



# wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 20, 2020 – 03:28 PM BST

PDB ID : 5YPK  
Title : Crystal structure of NDM-1 bound to hydrolyzed imipenem representing an EI2 complex  
Authors : Feng, H.; Wang, D.; Liu, W.  
Deposited on : 2017-11-02  
Resolution : 2.00 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 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 : 2.13.1  
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) : 2.13.1

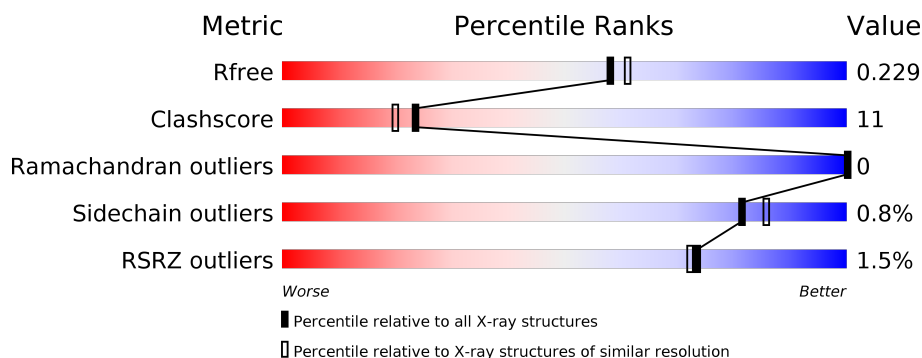
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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))
$R_{free}$	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 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	242	<div> <div></div> <div>84%10%6%</div> </div>
1	B	242	<div> <div>%</div> <div>81%12%6%</div> </div>
1	C	242	<div> <div>%</div> <div>82%12%6%</div> </div>
1	D	242	<div> <div>%</div> <div>83%11%5%</div> </div>
1	E	242	<div> <div>%</div> <div>81%13%6%</div> </div>
1	F	242	<div> <div></div> <div>78%16%6%</div> </div>

Continued on next page...

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
1	G	242	
1	H	242	

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:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SO4	C	305	-	-	X	-

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 15722 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 Metallo-beta-lactamase NDM-1.

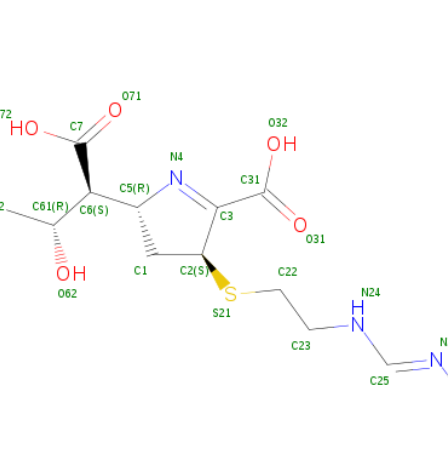
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	228	Total	C	N	O	S	0	0	0
			1697	1067	302	320	8			
1	B	228	Total	C	N	O	S	0	0	0
			1697	1067	302	320	8			
1	C	228	Total	C	N	O	S	0	0	0
			1697	1067	302	320	8			
1	D	229	Total	C	N	O	S	0	0	0
			1701	1069	303	321	8			
1	E	228	Total	C	N	O	S	0	0	0
			1697	1067	302	320	8			
1	F	228	Total	C	N	O	S	0	0	0
			1697	1067	302	320	8			
1	G	229	Total	C	N	O	S	0	0	0
			1701	1069	303	321	8			
1	H	228	Total	C	N	O	S	0	0	0
			1697	1067	302	320	8			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	2	Total	Zn	0	0
			2	2		
2	D	2	Total	Zn	0	0
			2	2		
2	E	2	Total	Zn	0	0
			2	2		
2	H	2	Total	Zn	0	0
			2	2		
2	B	2	Total	Zn	0	0
			2	2		
2	C	2	Total	Zn	0	0
			2	2		

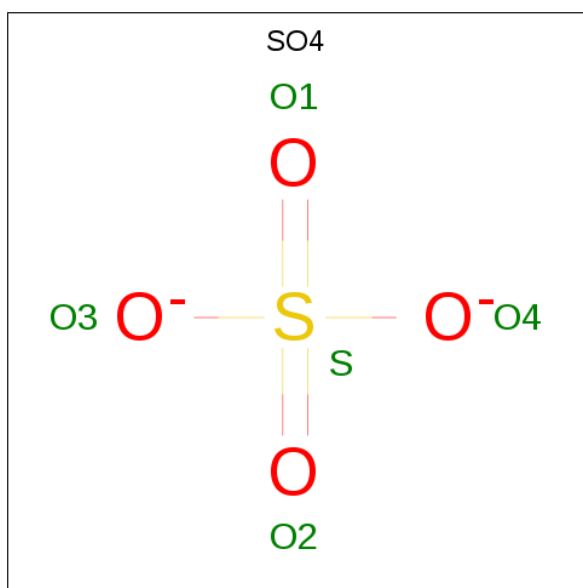
*Continued on next page...*

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Zn 2 2	0	0
2	F	2	Total Zn 2 2	0	0

- # HIW
- 
- ORTEP diagram of the chemical structure of 2,2-dimethyl-4,5-dihydro-4H-pyrimidin-5-yl 2-methylpropanoate. The structure shows a pyrimidine ring with a methyl group at C2, a methyl ester group at C5, and a 2-methylpropanoate group at C4. The atoms are labeled with their respective element symbols and numbers. The structure is shown in a 3D perspective view with thermal ellipsoids at the 50% probability level.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 21	C 12	N 3	O 5	S 1	0	0
3	B	1	Total 21	C 12	N 3	O 5	S 1	0	0
3	C	1	Total 21	C 12	N 3	O 5	S 1	0	0
3	D	1	Total 21	C 12	N 3	O 5	S 1	0	0
3	E	1	Total 21	C 12	N 3	O 5	S 1	0	0
3	F	1	Total 21	C 12	N 3	O 5	S 1	0	0
3	G	1	Total 21	C 12	N 3	O 5	S 1	0	0
3	H	1	Total 21	C 12	N 3	O 5	S 1	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	C	1	Total	O	S	0	0
			5	4	1		
4	D	1	Total	O	S	0	0
			5	4	1		
4	E	1	Total	O	S	0	0
			5	4	1		
4	F	1	Total	O	S	0	0
			5	4	1		

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	G	1	Total	Na	0	0
			1	1		
5	D	2	Total	Na	0	0
			2	2		
5	C	2	Total	Na	0	0
			2	2		
5	F	1	Total	Na	0	0
			1	1		

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	E	1	Total	Na	0	0
			1	1		

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	1	Total	Cl	0	0
			1	1		


- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	223	Total	O	0	0
			223	223		
7	B	247	Total	O	0	0
			247	247		
7	C	259	Total	O	0	0
			259	259		
7	D	267	Total	O	0	0
			267	267		
7	E	227	Total	O	0	0
			227	227		
7	F	250	Total	O	0	0
			250	250		
7	G	219	Total	O	0	0
			219	219		
7	H	219	Total	O	0	0
			219	219		

### 3 Residue-property plots


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.

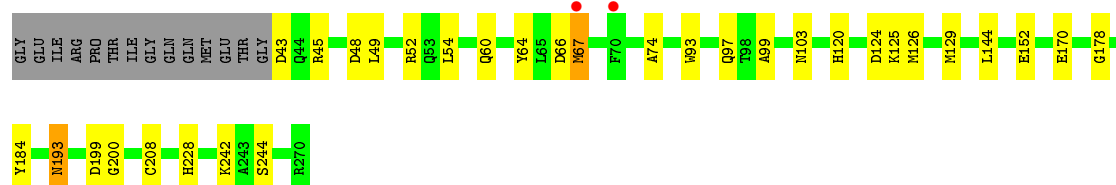
#### • Molecule 1: Metallo-beta-lactamase NDM-1

Chain A: 




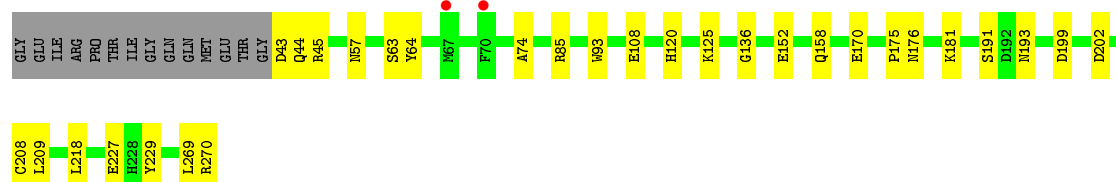
#### • Molecule 1: Metallo-beta-lactamase NDM-1

