



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jun 24, 2024 – 04:37 PM EDT

PDB ID : 6GIX  
Title : Water-soluble Chlorophyll Protein (WSCP) from *Lepidium virginicum* (Mutation L91P) with Chlorophyll-b  
Authors : Palm, D.M.; Agostini, A.; Averagesch, V.; Girr, P.; Werwie, M.; Takahashi, S.; Satoh, H.; Jaenicke, E.; Paulsen, H.  
Deposited on : 2018-05-15  
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

<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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.37.1  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.37.1

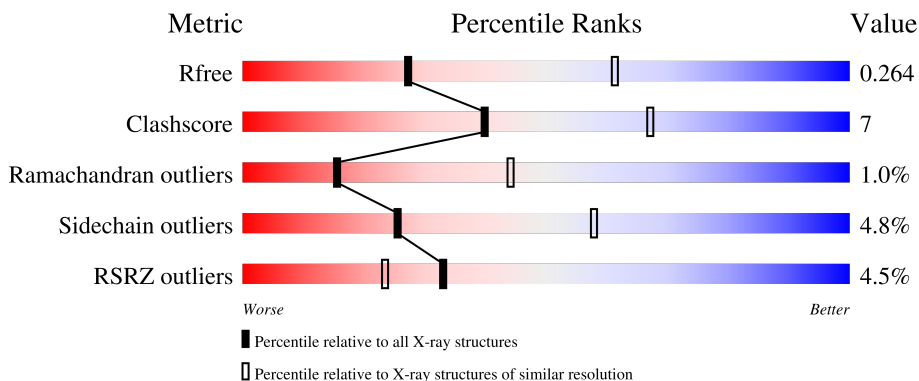
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
$R_{free}$	130704	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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  $\geq 3$ , 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	180	 3% (poor fit), 69% (0-1 outliers), 18% (2 outliers), 13% (3+ outliers or not modelled)
1	B	180	 4% (poor fit), 70% (0-1 outliers), 17% (2 outliers), 12% (3+ outliers or not modelled)
1	C	180	 6% (poor fit), 74% (0-1 outliers), 14% (2 outliers), 11% (3+ outliers or not modelled)
1	D	180	 3% (poor fit), 72% (0-1 outliers), 17% (2 outliers), 9% (3+ outliers or not modelled)

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 criteria:

<b>Mol</b>	<b>Type</b>	<b>Chain</b>	<b>Res</b>	<b>Chirality</b>	<b>Geometry</b>	<b>Clashes</b>	<b>Electron density</b>
2	CHL	A	1001	X	-	-	-
2	CHL	B	1001	X	-	-	-
2	CHL	C	1001	X	-	-	-
2	CHL	D	1001	X	-	-	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5231 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.

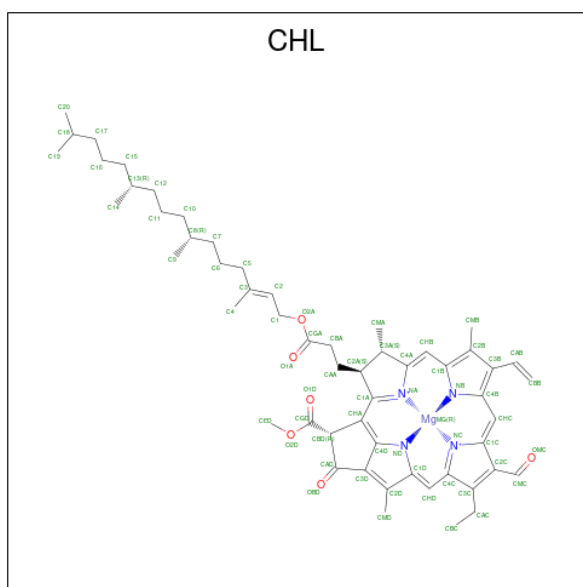
- Molecule 1 is a protein called Water-soluble chlorophyll protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	157	1209	775	199	233	2	0	0	0
1	B	158	1211	773	198	238	2	0	0	0
1	C	161	1242	792	204	244	2	0	0	0
1	D	163	1255	798	206	249	2	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	91	PRO	LEU	engineered mutation	UNP O04797
B	91	PRO	LEU	engineered mutation	UNP O04797
C	91	PRO	LEU	engineered mutation	UNP O04797
D	91	PRO	LEU	engineered mutation	UNP O04797

- Molecule 2 is CHLOROPHYLL B (three-letter code: CHL) (formula: C<sub>55</sub>H<sub>70</sub>MgN<sub>4</sub>O<sub>6</sub>).

