



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:36 AM UTC

PDB ID : 5F83 / pdb_00005f83
Title : Imipenem complex of the GES-5 C69G mutant
Authors : Smith, C.A.; Vakulenko, S.B.
Deposited on : 2015-12-09
Resolution : 1.38 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

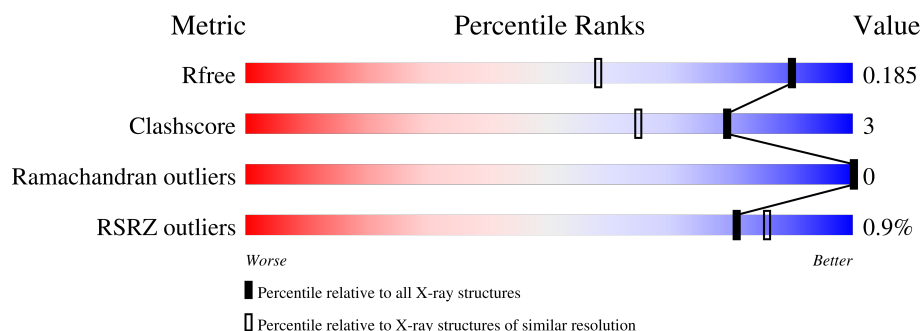
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.38 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	4403 (1.40-1.36)
Clashscore	190562	4528 (1.40-1.36)
Ramachandran outliers	187476	4459 (1.40-1.36)
RSRZ outliers	180081	4399 (1.40-1.36)

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4506 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.

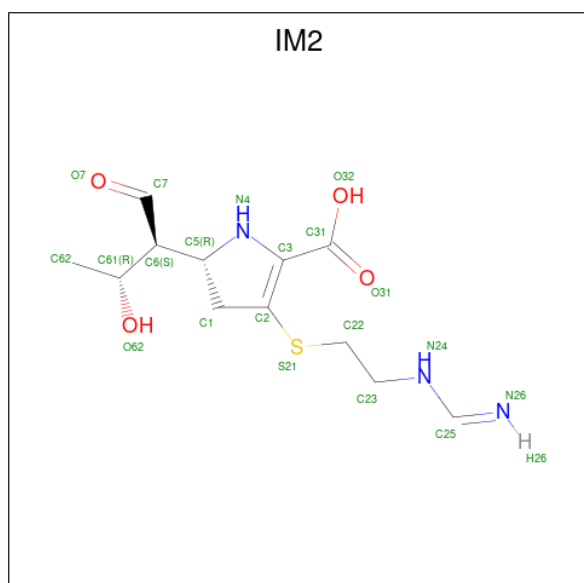
- Molecule 1 is a protein called Beta-lactamase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	268	Total	C	N	O	S	0	2	0
			2045	1278	360	399	8			
1	B	266	Total	C	N	O	S	0	2	0
			2030	1270	358	394	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	63	GLY	CYS	engineered mutation	UNP Q0Z8S4
A	286	ALA	-	expression tag	UNP Q0Z8S4
B	63	GLY	CYS	engineered mutation	UNP Q0Z8S4
B	286	ALA	-	expression tag	UNP Q0Z8S4

- Molecule 2 is (5R)-5-[(1S,2R)-1-formyl-2-hydroxypropyl]-3-[(2-{[(E)-iminomethyl]amino}ethyl)sulfanyl]-4,5-dihydro-1H-pyrrole-2-carboxylic acid (CCD ID: IM2) (formula: C₁₂H₁₉N₃O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			20	12	3	4	1		
2	B	1	Total	C	N	O	S	0	0
			20	12	3	4	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	207	Total	O	0	0
			207	207		
3	B	184	Total	O	0	0
			184	184		

SEQUENCE-PLOTS INFOmissingINFO

3 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.92Å 81.48Å 71.72Å 90.00° 101.97° 90.00°	Depositor
Resolution (Å)	35.80 – 1.38 35.80 – 1.38	Depositor EDS
% Data completeness (in resolution range)	97.5 (35.80-1.38) 97.4 (35.80-1.38)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 1.36Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.140 , 0.171 (Not available) , 0.185	Depositor DCC
R_{free} test set	4961 reflections (4.78%)	wwPDB-VP
Wilson B-factor (Å ²)	11.6	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	4506	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.65% 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.