Chain B: 




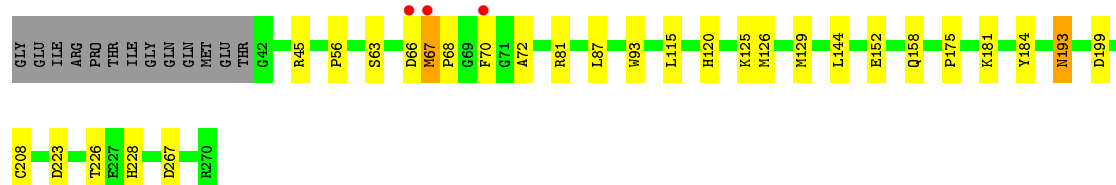
#### • Molecule 1: Metallo-beta-lactamase NDM-1

Chain C: 




#### • Molecule 1: Metallo-beta-lactamase NDM-1

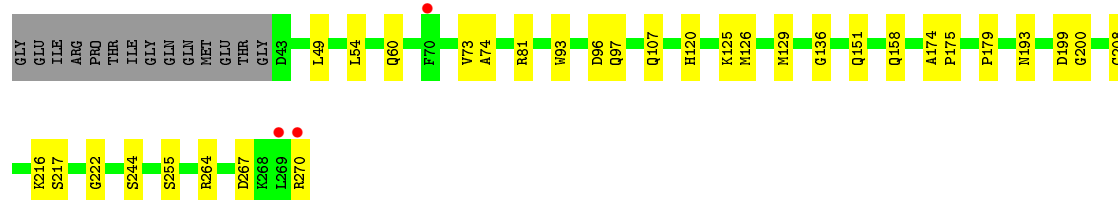
Chain D: 






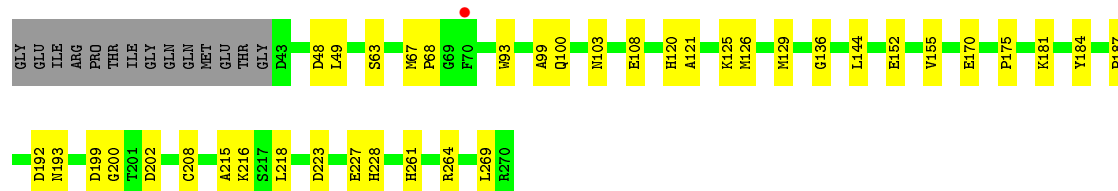
- Molecule 1: Metallo-beta-lactamase NDM-1

Chain E: 




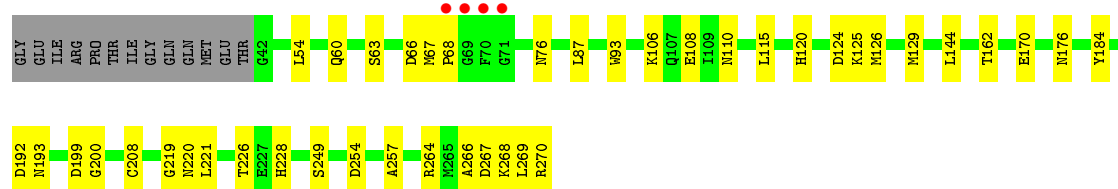
- Molecule 1: Metallo-beta-lactamase NDM-1

Chain F: 



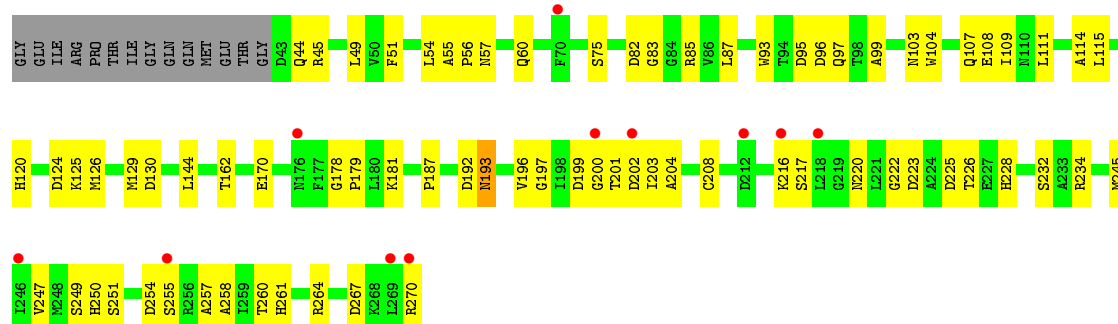
- Molecule 1: Metallo-beta-lactamase NDM-1

Chain G: 



- Molecule 1: Metallo-beta-lactamase NDM-1

Chain H: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	69.24Å 73.61Å 154.52Å 90.00° 90.08° 90.00°	Depositor
Resolution (Å)	42.21 – 2.00 47.96 – 2.00	Depositor EDS
% Data completeness (in resolution range)	90.0 (42.21-2.00) 94.7 (47.96-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.23 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.9-1692	Depositor
R, $R_{free}$	0.186 , 0.229 0.187 , 0.229	Depositor DCC
$R_{free}$ test set	2021 reflections (2.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.5	Xtriage
Anisotropy	0.287	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 33.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.267 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	15722	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 67.22 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 5.3322e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: NA, ZN, CL, SO4, HIW

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	0.27	0/1737	0.43	0/2366
1	B	0.27	0/1737	0.47	0/2366
1	C	0.27	0/1737	0.42	0/2366
1	D	0.27	0/1741	0.42	0/2371
1	E	0.27	0/1737	0.47	0/2366
1	F	0.26	0/1737	0.44	0/2366
1	G	0.30	0/1741	0.47	0/2371
1	H	0.26	0/1737	0.48	0/2366
All	All	0.27	0/13904	0.45	0/18938

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 [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	1697	0	1647	20	0
1	B	1697	0	1647	25	0
1	C	1697	0	1647	30	0
1	D	1701	0	1650	32	0
1	E	1697	0	1647	35	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	1697	0	1647	35	0
1	G	1701	0	1650	36	1
1	H	1697	0	1647	93	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0
3	A	21	0	0	0	0
3	B	21	0	0	0	0
3	C	21	0	0	2	0
3	D	21	0	0	0	0
3	E	21	0	0	0	0
3	F	21	0	0	0	0
3	G	21	0	0	3	0
3	H	21	0	0	2	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	C	10	0	0	4	0
4	D	5	0	0	0	0
4	E	5	0	0	1	0
4	F	5	0	0	1	0
5	C	2	0	0	0	0
5	D	2	0	0	0	0
5	E	1	0	0	0	0
5	F	1	0	0	0	0
5	G	1	0	0	0	0
6	D	1	0	0	1	0
7	A	223	0	0	10	4
7	B	247	0	0	10	4
7	C	259	0	0	27	4
7	D	267	0	0	15	2
7	E	227	0	0	26	7
7	F	250	0	0	13	2
7	G	219	0	0	16	3
7	H	219	0	0	69	6
All	All	15722	0	13182	302	17

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 11.

The worst 5 of 302 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:43:ASP:N	7:A:402:HOH:O	1.56	1.32
1:H:234:ARG:NH2	7:H:403:HOH:O	1.60	1.31
1:H:95:ASP:OD2	7:H:402:HOH:O	1.53	1.23
1:C:158:GLN:OE1	7:C:401:HOH:O	1.56	1.21
1:E:74:ALA:O	7:E:404:HOH:O	1.52	1.21

The worst 5 of 17 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)
7:E:591:HOH:O	7:H:571:HOH:O[2_748]	1.08	1.12
7:A:494:HOH:O	7:H:557:HOH:O[2_748]	1.13	1.07
7:B:623:HOH:O	7:B:631:HOH:O[2_849]	1.28	0.92
7:C:618:HOH:O	7:E:605:HOH:O[1_565]	1.36	0.84
7:D:538:HOH:O	7:F:584:HOH:O[1_545]	1.38	0.82

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	226/242 (93%)	223 (99%)	3 (1%)	0	100	100
1	B	226/242 (93%)	223 (99%)	3 (1%)	0	100	100
1	C	226/242 (93%)	222 (98%)	4 (2%)	0	100	100
1	D	227/242 (94%)	223 (98%)	4 (2%)	0	100	100
1	E	226/242 (93%)	221 (98%)	5 (2%)	0	100	100
1	F	226/242 (93%)	222 (98%)	4 (2%)	0	100	100
1	G	227/242 (94%)	224 (99%)	3 (1%)	0	100	100
1	H	226/242 (93%)	223 (99%)	3 (1%)	0	100	100

