



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 13, 2020 – 01:54 pm BST

PDB ID : 3W7G  
Title : Structure of Trypanosoma cruzi dihydroorotate dehydrogenase in complex with MII-5-013  
Authors : Inaoka, D.K.; Iida, M.; Tabuchi, T.; Lee, N.; Hashimoto, S.; Matsuoka, S.; Kuranaga, T.; Shiba, T.; Sakamoto, K.; Suzuki, S.; Balogun, E.O.; Nara, T.; Aoki, T.; Inoue, M.; Honma, T.; Tanaka, A.; Harada, S.; Kita, K.  
Deposited on : 2013-02-28  
Resolution : 1.55 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.

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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

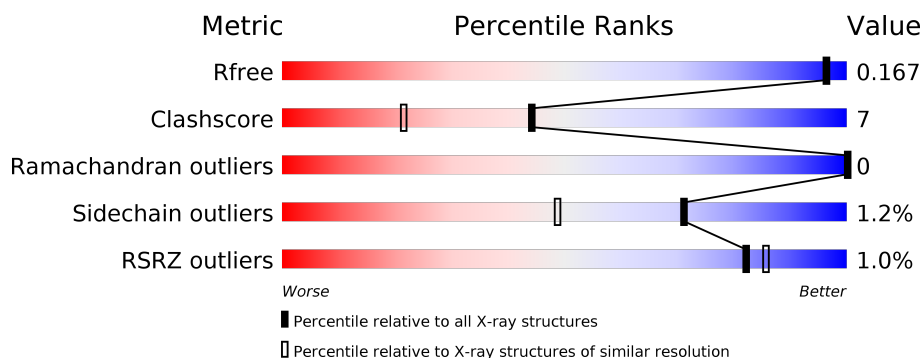
# 1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.55 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	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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	313	<div> <div style="width: 86%;"></div> <div style="width: 13%;"></div> <div style="width: 1%;"></div> </div> <div>86% 13% .</div>
1	B	313	<div> <div style="width: 2%;"></div> <div style="width: 87%;"></div> <div style="width: 11%;"></div> </div> <div>2% 87% 11% .</div>

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
2	W7G	B	401[B]	-	-	X	-
3	GOL	B	407	-	-	X	-

## 2 Entry composition [i](#)

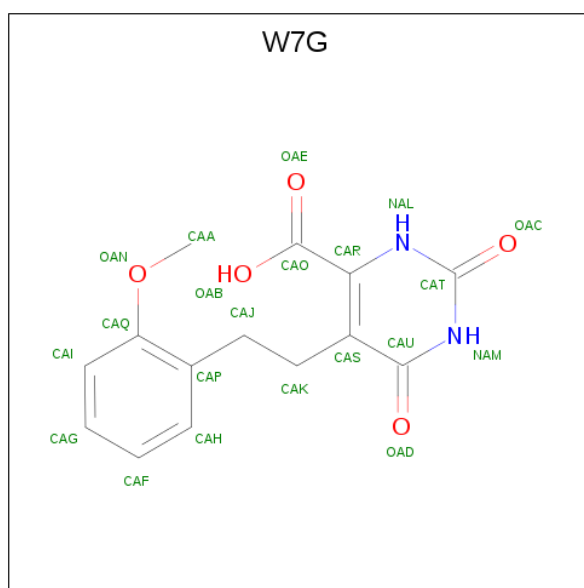
There are 6 unique types of molecules in this entry. The entry contains 5767 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Dihydroorotate dehydrogenase (fumarate).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	313	Total	C	N	O	S	0	5	0
			2415	1547	402	448	18			
1	B	313	Total	C	N	O	S	0	3	0
			2402	1539	399	445	19			

- Molecule 2 is 5-[2-(2-methoxyphenyl)ethyl]-2,6-dioxo-1,2,3,6-tetrahydropyrimidine-4-carboxylic acid (three-letter code: W7G) (formula: C<sub>14</sub>H<sub>14</sub>N<sub>2</sub>O<sub>5</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			21	14	2	5		
2	B	1	Total	C	N	O	0	1
			42	28	4	10		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	1
			12	6	6		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

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- The diagram illustrates the chemical structure of Flavin Mononucleotide (FMN). It features an isoalloxazine ring system (rings C4A, C5A, C8A, C9A) with nitrogen atoms N1, N3, N5, and N10. The ring is substituted with a ribitol chain at C10A and a phosphate group at C2. The ribitol chain consists of a ribose ring (C3', C4', C5') and a terminal CH<sub>2</sub> group (C5'). The phosphate group is shown as a phosphorus atom (P) double-bonded to one oxygen (O1P) and single-bonded to three others (O2P, O3P, O5'). The ribose ring has hydroxyl groups at C3' and C4'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 31	C 17	N 4	O 9	P 1	0	0
4	B	1	Total 31	C 17	N 4	O 9	P 1	0	0

- 
- Chemical structure of hexamminecobalt(III) ion,  $[Co(NH_3)_6]^{3+}$ . The central cobalt atom (Co) is bonded to six ammonia (NH<sub>3</sub>) ligands in an octahedral geometry. The ligands are labeled N1 through N6, and the central cobalt is labeled CO. The overall charge is +3.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total 7	Co 1	N 6	0	0