4 Model quality

4.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: IM2

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	0.32	0/2084	0.55	0/2820
1	B	0.30	0/2069	0.54	0/2800
All	All	0.31	0/4153	0.54	0/5620

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

4.2 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	2045	0	2047	10	1
1	B	2030	0	2033	12	0
2	A	20	0	14	4	0
2	B	20	0	14	1	0
3	A	207	0	0	0	0
3	B	184	0	0	5	0
All	All	4506	0	4108	22	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 3.

All (22) 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:210:GLN:HE22	1:B:210:GLN:HE22	1.24	0.82
1:A:161:GLU:OE2	2:A:301:IM2:H621	1.81	0.81
1:A:150:ILE:O	1:A:187:LYS:HD3	1.90	0.71
2:B:301:IM2:H61	3:B:402:HOH:O	1.94	0.68
1:B:30:LYS:HE3	1:B:34:GLU:OE2	2.00	0.62
1:B:22:LEU:O	1:B:26:THR:HG23	2.00	0.62
1:B:89:LEU:HD22	1:B:136:GLU:HG3	1.83	0.60
1:B:222:LYS:NZ	3:B:403:HOH:O	2.36	0.58
1:A:73:LEU:C	1:A:73:LEU:HD23	2.31	0.55
1:A:165[A]:SER:OG	2:A:301:IM2:C62	2.56	0.54
1:B:56:MET:O	1:B:179:ILE:HG22	2.08	0.53
1:B:247:GLN:HG3	3:B:558:HOH:O	2.09	0.52
1:A:161:GLU:CD	2:A:301:IM2:H621	2.35	0.51
1:B:223:ASP:HB3	3:B:566:HOH:O	2.11	0.49
1:A:249:ARG:HG2	1:A:251:TYR:OH	2.12	0.49
1:B:73:LEU:C	1:B:73:LEU:HD23	2.38	0.49
1:B:99:TRP:HZ2	3:B:496:HOH:O	1.95	0.49
1:A:20:GLU:HG2	1:A:284:SER:HB3	1.97	0.46
1:A:165[A]:SER:OG	2:A:301:IM2:H622	2.17	0.44
1:B:96:ILE:HD11	1:B:107:LEU:HD22	2.00	0.43
1:B:140:PRO:HA	1:B:157:LEU:HD23	2.01	0.43
1:A:95:MET:SD	1:A:135:ARG:NH1	2.94	0.41

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:20:GLU:OE1	1:A:80[A]:SER:OG[2_546]	2.16	0.04

4.3 Torsion angles [i](#)

4.3.1 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	268/2 (13400%)	264 (98%)	4 (2%)	0	100	100
1	B	266/2 (13300%)	263 (99%)	3 (1%)	0	100	100
All	All	534/4 (13350%)	527 (99%)	7 (1%)	0	100	100

There are no Ramachandran outliers to report.

4.3.2 Protein sidechains ⓘ

There are no protein residues with a non-rotameric sidechain to report in this entry.

4.3.3 RNA ⓘ

There are no RNA molecules in this entry.

4.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

4.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

4.6 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	IM2	A	301	1	15,20,20	2.11	3 (20%)	8,26,26	4.09	2 (25%)
2	IM2	B	301	1	15,20,20	2.05	3 (20%)	8,26,26	3.70	2 (25%)

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	IM2	A	301	1	-	8/18/32/32	0/1/1/1
2	IM2	B	301	1	-	13/18/32/32	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	IM2	C25-N24	-5.73	1.21	1.34
2	B	301	IM2	C25-N24	-5.55	1.21	1.34
2	A	301	IM2	C3-N4	-4.62	1.24	1.36
2	B	301	IM2	C3-N4	-4.38	1.25	1.36
2	B	301	IM2	O32-C31	-2.16	1.24	1.30
2	A	301	IM2	O32-C31	-2.05	1.25	1.30

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	IM2	C3-C2-S21	-11.19	108.63	124.37
2	B	301	IM2	C3-C2-S21	-10.11	110.15	124.37
2	A	301	IM2	O7-C7-C6	-2.51	119.29	125.27
2	B	301	IM2	O7-C7-C6	-2.43	119.46	125.27