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	1810/1936 (94%)	1781 (98%)	29 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	171/182 (94%)	170 (99%)	1 (1%)	86	90
1	B	171/182 (94%)	168 (98%)	3 (2%)	59	63
1	C	171/182 (94%)	170 (99%)	1 (1%)	86	90
1	D	171/182 (94%)	169 (99%)	2 (1%)	71	76
1	E	171/182 (94%)	170 (99%)	1 (1%)	86	90
1	F	171/182 (94%)	170 (99%)	1 (1%)	86	90
1	G	171/182 (94%)	170 (99%)	1 (1%)	86	90
1	H	171/182 (94%)	170 (99%)	1 (1%)	86	90
All	All	1368/1456 (94%)	1357 (99%)	11 (1%)	81	86

5 of 11 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	193	ASN
1	D	67	MET
1	F	193	ASN
1	B	193	ASN
1	E	193	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 39 ligands modelled in this entry, 24 are monoatomic - leaving 15 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$
4	SO4	D	304	-	4,4,4	0.14	0	6,6,6	0.05	0
4	SO4	F	304	-	4,4,4	0.15	0	6,6,6	0.05	0
3	HIW	C	303	2	12,21,21	4.05	2 (16%)	8,28,28	1.26	2 (25%)
3	HIW	B	303	2	12,21,21	4.09	2 (16%)	8,28,28	1.02	1 (12%)
3	HIW	E	303	2	12,21,21	4.18	2 (16%)	8,28,28	1.49	1 (12%)
3	HIW	D	303	2	12,21,21	4.18	3 (25%)	8,28,28	0.89	1 (12%)
3	HIW	G	303	2	12,21,21	4.10	2 (16%)	8,28,28	1.11	1 (12%)
3	HIW	F	303	2	12,21,21	4.05	2 (16%)	8,28,28	1.26	1 (12%)
4	SO4	E	304	-	4,4,4	0.14	0	6,6,6	0.07	0
4	SO4	A	304	-	4,4,4	0.13	0	6,6,6	0.06	0
4	SO4	C	304	-	4,4,4	0.15	0	6,6,6	0.04	0
3	HIW	A	303	2	12,21,21	4.07	2 (16%)	8,28,28	1.38	2 (25%)
4	SO4	C	305	-	4,4,4	0.13	0	6,6,6	0.06	0
3	HIW	H	303	2	12,21,21	4.10	2 (16%)	8,28,28	0.98	0
4	SO4	B	304	-	4,4,4	0.15	0	6,6,6	0.06	0

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
3	HIW	C	303	2	-	6/13/34/34	0/1/1/1
3	HIW	B	303	2	-	5/13/34/34	0/1/1/1
3	HIW	E	303	2	-	7/13/34/34	0/1/1/1
3	HIW	D	303	2	-	9/13/34/34	0/1/1/1
3	HIW	G	303	2	-	10/13/34/34	0/1/1/1
3	HIW	F	303	2	-	5/13/34/34	0/1/1/1
3	HIW	A	303	2	-	7/13/34/34	0/1/1/1
3	HIW	H	303	2	-	6/13/34/34	0/1/1/1

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	303	HIW	C31-C3	-12.48	1.33	1.52
3	C	303	HIW	C31-C3	-12.13	1.33	1.52
3	B	303	HIW	C31-C3	-12.10	1.33	1.52
3	A	303	HIW	C31-C3	-12.06	1.33	1.52
3	G	303	HIW	C31-C3	-12.06	1.33	1.52

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	E	303	HIW	C1-C5-C6	-3.04	108.41	115.10
3	A	303	HIW	C6-C5-N4	-2.59	106.93	111.93
3	C	303	HIW	C6-C5-N4	-2.44	107.21	111.93
3	F	303	HIW	C6-C5-N4	-2.38	107.34	111.93
3	B	303	HIW	C6-C5-N4	-2.23	107.63	111.93

There are no chirality outliers.

5 of 55 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	303	HIW	N4-C5-C6-C7
3	C	303	HIW	N4-C5-C6-C61
3	C	303	HIW	C1-C5-C6-C7
3	C	303	HIW	C1-C5-C6-C61
3	E	303	HIW	N4-C5-C6-C7

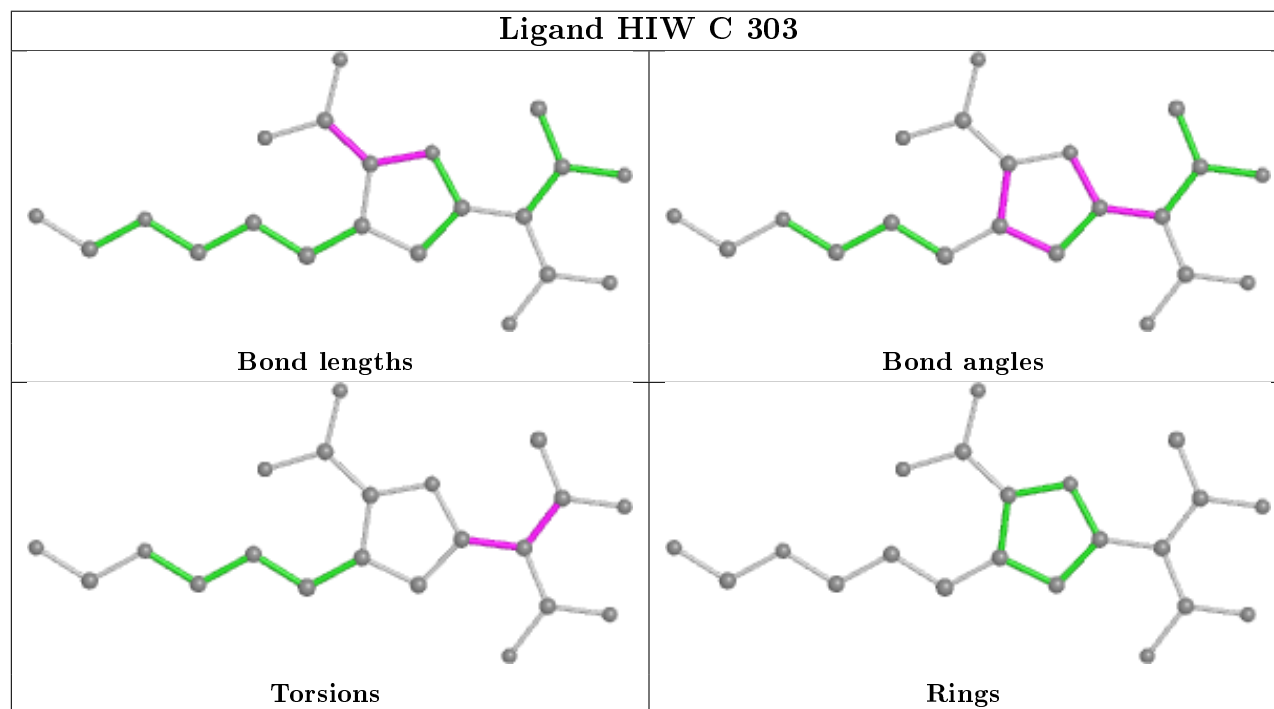


There are no ring outliers.