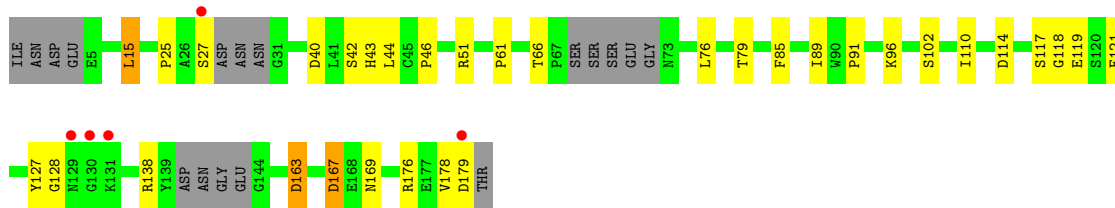


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	Mg	N			O
2	A	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
2	B	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
2	C	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
2	D	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	15	Total	O	0	0
			15	15		
3	B	10	Total	O	0	0
			10	10		
3	C	11	Total	O	0	0
			11	11		
3	D	14	Total	O	0	0
			14	14		





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	72.86Å 82.13Å 122.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.73 – 2.80 19.81 – 2.80	Depositor EDS
% Data completeness (in resolution range)	98.9 (19.73-2.80) 99.2 (19.81-2.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.28 (at 2.79Å)	Xtrriage
Refinement program	REFMAC 5.8.0158	Depositor
R, $R_{free}$	0.214 , 0.259 0.225 , 0.264	Depositor DCC
$R_{free}$ test set	925 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	44.0	Xtrriage
Anisotropy	0.075	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 39.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	5231	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.50% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.50	0/1238	0.75	1/1684 (0.1%)
1	B	0.49	0/1239	0.70	0/1686
1	C	0.49	0/1271	0.73	1/1729 (0.1%)
1	D	0.51	0/1285	0.79	3/1750 (0.2%)
All	All	0.50	0/5033	0.74	5/6849 (0.1%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	15	LEU	CA-CB-CG	6.64	130.56	115.30
1	C	15	LEU	CA-CB-CG	6.13	129.40	115.30
1	D	76	LEU	CB-CG-CD2	5.89	121.01	111.00
1	D	167	ASP	CB-CG-OD2	-5.73	113.14	118.30
1	A	66	THR	CA-CB-CG2	5.24	119.73	112.40

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	1209	0	1196	16	1
1	B	1211	0	1190	19	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1242	0	1216	16	1
1	D	1255	0	1224	12	0
2	A	66	0	70	7	0
2	B	66	0	70	10	0
2	C	66	0	70	5	0
2	D	66	0	70	5	0
3	A	15	0	0	1	0
3	B	10	0	0	0	0
3	C	11	0	0	0	0
3	D	14	0	0	0	0
All	All	5231	0	5106	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.

The worst 5 of 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:B:109:ILE:HD12	1:B:158:LEU:HB3	1.38	1.06
1:B:109:ILE:CD1	1:B:158:LEU:HB3	1.96	0.94
2:C:1001:CHL:H191	2:D:1001:CHL:H91	1.56	0.88
1:B:109:ILE:HD12	1:B:158:LEU:CB	2.03	0.87
2:B:1001:CHL:HBB2	1:C:44:LEU:HD11	1.58	0.85

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:129:ASN:ND2	1:C:106:GLU:O[4_547]	2.15	0.05

