.48
1:B:63:ILE:HB	1:B:68:TYR:HE1	1.79	0.47
1:B:65:GLY:HA2	1:B:176:LYS:O	2.14	0.47
1:C:46:ALA:HB2	2:C:500:BLV:CBB	2.43	0.47
2:B:500:BLV:CBB	2:B:500:BLV:HMB1	2.45	0.47
1:D:103:PHE:CZ	1:D:116:GLY:HA3	2.49	0.46
1:A:13:LYS:HG2	1:A:13:LYS:H	1.52	0.46
1:A:8:ALA:HA	1:A:121:TYR:CE1	2.51	0.46
2:B:500:BLV:HBB1	2:B:500:BLV:HMB1	1.98	0.46
1:B:26:LYS:HE2	1:B:28:TRP:CZ2	2.51	0.46
1:D:29:GLU:HG3	1:D:29:GLU:O	2.16	0.45
1:A:154:ILE:HG21	1:A:154:ILE:HD12	1.67	0.45
1:D:38:GLU:HA	1:D:61:HIS:HE1	1.82	0.44
1:D:58:SER:HB3	1:D:60:TYR:CE2	2.53	0.44
1:A:58:SER:HB3	1:A:71:GLU:HG2	2.00	0.44
1:A:39:LYS:HG3	1:C:4:TYR:CD1	2.53	0.44
1:C:149:VAL:HG12	1:C:153:LEU:HD22	2.00	0.44
1:A:59:ASN:HB2	1:A:70:ILE:HG22	2.00	0.43
1:A:125:LYS:HD2	1:C:63:ILE:CD1	2.49	0.43
1:A:12:VAL:HG22	1:A:158:VAL:HG21	1.99	0.43
1:B:38:GLU:HA	1:B:61:HIS:HE1	1.84	0.43

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:ASN:HA	1:A:151:ASN:HD22	1.58	0.43
1:C:151:ASN:HD22	1:C:151:ASN:HA	1.61	0.43
1:D:72:GLY:HA3	1:D:92:LEU:HD23	2.01	0.43
1:C:41:GLY:HA2	1:C:63:ILE:HG12	2.00	0.43
1:C:18:PHE:O	1:C:84:LYS:HA	2.19	0.43
1:A:125:LYS:HD2	1:C:63:ILE:HD13	2.01	0.42
1:A:156:SER:HA	1:A:157:PRO:HD3	1.82	0.42
1:A:65:GLY:HA2	1:A:176:LYS:O	2.19	0.42
1:D:121:TYR:CE2	1:D:123:GLU:HA	2.54	0.42
1:B:49:THR:HB	1:B:56:LYS:HB2	2.01	0.41
1:D:150:GLU:O	1:D:154:ILE:HG12	2.20	0.41
1:B:110:ASN:HA	3:B:539:HOH:O	2.21	0.41
1:B:163:GLN:HE22	1:C:163:GLN:NE2	2.18	0.40
1:C:35:ASN:HD21	1:C:38:GLU:HG3	1.87	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:B:49:THR:OG1	1:D:24:HIS:NE2[1_554]	2.08	0.12

## 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	171/173 (99%)	165 (96%)	6 (4%)	0	100	100
1	B	171/173 (99%)	167 (98%)	4 (2%)	0	100	100
1	C	171/173 (99%)	165 (96%)	5 (3%)	1 (1%)	30	22
1	D	171/173 (99%)	165 (96%)	6 (4%)	0	100	100
All	All	684/692 (99%)	662 (97%)	21 (3%)	1 (0%)	56	53

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	52	GLY

### 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	150/150 (100%)	139 (93%)	11 (7%)	17	11
1	B	149/150 (99%)	144 (97%)	5 (3%)	44	41
1	C	150/150 (100%)	143 (95%)	7 (5%)	32	27
1	D	149/150 (99%)	140 (94%)	9 (6%)	24	17
All	All	598/600 (100%)	566 (95%)	32 (5%)	27	21

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	VAL
1	A	37	VAL
1	A	55	VAL
1	A	61	HIS
1	A	79	ASP
1	A	85	ILE
1	A	142	THR
1	A	153	LEU
1	A	154	ILE
1	A	165	LEU
1	A	178	ASN
1	B	22	ASN
1	B	55	VAL
1	B	142	THR
1	B	153	LEU
1	B	163	GLN
1	C	2	ASN
1	C	35	ASN
1	C	61	HIS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	C	102	VAL
1	C	142	THR
1	C	153	LEU
1	C	178	ASN
1	D	2	ASN
1	D	12	VAL
1	D	55	VAL
1	D	61	HIS
1	D	123	GLU
1	D	126	LYS
1	D	153	LEU
1	D	162	SER
1	D	163	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	2	ASN
1	A	17	ASN
1	A	59	ASN
1	A	151	ASN
1	A	163	GLN
1	B	2	ASN
1	B	5	HIS
1	B	22	ASN
1	B	59	ASN
1	B	129	GLN
1	B	151	ASN
1	B	163	GLN
1	B	178	ASN
1	C	35	ASN
1	C	112	ASN
1	C	151	ASN
1	C	163	GLN
1	C	178	ASN
1	D	2	ASN
1	D	5	HIS
1	D	59	ASN
1	D	151	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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z  > 2$	Counts	RMSZ	# $ Z  > 2$
2	BLV	A	500	-	33,46,46	1.59	8 (24%)	30,67,67	2.44	11 (36%)
2	BLV	B	500	-	33,46,46	1.50	5 (15%)	30,67,67	2.54	8 (26%)
2	BLV	C	500	-	33,46,46	1.48	6 (18%)	30,67,67	2.63	12 (40%)
2	BLV	D	500	-	33,46,46	1.56	7 (21%)	30,67,67	2.52	8 (26%)