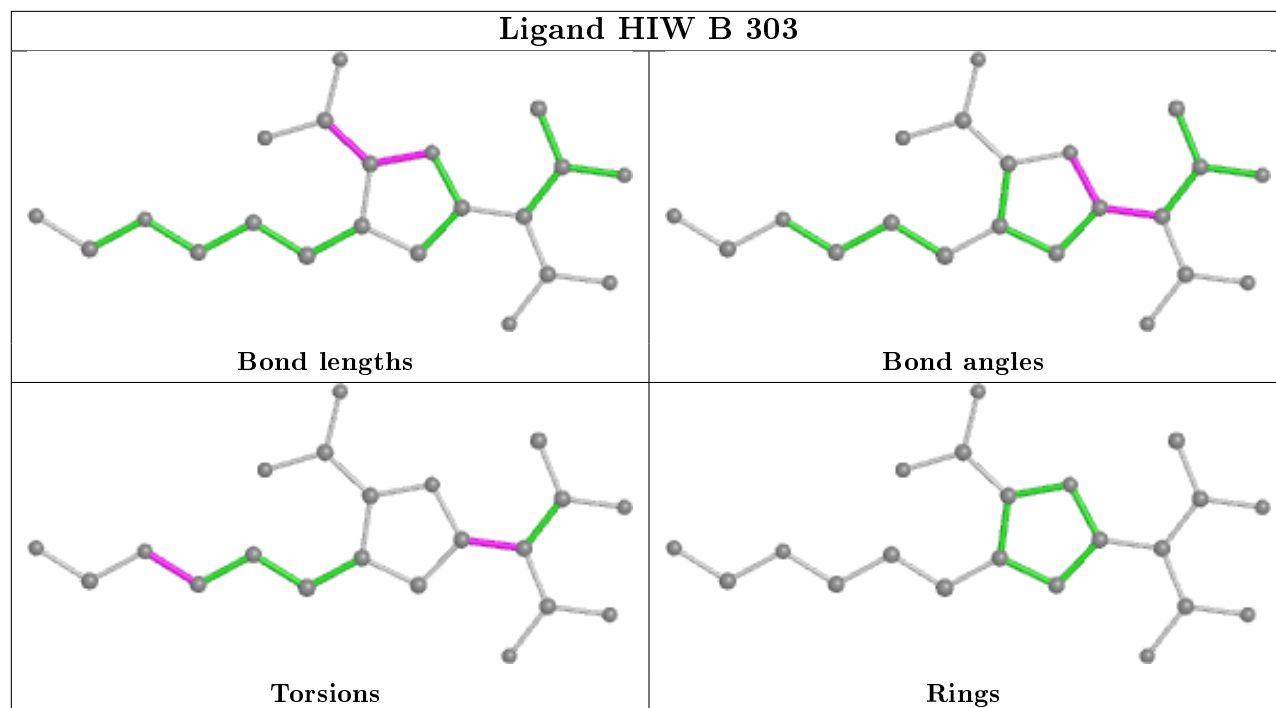
7 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	F	304	SO4	1	0
3	C	303	HIW	2	0
3	G	303	HIW	3	0
4	E	304	SO4	1	0
4	C	304	SO4	1	0
4	C	305	SO4	3	0
3	H	303	HIW	2	0

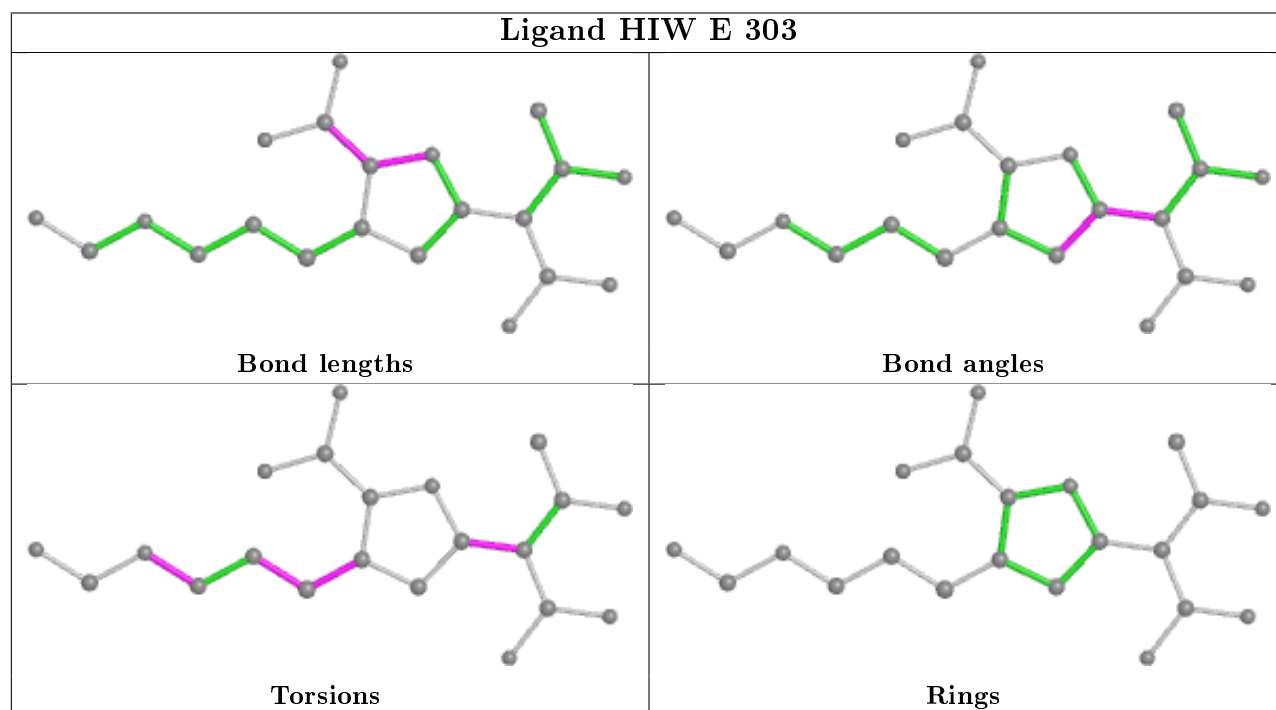
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

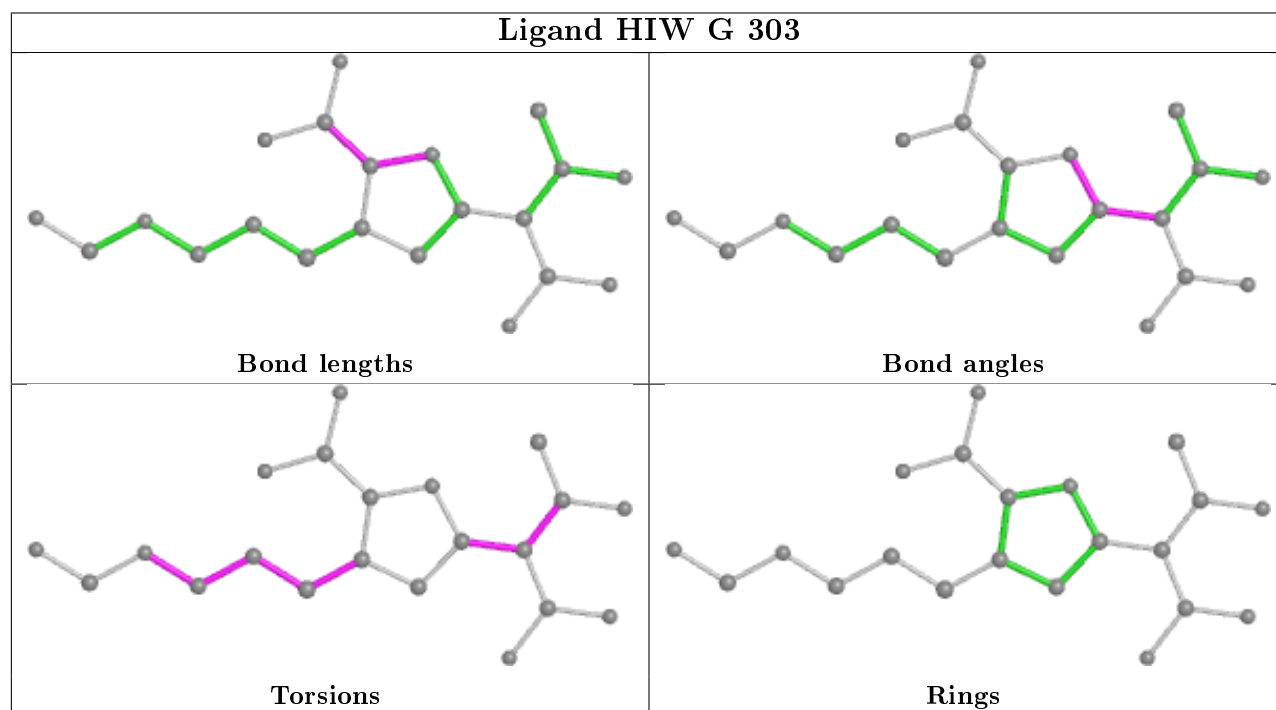
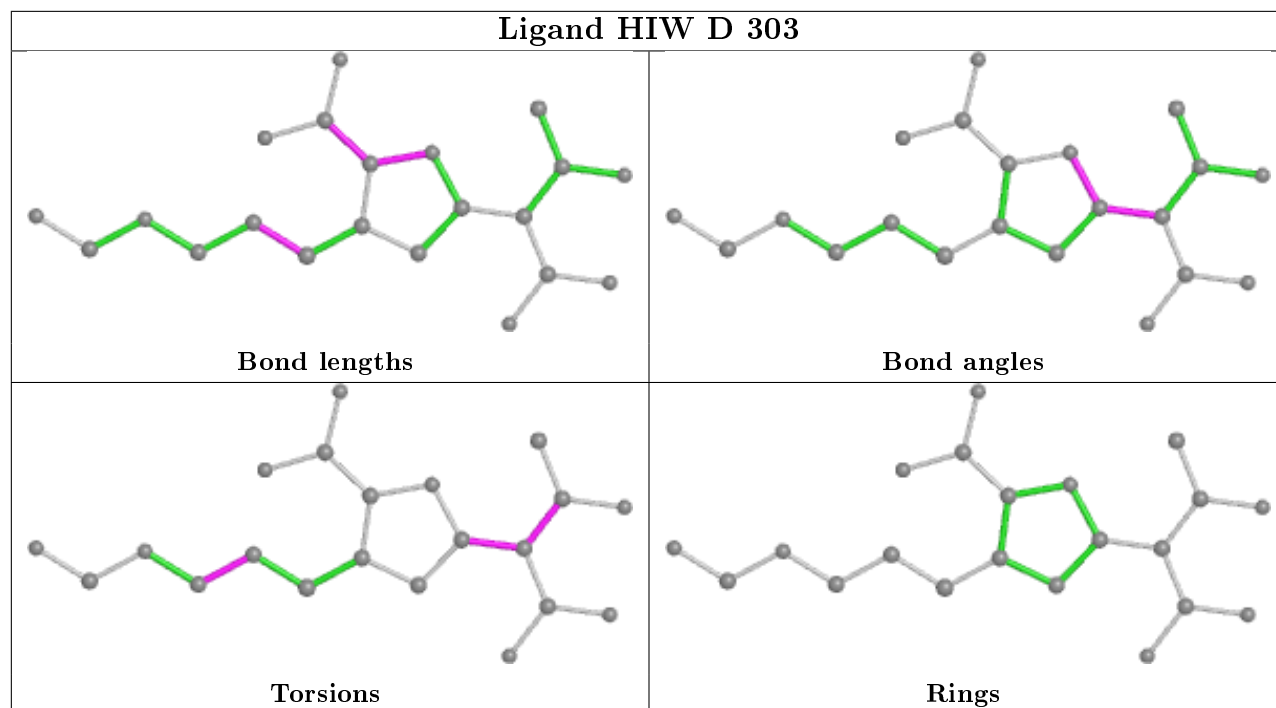