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	IM2	C1-C5-C6-C7
2	A	301	IM2	C1-C5-C6-C61
2	A	301	IM2	C2-C3-C31-O31
2	A	301	IM2	C2-C3-C31-O32
2	A	301	IM2	N4-C3-C31-O31
2	A	301	IM2	N4-C3-C31-O32
2	B	301	IM2	C1-C2-S21-C22
2	B	301	IM2	C1-C5-C6-C7
2	B	301	IM2	C1-C5-C6-C61
2	B	301	IM2	C7-C6-C61-O62
2	B	301	IM2	C7-C6-C61-C62
2	B	301	IM2	C5-C6-C61-O62

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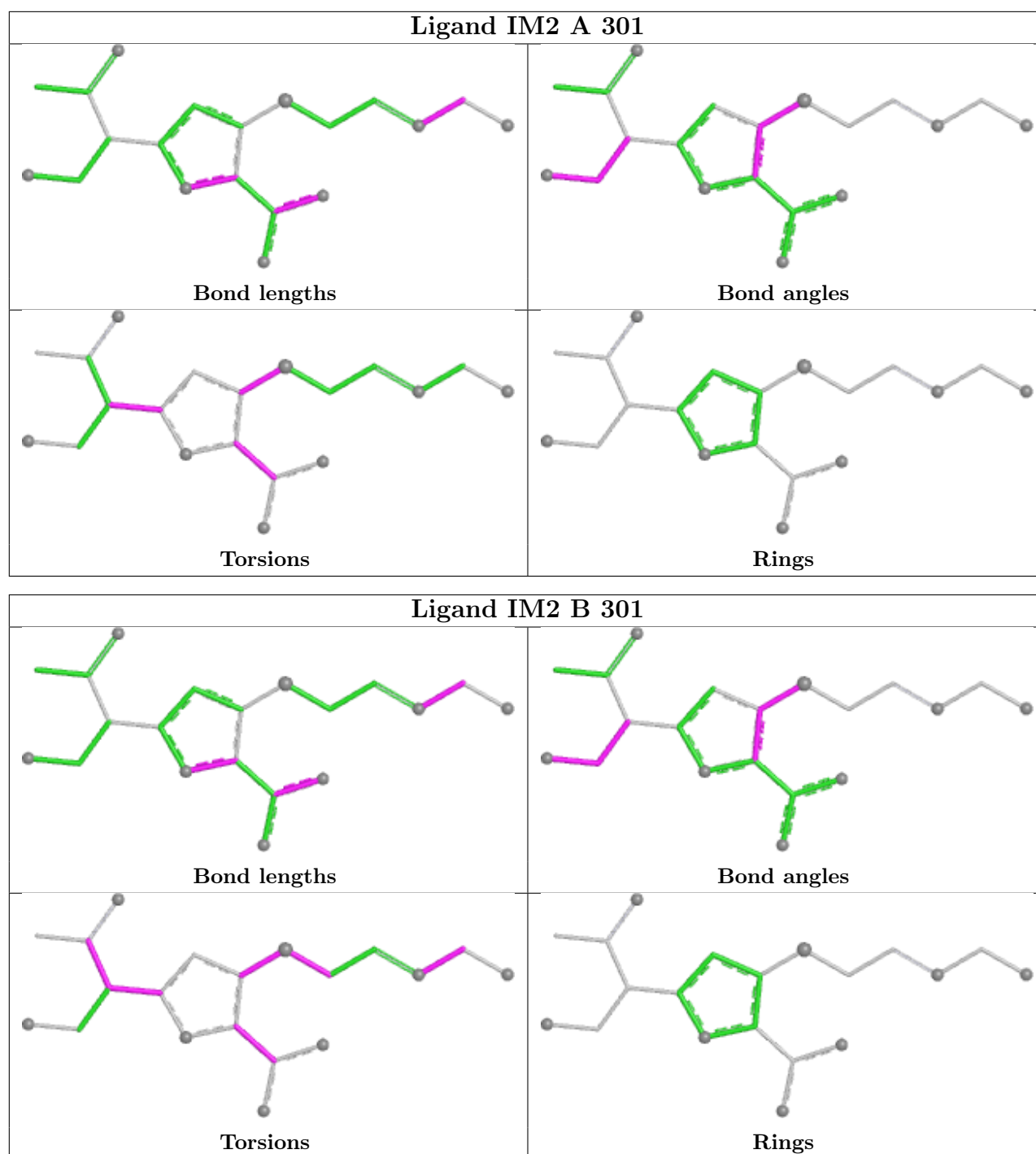
Mol	Chain	Res	Type	Atoms
2	B	301	IM2	C5-C6-C61-C62
2	B	301	IM2	C2-C3-C31-O31
2	B	301	IM2	C2-C3-C31-O32
2	B	301	IM2	N4-C3-C31-O31
2	B	301	IM2	N4-C3-C31-O32
2	B	301	IM2	C23-C22-S21-C2
2	B	301	IM2	N26-C25-N24-C23
2	A	301	IM2	C1-C2-S21-C22
2	A	301	IM2	C3-C2-S21-C22

