



# Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 07:36 am BST

PDB ID : 3FE9  
Title : Crystal structure of a pheromone binding protein from *Apis mellifera* with a serendipitous ligand soaked at pH 7.0  
Authors : Pesenti, M.E.; Spinelli, S.; Bezirard, V.; Briand, L.; Pernollet, J.C.; Tegoni, M.; Cambillau, C.  
Deposited on : 2008-11-28  
Resolution : 1.75 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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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 : 2.11  
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) : 2.11

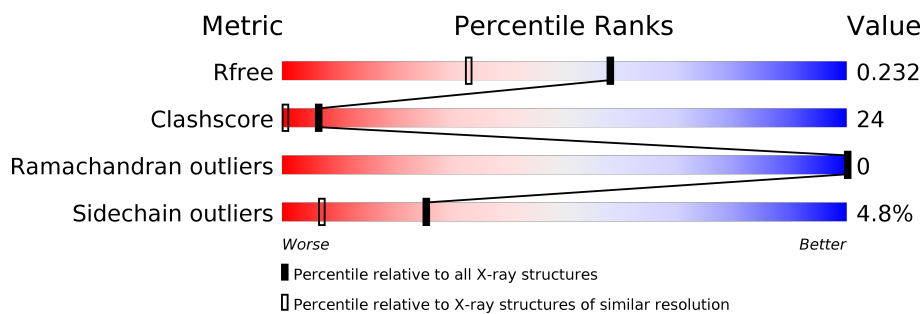
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)

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 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  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	119	

## 2 Entry composition [i](#)

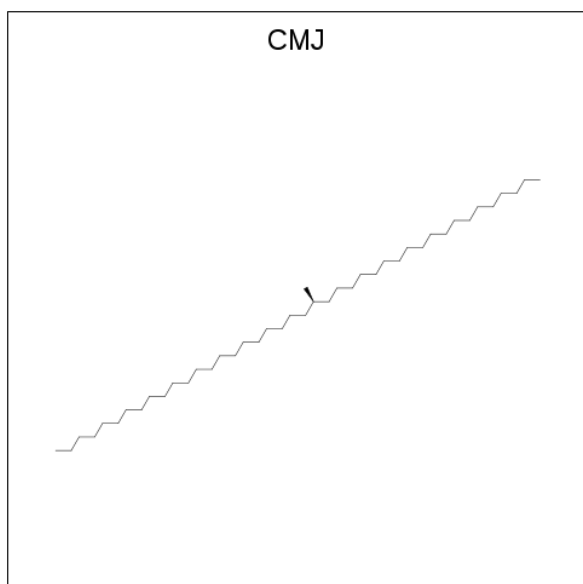
There are 4 unique types of molecules in this entry. The entry contains 1137 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Pheromone-binding protein ASP1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	117	Total	C	N	O	S	0	6	0
			934	586	145	192	11			

- Molecule 2 is (20S)-20-methyldotetracontane (three-letter code: CMJ) (formula: C<sub>43</sub>H<sub>88</sub>).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	C	0	0
			43	43		

- Molecule 3 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cl	0	0
			1	1		

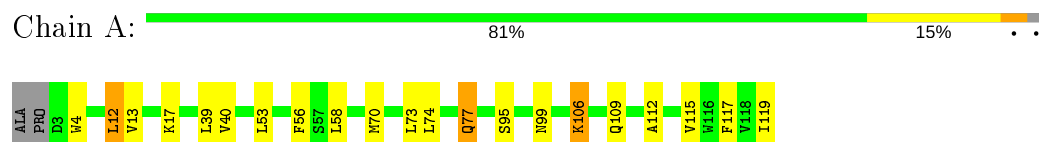
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	159	Total 159	O 159	0	0

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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. 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. 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: Pheromone-binding protein ASP1



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.30Å 83.91Å 46.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.75 29.56 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.00-1.75) 99.9 (29.56-1.75)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	0.09	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.42 (at 1.75Å)	Xtriage
Refinement program	REFMAC 5.4.0077	Depositor
R, $R_{free}$	0.161 , 0.187 0.204 , 0.232	Depositor DCC
$R_{free}$ test set	992 reflections (5.87%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.6	Xtriage
Anisotropy	0.069	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 35.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.031 for -k,-h,-l	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	1137	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.39% 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

### 5.1 Standard geometry

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

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	0.59	0/975	0.65	0/1324

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	934	0	898	39	0
2	A	43	0	88	12	0
3	A	1	0	0	0	0
4	A	159	0	0	2	0
All	All	1137	0	986	47	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 24.