## Ligand HIW B 303

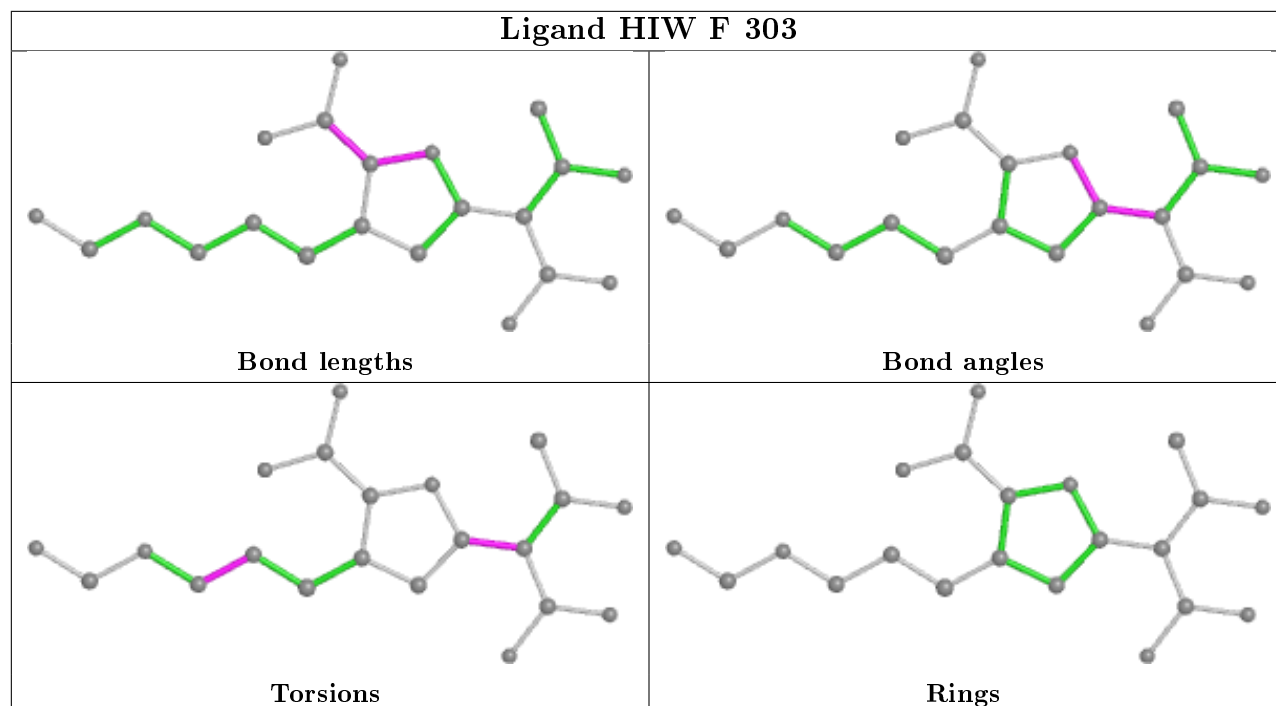


## Ligand HIW E 303

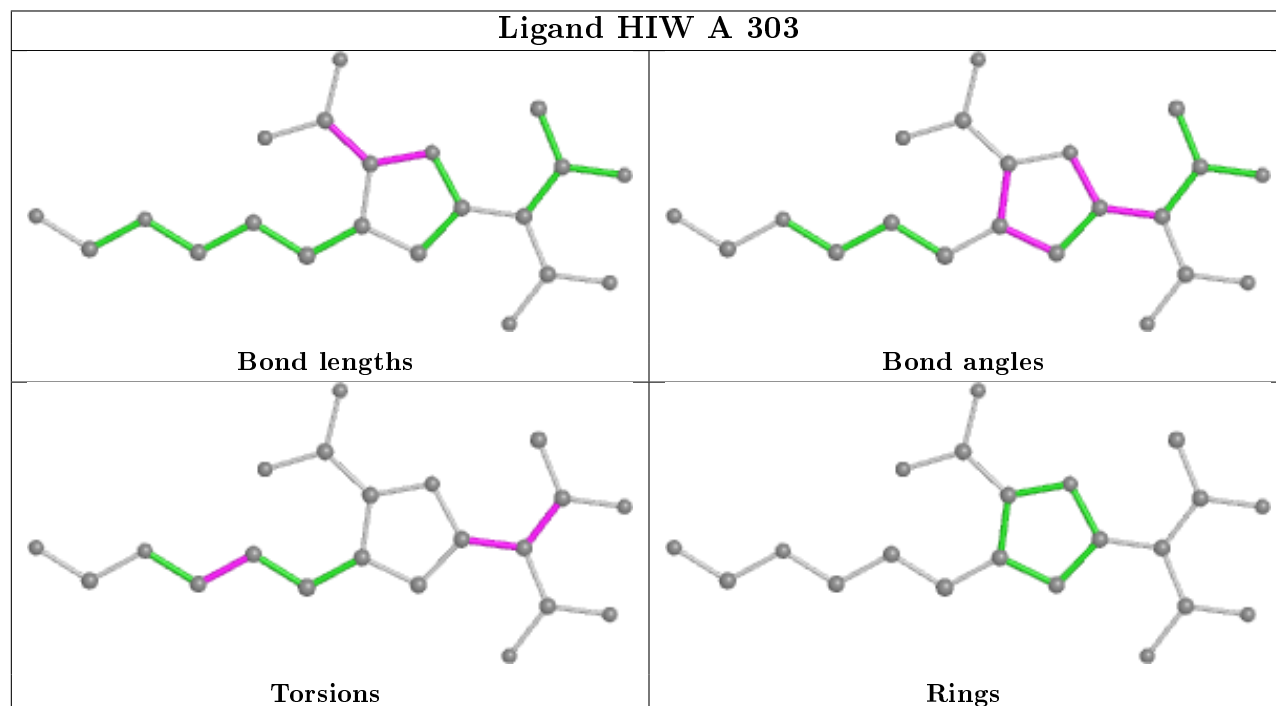


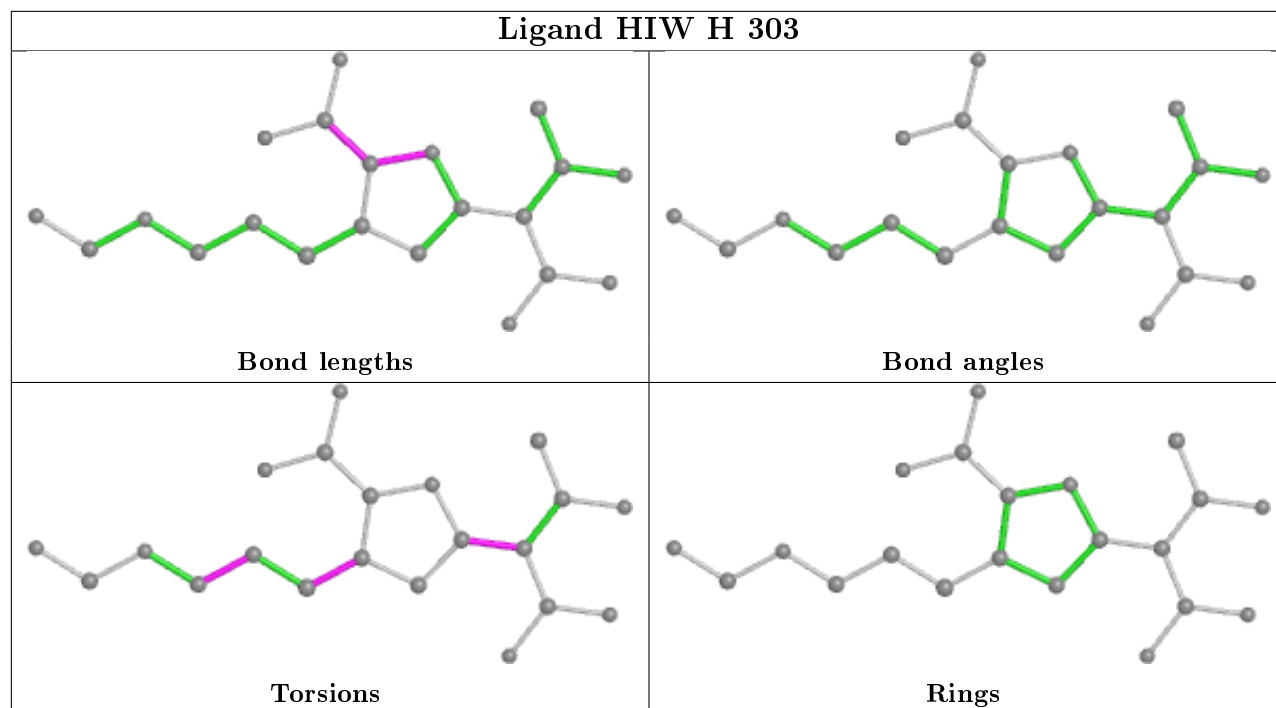


## Ligand HIW F 303



## Ligand HIW A 303





## 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 ⓘ

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	228/242 (94%)	-0.65	1 (0%) 92 92	11, 19, 38, 65	0
1	B	228/242 (94%)	-0.50	2 (0%) 84 83	12, 22, 40, 57	0
1	C	228/242 (94%)	-0.60	2 (0%) 84 83	12, 21, 37, 63	0
1	D	229/242 (94%)	-0.57	3 (1%) 77 76	11, 20, 42, 70	0
1	E	228/242 (94%)	-0.32	3 (1%) 77 76	14, 26, 46, 60	0
1	F	228/242 (94%)	-0.58	1 (0%) 92 92	13, 21, 36, 50	0
1	G	229/242 (94%)	-0.42	4 (1%) 70 68	15, 24, 43, 66	0
1	H	228/242 (94%)	0.10	11 (4%) 30 29	14, 31, 49, 70	0
All	All	1826/1936 (94%)	-0.44	27 (1%) 73 72	11, 23, 44, 70	0

The worst 5 of 27 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	270	ARG	5.8
1	E	269	LEU	4.5
1	B	70	PHE	4.2
1	D	70	PHE	3.9
1	H	200	GLY	3.8

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

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

### 6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 6.4 Ligands ⓘ

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.

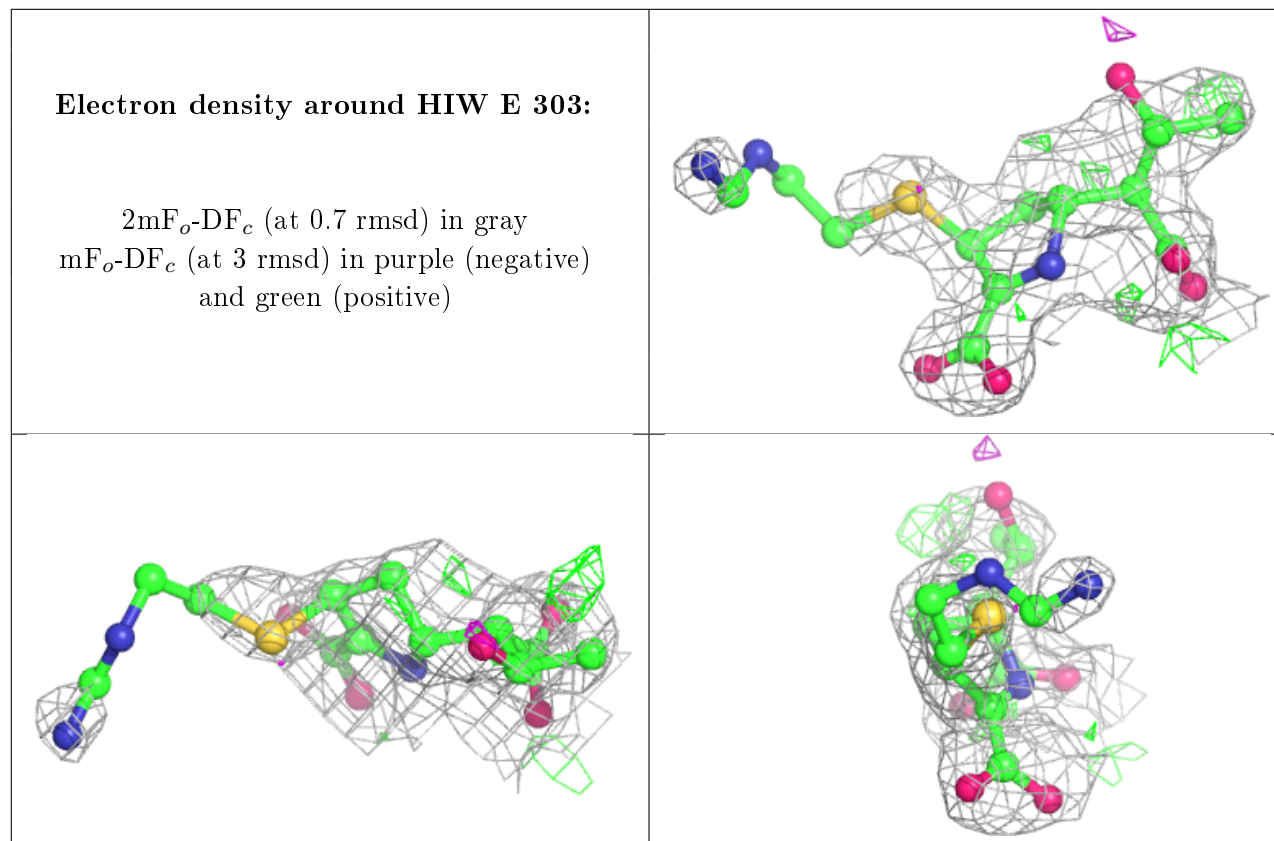
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	NA	F	305	1/1	0.69	0.11	46,46,46,46	0
3	HIW	E	303	21/21	0.79	0.21	24,39,53,57	21