## 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	147/180 (82%)	143 (97%)	3 (2%)	1 (1%)	22	53
1	B	148/180 (82%)	144 (97%)	2 (1%)	2 (1%)	11	34
1	C	151/180 (84%)	147 (97%)	3 (2%)	1 (1%)	22	53
1	D	155/180 (86%)	150 (97%)	3 (2%)	2 (1%)	12	36
All	All	601/720 (84%)	584 (97%)	11 (2%)	6 (1%)	15	44

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	163	ASP
1	A	128	GLY
1	C	128	GLY
1	B	128	GLY
1	B	129	ASN

### 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	137/158 (87%)	133 (97%)	4 (3%)	42	76
1	B	138/158 (87%)	132 (96%)	6 (4%)	29	62
1	C	141/158 (89%)	134 (95%)	7 (5%)	24	56
1	D	143/158 (90%)	133 (93%)	10 (7%)	15	40
All	All	559/632 (88%)	532 (95%)	27 (5%)	25	58

5 of 27 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	129	ASN
1	D	15	LEU
1	D	169	ASN
1	C	163	ASP
1	D	27	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	11	ASN
1	C	13	ASN
1	D	73	ASN
1	C	43	HIS
1	B	43	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

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	CHL	C	1001	1	66,74,74	1.84	11 (16%)	73,114,114	2.11	20 (27%)
2	CHL	D	1001	1	66,74,74	1.77	13 (19%)	73,114,114	2.02	18 (24%)
2	CHL	B	1001	1	66,74,74	1.75	13 (19%)	73,114,114	2.06	23 (31%)
2	CHL	A	1001	1	66,74,74	1.76	14 (21%)	73,114,114	2.10	19 (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	CHL	C	1001	1	4/4/20/26	15/39/137/137	-
2	CHL	D	1001	1	3/3/20/26	13/39/137/137	-
2	CHL	B	1001	1	4/4/20/26	22/39/137/137	-
2	CHL	A	1001	1	4/4/20/26	19/39/137/137	-

The worst 5 of 51 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1001	CHL	O2D-CGD	5.74	1.47	1.33
2	C	1001	CHL	CHC-C1C	5.36	1.48	1.35
2	C	1001	CHL	CHD-C1D	5.28	1.48	1.38
2	D	1001	CHL	CHD-C1D	5.26	1.48	1.38
2	B	1001	CHL	CHD-C1D	5.14	1.48	1.38

The worst 5 of 80 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1001	CHL	CMD-C2D-C1D	7.33	137.64	124.71
2	B	1001	CHL	CMD-C2D-C1D	6.81	136.72	124.71
2	D	1001	CHL	CMD-C2D-C1D	6.79	136.68	124.71
2	A	1001	CHL	O2D-CGD-CBD	6.19	122.26	111.27
2	D	1001	CHL	O2D-CGD-CBD	6.06	122.03	111.27

5 of 15 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	1001	CHL	NA
2	A	1001	CHL	NC
2	A	1001	CHL	C8
2	A	1001	CHL	ND
2	B	1001	CHL	NA

5 of 69 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	CHL	CBD-CGD-O2D-CED
2	A	1001	CHL	C4-C3-C5-C6
2	B	1001	CHL	C2-C3-C5-C6