All (47) 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:39[B]:LEU:HD11	1:A:106:LYS:CE	1.84	1.05
2:A:120:CMJ:H41A	2:A:120:CMJ:C1	1.98	0.93
1:A:39[B]:LEU:CD1	1:A:106:LYS:CD	2.51	0.88
2:A:120:CMJ:H1A	2:A:120:CMJ:C41	2.04	0.87
1:A:39[B]:LEU:CD1	1:A:106:LYS:HD2	2.05	0.86
1:A:39[B]:LEU:HD11	1:A:106:LYS:CD	2.04	0.86
1:A:39[B]:LEU:HD11	1:A:106:LYS:HE3	1.61	0.83
2:A:120:CMJ:H41A	2:A:120:CMJ:H1A	1.62	0.81
1:A:39[B]:LEU:CD1	1:A:106:LYS:CE	2.61	0.75
1:A:95:SER:H	1:A:99:ASN:HD22	1.33	0.75
1:A:39[B]:LEU:CD1	1:A:106:LYS:NZ	2.51	0.73
2:A:120:CMJ:H1A	2:A:120:CMJ:H41	1.70	0.72
1:A:13:VAL:HG21	1:A:119:ILE:HD13	1.71	0.71
1:A:39[B]:LEU:HD12	1:A:106:LYS:NZ	2.06	0.70
1:A:39[B]:LEU:HD13	1:A:106:LYS:HD2	1.78	0.64
2:A:120:CMJ:H36A	2:A:120:CMJ:H40	1.82	0.61
1:A:39[B]:LEU:CD1	1:A:106:LYS:HG2	2.33	0.58
1:A:39[B]:LEU:HD11	1:A:106:LYS:NZ	2.14	0.58
1:A:39[B]:LEU:HD12	1:A:106:LYS:HZ2	1.68	0.57
1:A:77:GLN:H	1:A:77:GLN:HE21	1.53	0.56
1:A:39[B]:LEU:HD11	1:A:106:LYS:CG	2.35	0.56
1:A:17[A]:LYS:HE3	1:A:119:ILE:OXT	2.06	0.56
1:A:77:GLN:H	1:A:77:GLN:NE2	2.04	0.54
1:A:39[B]:LEU:CD1	1:A:106:LYS:CG	2.86	0.54
1:A:4:TRP:CD1	1:A:115[A]:VAL:HG23	2.48	0.49
1:A:39[B]:LEU:HD13	1:A:106:LYS:CD	2.38	0.47
1:A:39[B]:LEU:HD11	1:A:106:LYS:HG2	1.96	0.47
1:A:13:VAL:HG21	1:A:119:ILE:CD1	2.41	0.46
1:A:56:PHE:HB3	2:A:120:CMJ:H21	1.98	0.46
1:A:39[B]:LEU:HD13	1:A:106:LYS:HG2	1.97	0.45
1:A:58:LEU:HD22	1:A:70[B]:MET:HE2	1.98	0.45
1:A:40:VAL:HG22	4:A:413:HOH:O	2.17	0.45
1:A:39[B]:LEU:CD1	1:A:106:LYS:HZ1	2.28	0.44
1:A:4:TRP:NE1	1:A:115[A]:VAL:CG2	2.82	0.43
1:A:12:LEU:HA	1:A:12:LEU:HD12	1.79	0.43
2:A:120:CMJ:H35	2:A:120:CMJ:H32	1.73	0.42
1:A:73:LEU:HD13	2:A:120:CMJ:H21A	2.01	0.42
1:A:53:LEU:HD11	2:A:120:CMJ:H9	2.01	0.42
2:A:120:CMJ:H41A	2:A:120:CMJ:H1	1.95	0.42
1:A:70[A]:MET:HE2	1:A:74:LEU:HG	2.01	0.42
1:A:73:LEU:HD22	2:A:120:CMJ:H23	2.02	0.41
1:A:112:ALA:HB1	1:A:115[A]:VAL:HG12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:120:CMJ:H1B	2:A:120:CMJ:H4A	1.64	0.41
1:A:95:SER:N	1:A:99:ASN:HD22	2.10	0.41
1:A:4:TRP:NE1	1:A:115[A]:VAL:HG23	2.35	0.41
1:A:39[B]:LEU:HG	4:A:162:HOH:O	2.20	0.41