3	HIW	H	303	21/21	0.83	0.24	35,42,50,54	21
5	NA	G	304	1/1	0.88	0.26	30,30,30,30	1
5	NA	C	307	1/1	0.88	0.18	32,32,32,32	1
3	HIW	C	303	21/21	0.90	0.16	19,27,34,38	21
6	CL	D	305	1/1	0.90	0.17	46,46,46,46	0
3	HIW	A	303	21/21	0.90	0.18	21,27,39,49	21
3	HIW	D	303	21/21	0.90	0.17	25,33,47,48	21
3	HIW	F	303	21/21	0.91	0.15	24,28,42,44	21
3	HIW	G	303	21/21	0.91	0.20	23,34,42,48	21
5	NA	D	306	1/1	0.92	0.16	46,46,46,46	0
5	NA	C	306	1/1	0.93	0.14	52,52,52,52	0
4	SO4	F	304	5/5	0.95	0.12	33,44,50,52	0
4	SO4	C	304	5/5	0.96	0.15	29,37,41,48	0
3	HIW	B	303	21/21	0.96	0.11	25,33,48,59	0
4	SO4	C	305	5/5	0.96	0.21	36,47,49,53	0
5	NA	E	305	1/1	0.97	0.41	28,28,28,28	1
2	ZN	H	302	1/1	0.97	0.08	44,44,44,44	0
4	SO4	E	304	5/5	0.98	0.12	34,40,44,47	0
2	ZN	B	302	1/1	0.98	0.06	32,32,32,32	0
4	SO4	A	304	5/5	0.98	0.06	28,31,37,40	0
2	ZN	A	302	1/1	0.99	0.03	23,23,23,23	0
5	NA	D	307	1/1	0.99	0.10	24,24,24,24	1
2	ZN	H	301	1/1	0.99	0.04	31,31,31,31	0
4	SO4	B	304	5/5	0.99	0.06	29,29,32,38	0
4	SO4	D	304	5/5	0.99	0.07	23,33,35,40	0
2	ZN	C	302	1/1	0.99	0.05	25,25,25,25	0
2	ZN	E	302	1/1	0.99	0.04	37,37,37,37	0
2	ZN	G	302	1/1	0.99	0.05	29,29,29,29	0
2	ZN	D	302	1/1	0.99	0.05	26,26,26,26	0
2	ZN	G	301	1/1	0.99	0.06	21,21,21,21	0
2	ZN	E	301	1/1	0.99	0.06	24,24,24,24	0
2	ZN	F	302	1/1	1.00	0.03	27,27,27,27	0
2	ZN	C	301	1/1	1.00	0.04	18,18,18,18	0
2	ZN	F	301	1/1	1.00	0.02	21,21,21,21	0
2	ZN	A	301	1/1	1.00	0.02	21,21,21,21	0