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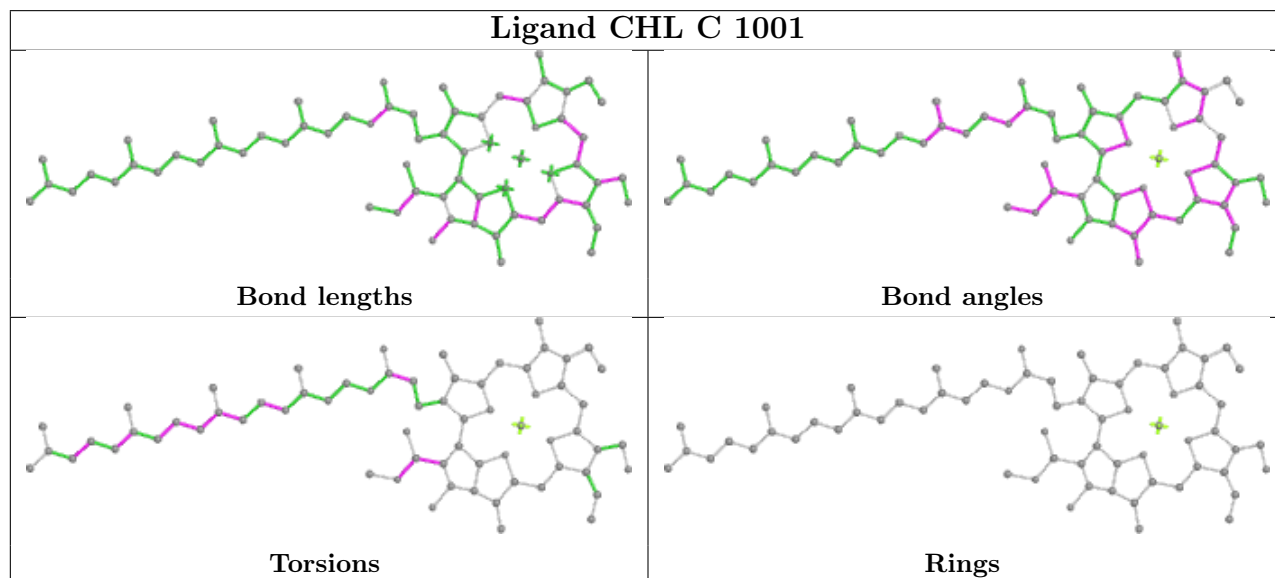
Mol	Chain	Res	Type	Atoms