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	122/119 (102%)	121 (99%)	1 (1%)	0	100	100

There are no Ramachandran outliers to report.

### 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	111/105 (106%)	106 (96%)	5 (4%)	27	8

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	77	GLN

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Mol	Chain	Res	Type
1	A	106	LYS
1	A	109	GLN
1	A	117	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	GLN
1	A	83	GLN
1	A	99	ASN
1	A	103	ASN
1	A	109	GLN

### 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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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	CMJ	A	120	-	42,42,42	0.19	0	42,42,42	0.81	1 (2%)

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	CMJ	A	120	-	-	21/41/41/41	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	120	CMJ	C23-C22-C20	-2.10	109.12	115.92

There are no chirality outliers.

All (21) torsion outliers are listed below:

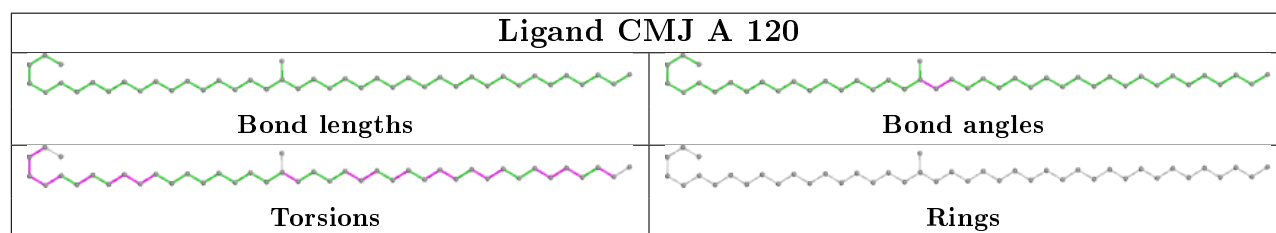
Mol	Chain	Res	Type	Atoms
2	A	120	CMJ	C21-C20-C22-C23
2	A	120	CMJ	C32-C33-C34-C35
2	A	120	CMJ	C9-C10-C11-C12
2	A	120	CMJ	C38-C39-C40-C41
2	A	120	CMJ	C4-C5-C6-C7
2	A	120	CMJ	C36-C37-C38-C39
2	A	120	CMJ	C25-C26-C27-C28
2	A	120	CMJ	C33-C34-C35-C36
2	A	120	CMJ	C11-C10-C9-C8
2	A	120	CMJ	C30-C31-C32-C33
2	A	120	CMJ	C1-C2-C3-C4
2	A	120	CMJ	C3-C4-C5-C6
2	A	120	CMJ	C27-C28-C29-C30
2	A	120	CMJ	C29-C30-C31-C32
2	A	120	CMJ	C24-C25-C26-C27
2	A	120	CMJ	C19-C20-C22-C23
2	A	120	CMJ	C37-C38-C39-C40
2	A	120	CMJ	C40-C41-C42-C43
2	A	120	CMJ	C2-C3-C4-C5
2	A	120	CMJ	C6-C7-C8-C9
2	A	120	CMJ	C10-C11-C12-C13

There are no ring outliers.