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	IM2	4	0
2	B	301	IM2	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.



4.7 Other polymers [i](#)

There are no such residues in this entry.

4.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

5.1 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	2/2 (100%)	0.55	0 100 100	19, 19, 19, 24	0
1	B	1/2 (50%)	1.46	0 100 100	29, 29, 29, 29	0
All	All	3/4 (75%)	0.85	0 81 100	19, 19, 24, 29	0

There are no RSRZ outliers to report.

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

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

5.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.4 Ligands [i](#)

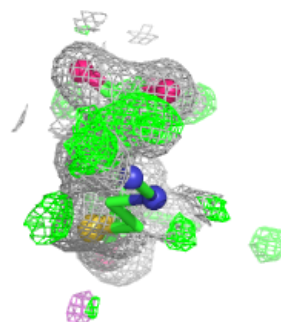
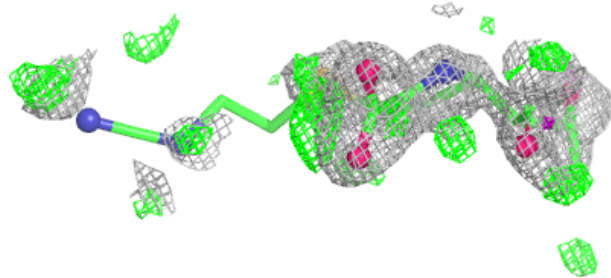
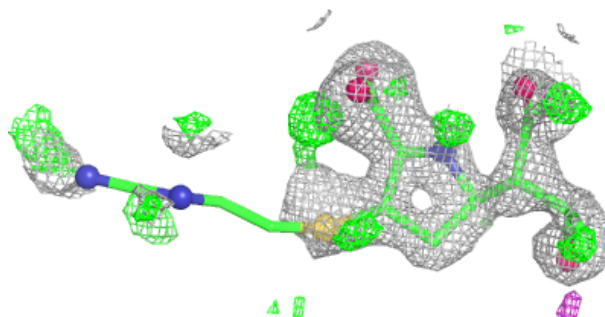
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists 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	IM2	B	301	20/20	0.81	0.18	17,27,44,56	20
2	IM2	A	301	20/20	0.86	0.16	16,27,54,54	20

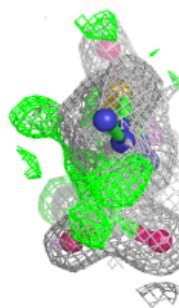
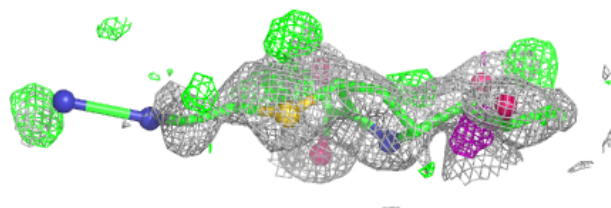
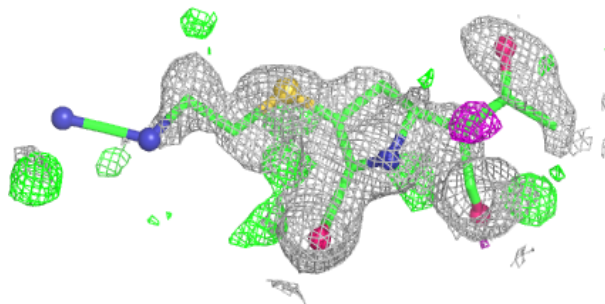
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around IM2 B 301:

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

**Electron density around IM2 A 301:**

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



5.5 Other polymers [i](#)

There are no such residues in this entry.