2	B	1001	CHL	C4-C3-C5-C6
2	D	1001	CHL	CBD-CGD-O2D-CED

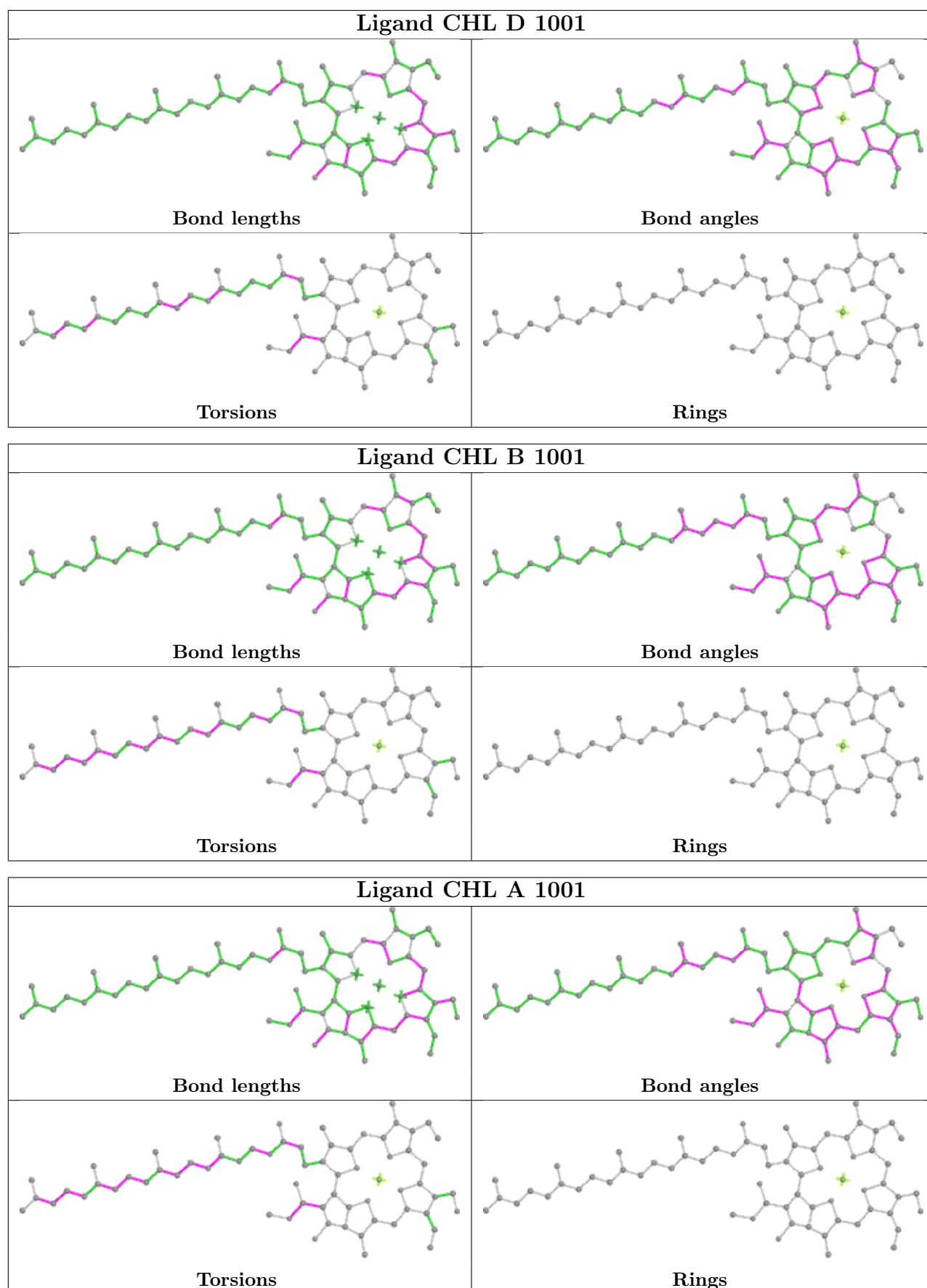
There are no ring outliers.