*Continued on next page...*

*Continued from previous page...*

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	B	301	1/1	1.00	0.04	21,21,21,21	0
2	ZN	D	301	1/1	1.00	0.04	19,19,19,19	0

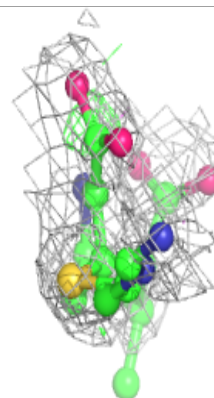
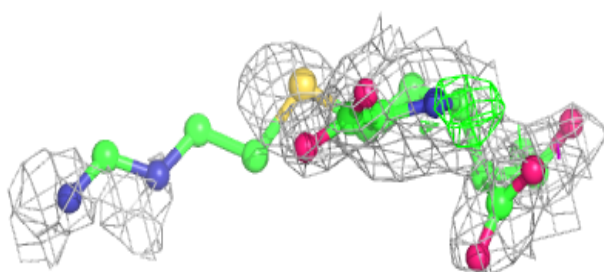
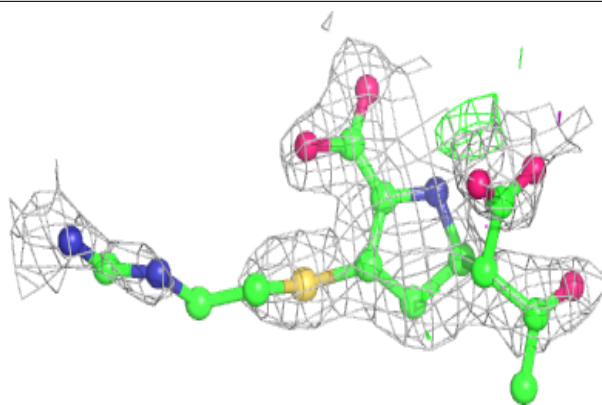
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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



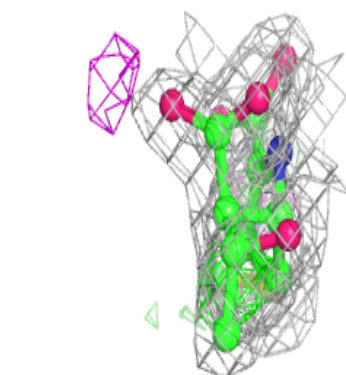
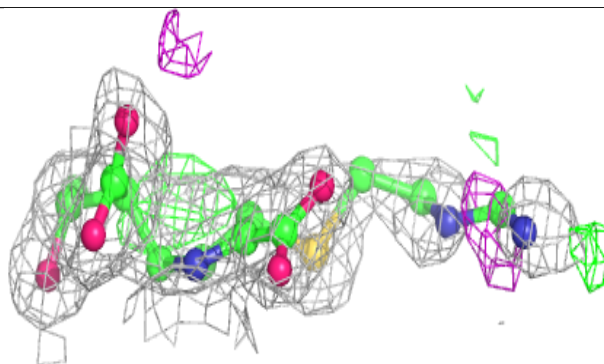
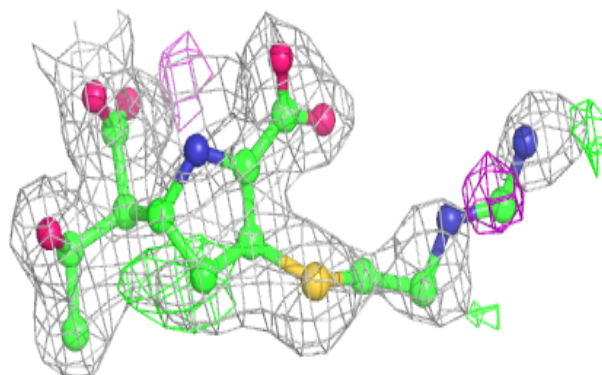


**Electron density around HIW H 303:**

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

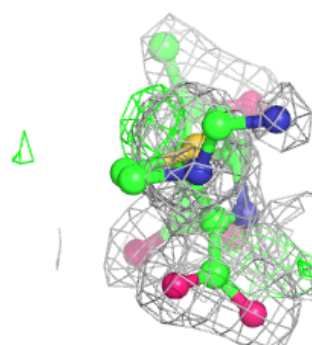
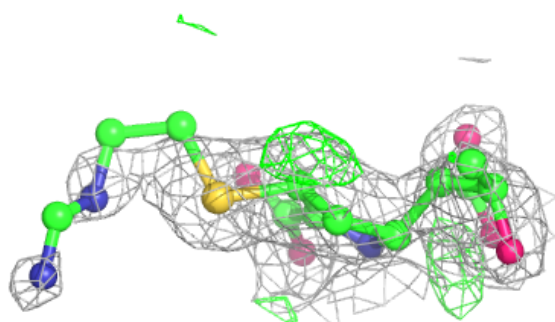
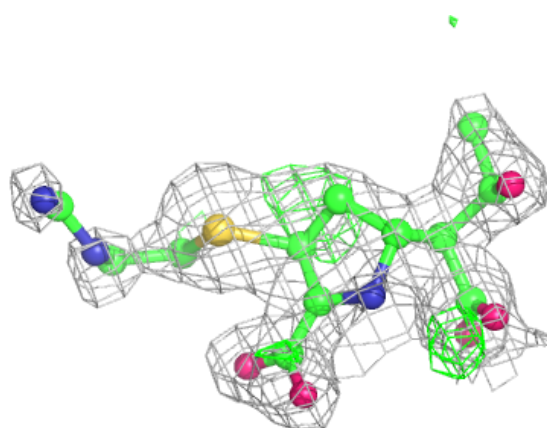
**Electron density around HIW C 303:**