1 monomer is involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	120	CMJ	12	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

### 6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

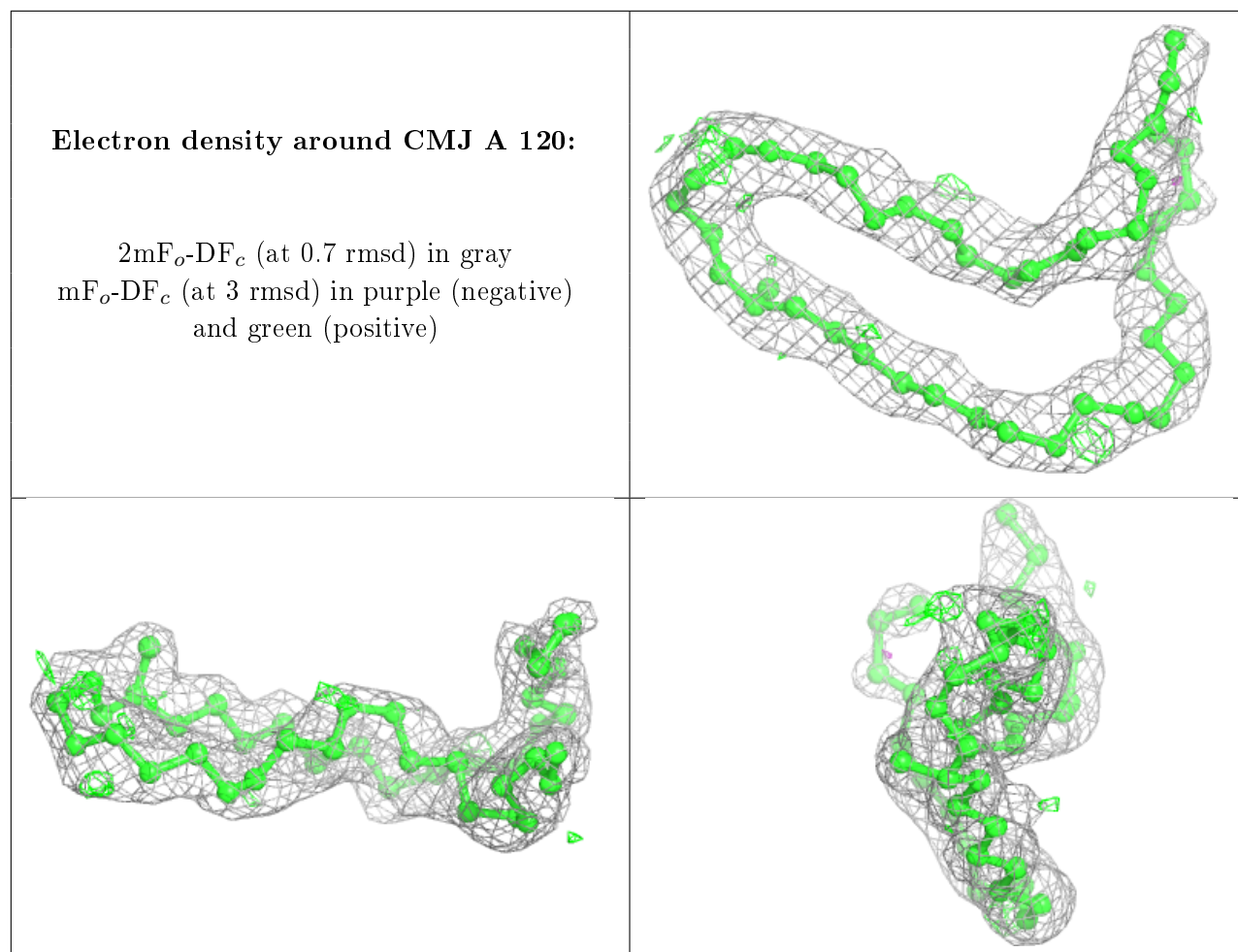
### 6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

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.



## 6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.