4 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1001	CHL	5	0
2	D	1001	CHL	5	0
2	B	1001	CHL	10	0
2	A	1001	CHL	7	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	157/180 (87%)	-0.01	5 (3%) 47 37	23, 37, 70, 90	0
1	B	158/180 (87%)	0.15	8 (5%) 28 19	26, 44, 83, 117	0
1	C	161/180 (89%)	0.06	11 (6%) 17 10	23, 39, 78, 104	0
1	D	163/180 (90%)	0.03	5 (3%) 49 39	21, 42, 71, 93	0
All	All	639/720 (88%)	0.05	29 (4%) 33 23	21, 41, 79, 117	0

The worst 5 of 29 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	129	ASN	5.2
1	B	163	ASP	5.1
1	D	179	ASP	3.9
1	C	129	ASN	3.5
1	B	105	GLU	3.3

### 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](#)

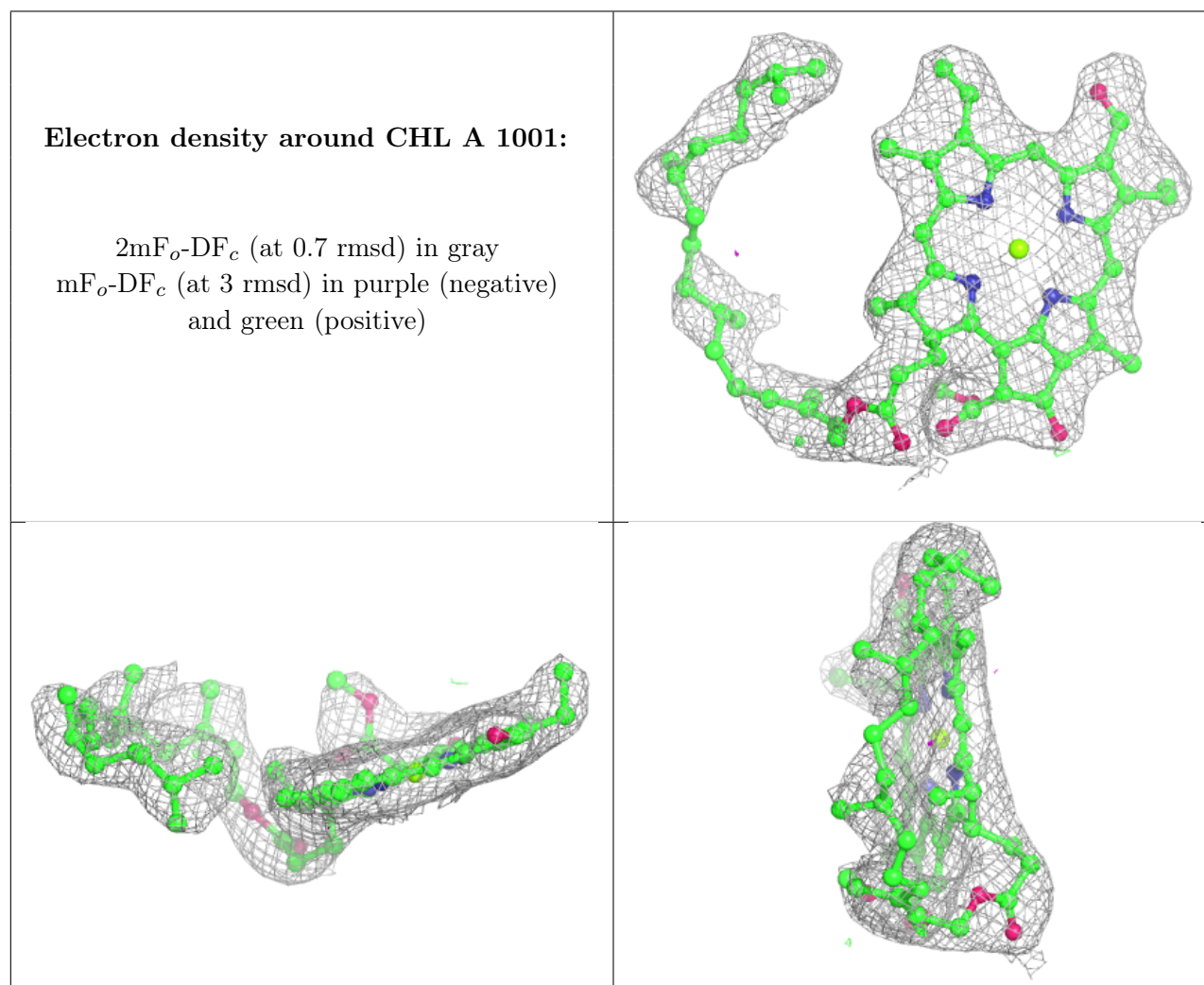
There are no monosaccharides in this entry.