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	BLV	A	500	-	-	0/14/58/58	0/4/4/4
2	BLV	B	500	-	-	0/14/58/58	0/4/4/4
2	BLV	C	500	-	-	0/14/58/58	0/4/4/4
2	BLV	D	500	-	-	0/14/58/58	0/4/4/4



All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	500	BLV	C4D-ND	-3.54	1.30	1.37
2	A	500	BLV	C4D-ND	-3.06	1.31	1.37
2	B	500	BLV	C4D-ND	-2.94	1.31	1.37
2	C	500	BLV	C4D-ND	-2.89	1.32	1.37
2	A	500	BLV	C1D-C2D	-2.49	1.40	1.45
2	D	500	BLV	C1A-NA	-2.19	1.32	1.37
2	C	500	BLV	C1D-C2D	-2.09	1.41	1.45
2	A	500	BLV	C1A-NA	-2.07	1.33	1.37
2	A	500	BLV	C3C-C2C	2.00	1.43	1.40
2	D	500	BLV	C3B-C2B	2.15	1.43	1.40
2	B	500	BLV	C4A-NA	2.16	1.41	1.37
2	C	500	BLV	C3D-C2D	2.22	1.41	1.36
2	D	500	BLV	C3D-C2D	2.37	1.41	1.36
2	A	500	BLV	C3D-C2D	2.38	1.41	1.36
2	A	500	BLV	CHB-C4A	2.65	1.40	1.34
2	D	500	BLV	CHB-C4A	2.77	1.40	1.34
2	C	500	BLV	CHB-C4A	2.78	1.40	1.34
2	B	500	BLV	CHB-C4A	3.12	1.41	1.34
2	C	500	BLV	CBC-CAC	3.18	1.51	1.28
2	B	500	BLV	CBC-CAC	3.20	1.52	1.28
2	D	500	BLV	CBC-CAC	3.32	1.52	1.28
2	C	500	BLV	CBB-CAB	3.36	1.53	1.28
2	D	500	BLV	CBB-CAB	3.37	1.53	1.28
2	A	500	BLV	CBB-CAB	3.39	1.53	1.28
2	B	500	BLV	CBB-CAB	3.40	1.53	1.28
2	A	500	BLV	CBC-CAC	3.45	1.53	1.28

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	BLV	C3C-CAC-CBC	-6.49	113.03	126.32
2	D	500	BLV	C3B-CAB-CBB	-6.17	113.69	126.32
2	D	500	BLV	CAD-CBD-CGD	-6.07	101.62	112.75
2	C	500	BLV	C3B-CAB-CBB	-6.04	113.96	126.32
2	A	500	BLV	CAD-CBD-CGD	-5.91	101.91	112.75
2	B	500	BLV	C3B-CAB-CBB	-5.86	114.33	126.32
2	C	500	BLV	C3C-CAC-CBC	-5.66	114.74	126.32
2	A	500	BLV	C3B-CAB-CBB	-5.17	115.75	126.32
2	D	500	BLV	C1A-C2A-C3A	-5.10	105.14	108.05
2	D	500	BLV	C3C-CAC-CBC	-4.78	116.54	126.32
2	C	500	BLV	OA-C1A-C2A	-4.76	122.39	128.09
2	A	500	BLV	C1A-C2A-C3A	-4.75	105.34	108.05

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	BLV	C1A-C2A-C3A	-3.78	105.89	108.05
2	C	500	BLV	CAA-CBA-CGA	-3.73	105.92	112.75
2	C	500	BLV	C1A-C2A-C3A	-3.67	105.95	108.05
2	A	500	BLV	C3C-CAC-CBC	-3.57	119.02	126.32
2	A	500	BLV	OA-C1A-C2A	-3.40	124.02	128.09
2	C	500	BLV	C4A-NA-C1A	-2.74	106.73	110.73
2	C	500	BLV	CAD-CBD-CGD	-2.66	107.87	112.75
2	D	500	BLV	OA-C1A-C2A	-2.64	124.93	128.09
2	A	500	BLV	CAA-CBA-CGA	-2.56	108.05	112.75
2	B	500	BLV	CAD-CBD-CGD	-2.41	108.32	112.75
2	A	500	BLV	CHB-C4A-C3A	-2.11	122.57	126.94
2	A	500	BLV	C4A-NA-C1A	-2.06	107.72	110.73
2	D	500	BLV	C3A-C4A-NA	2.03	109.94	107.00
2	A	500	BLV	C3A-C4A-NA	2.19	110.17	107.00
2	B	500	BLV	C2A-C1A-NA	2.22	108.75	106.74
2	B	500	BLV	CMA-C3A-C4A	2.36	127.33	124.20
2	C	500	BLV	C3A-C4A-NA	2.36	110.43	107.00
2	B	500	BLV	CAD-C3D-C4D	2.72	127.40	122.53
2	D	500	BLV	C2A-C1A-NA	2.99	109.46	106.74
2	C	500	BLV	CMA-C3A-C4A	3.09	128.31	124.20
2	C	500	BLV	CAD-C3D-C4D	3.32	128.48	122.53
2	A	500	BLV	C2A-C1A-NA	3.42	109.84	106.74
2	C	500	BLV	C2A-C1A-NA	3.92	110.30	106.74
2	C	500	BLV	CAA-C2A-C1A	4.42	125.72	121.37
2	D	500	BLV	CAA-C2A-C1A	5.08	126.37	121.37
2	A	500	BLV	CAA-C2A-C1A	5.14	126.43	121.37
2	B	500	BLV	CAA-C2A-C1A	7.59	128.84	121.37

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	500	BLV	2	0
2	B	500	BLV	2	0
2	C	500	BLV	1	0
2	D	500	BLV	1	0

## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

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

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section will therefore be empty.