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

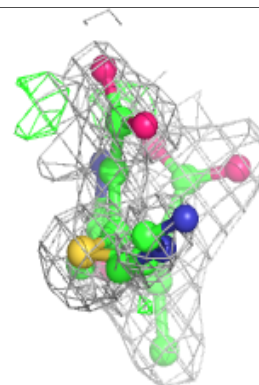
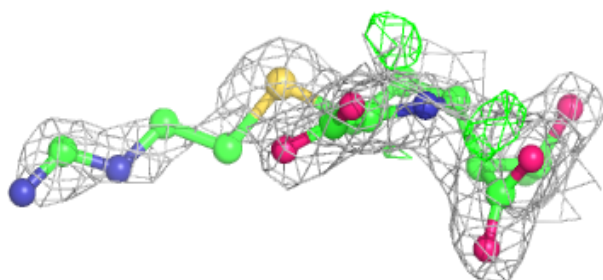
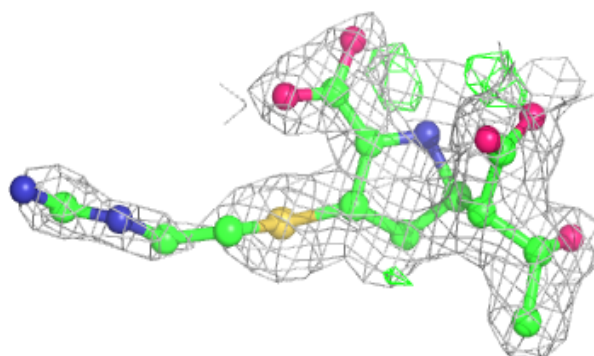


**Electron density around HIW A 303:**

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

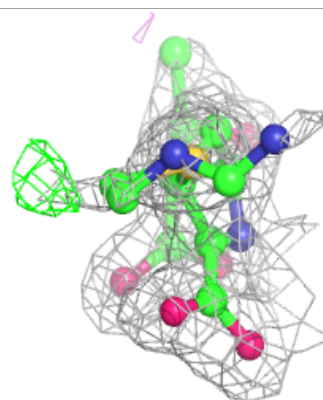
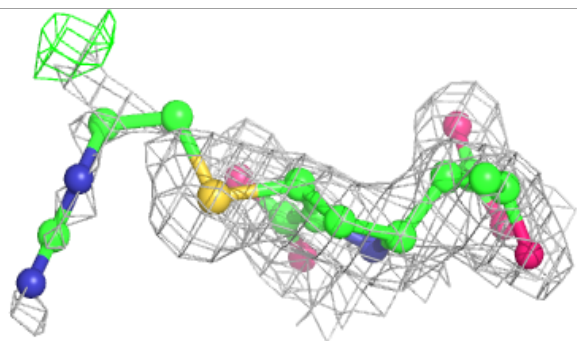
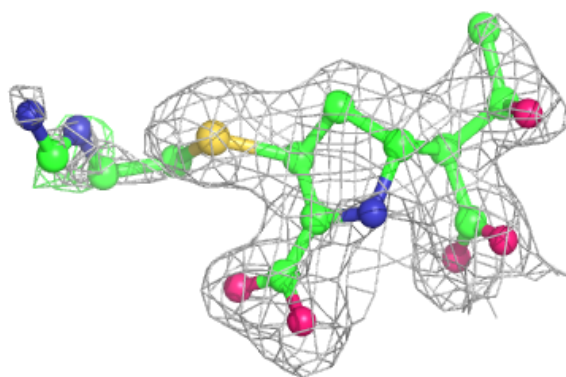
**Electron density around HIW D 303:**

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

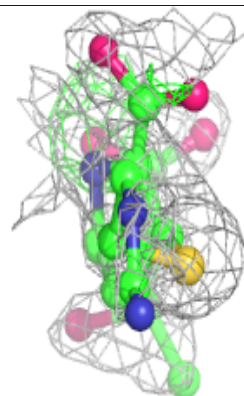
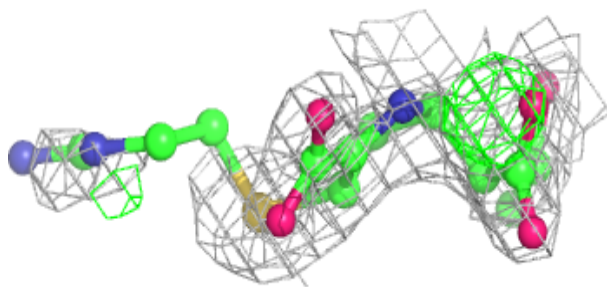
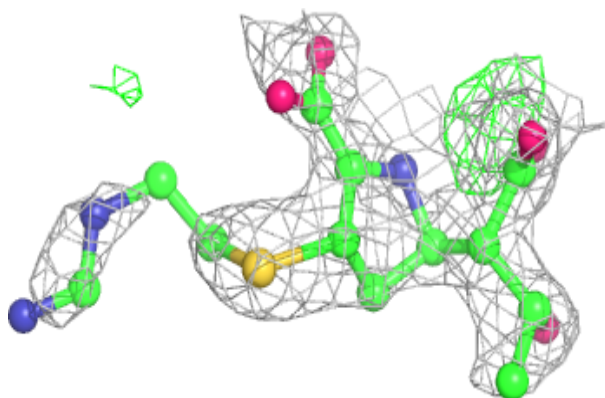


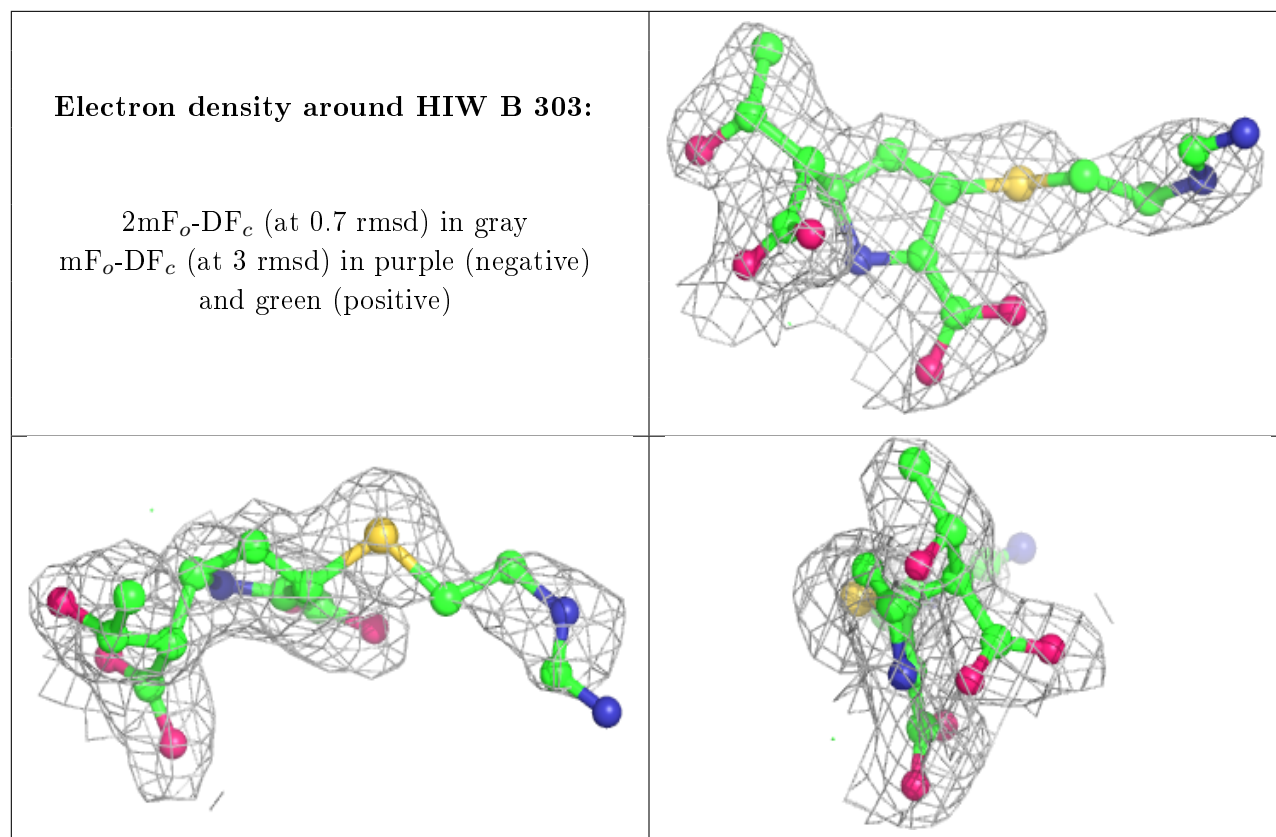
**Electron density around HIW F 303:**

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

**Electron density around HIW G 303:**

$2mF_o-DF_c$  (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.