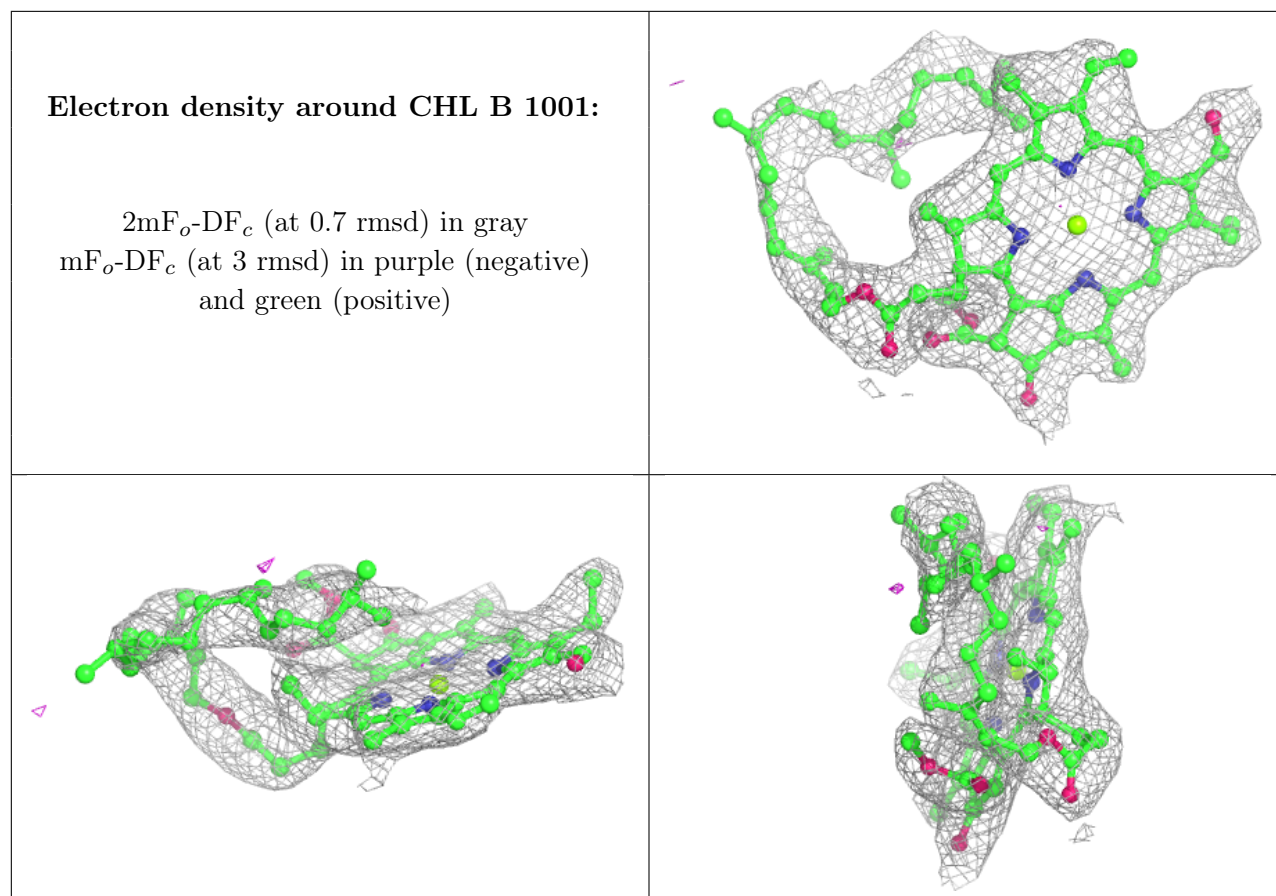
### 6.4 Ligands [i](#)

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, 95<sup>th</sup> 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
2	CHL	A	1001	66/66	0.94	0.19	24,29,71,74	0
2	CHL	B	1001	66/66	0.94	0.21	26,30,77,79	0
2	CHL	D	1001	66/66	0.94	0.18	21,24,54,59	0
2	CHL	C	1001	66/66	0.96	0.17	23,26,65,69	0

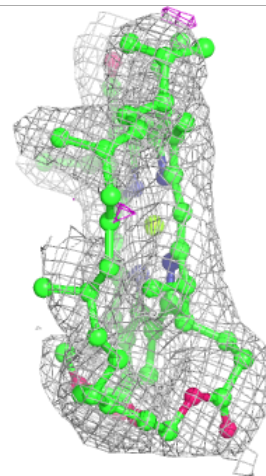
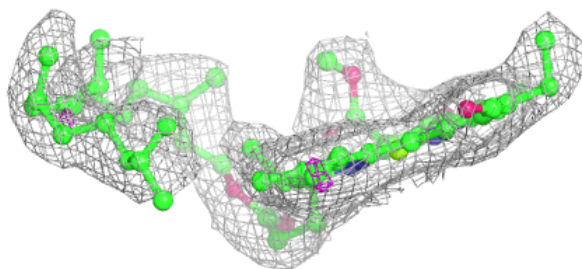
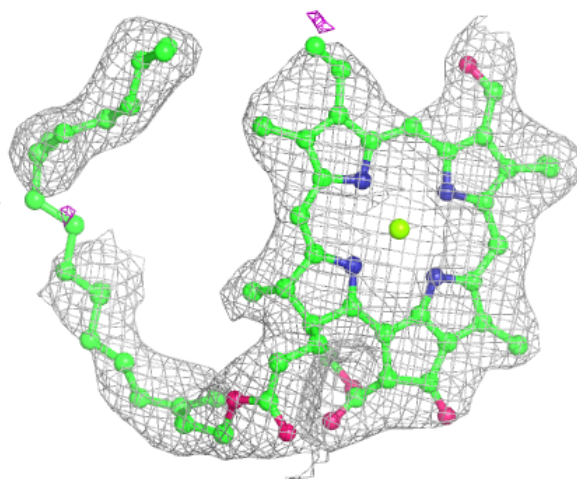
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

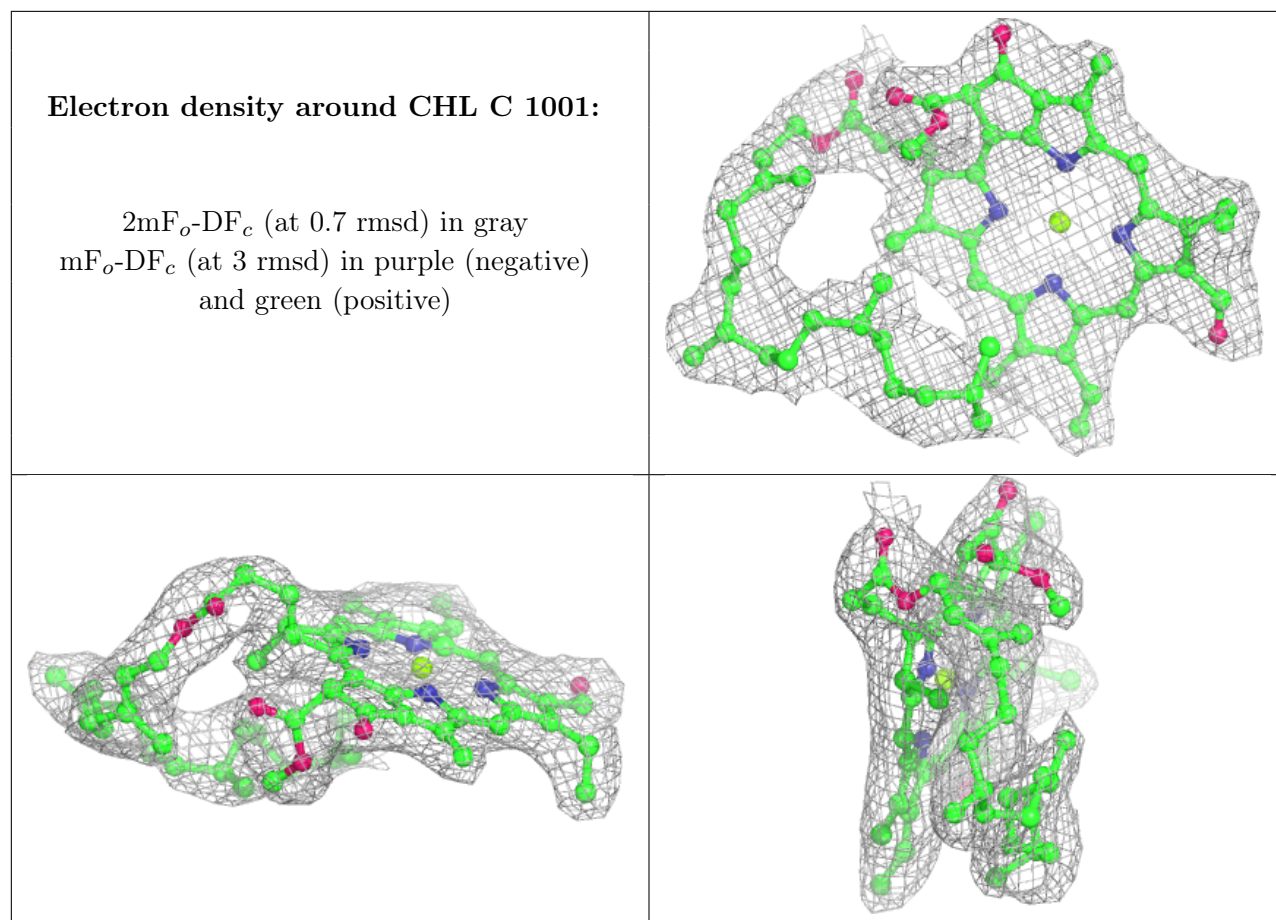




**Electron density around CHL D 1001:**

$2mF_o-DF_c$  (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.