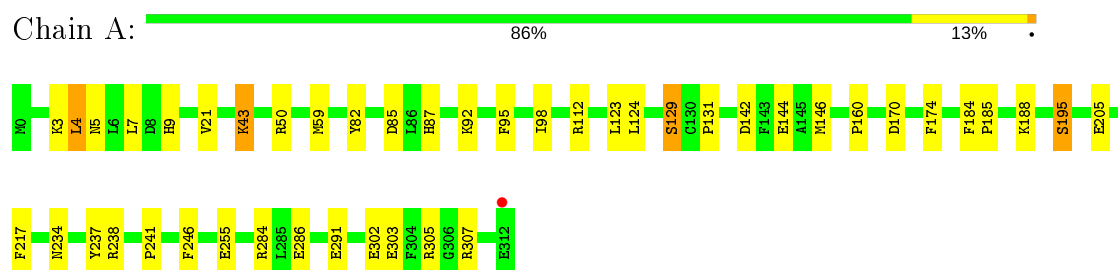
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	374	Total 374	O 374	0	0
6	B	354	Total 354	O 354	0	0

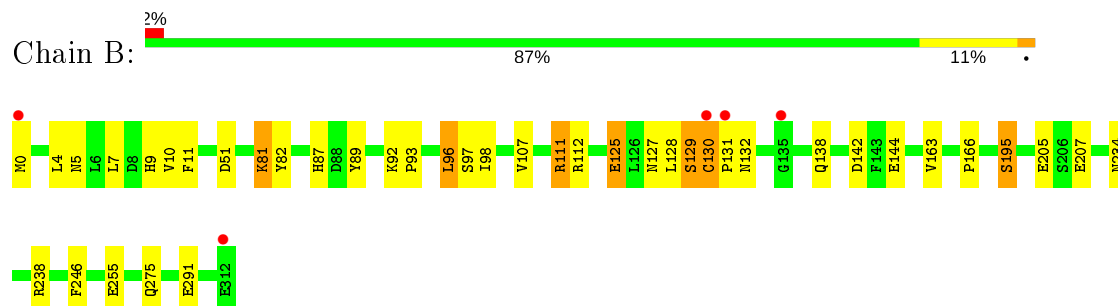
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Dihydroorotate dehydrogenase (fumarate)



- Molecule 1: Dihydroorotate dehydrogenase (fumarate)





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	67.95Å 71.83Å 129.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.88 – 1.55 29.88 – 1.55	Depositor EDS
% Data completeness (in resolution range)	93.0 (29.88-1.55) 93.0 (29.88-1.55)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.79 (at 1.55Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, $R_{free}$	0.137 , 0.165 0.138 , 0.167	Depositor DCC
$R_{free}$ test set	4295 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	11.5	Xtriage
Anisotropy	0.017	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 53.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	5767	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	14.0	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 25.53 % 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 3.1173e-03. 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: FMN, GOL, W7G, NCO

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	1.26	14/2482 (0.6%)	1.17	14/3360 (0.4%)
1	B	1.24	6/2463 (0.2%)	1.13	9/3336 (0.3%)
All	All	1.25	20/4945 (0.4%)	1.15	23/6696 (0.3%)

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	1
All	All	0	3

The worst 5 of 20 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	255	GLU	CD-OE1	-7.42	1.17	1.25
1	A	205	GLU	CD-OE1	-6.15	1.18	1.25
1	A	21	VAL	C-O	-5.92	1.12	1.23
1	A	195	SER	CB-OG	-5.86	1.34	1.42
1	B	195	SER	CB-OG	-5.68	1.34	1.42

The worst 5 of 23 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	3	LYS	CD-CE-NZ	8.36	130.92	111.70
1	A	4	LEU	CA-CB-CG	6.22	129.62	115.30
1	B	51	ASP	CB-CG-OD2	-6.22	112.70	118.30
1	A	246	PHE	CB-CG-CD1	6.06	125.04	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	246	PHE	CB-CG-CD2	-5.95	116.64	120.80

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	129[A]	SER	Mainchain
1	A	129[B]	SER	Mainchain
1	B	129	SER	Peptide

## 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	2415	0	2431	19	0
1	B	2402	0	2414	43	0
2	A	21	0	13	0	0
2	B	42	0	26	7	0
3	A	54	0	72	4	0
3	B	36	0	47	10	0
4	A	31	0	19	1	0
4	B	31	0	19	0	0
5	A	7	0	0	0	0
6	A	374	0	0	7	0
6	B	354	0	0	10	0
All	All	5767	0	5041	66	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 7.

The worst 5 of 66 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:131:PRO:HG3	2:B:401[B]:W7G:CAA	1.74	1.17
1:B:131:PRO:HG3	2:B:401[B]:W7G:H14	1.22	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:132:ASN:HB2	6:B:839:HOH:O	1.61	0.97
1:B:131:PRO:CG	2:B:401[B]:W7G:H14	2.05	0.86
1:B:9:HIS:CE1	3:B:407:GOL:H32	2.11	0.84

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	316/313 (101%)	307 (97%)	9 (3%)	0	100	100
1	B	314/313 (100%)	308 (98%)	6 (2%)	0	100	100
All	All	630/626 (101%)	615 (98%)	15 (2%)	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	263/258 (102%)	260 (99%)	3 (1%)	73	53
1	B	261/258 (101%)	257 (98%)	4 (2%)	65	37
All	All	524/516 (102%)	517 (99%)	7 (1%)	71	44

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	81	LYS
1	B	166	PRO
1	B	130[A]	CYS
1	A	5	ASN
1	B	130[B]	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	215	GLN
1	A	234	ASN
1	B	215	GLN
1	A	152	GLN
1	B	152	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](#)

21 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	W7G	B	401[A]	-	18,22,22	2.16	6 (33%)	17,30,30	2.41	5 (29%)
3	GOL	B	403	-	5,5,5	1.26	1 (20%)	5,5,5	1.45	1 (20%)
3	GOL	B	405	-	5,5,5	0.68	0	5,5,5	0.31	0
3	GOL	A	409	-	5,5,5	0.46	0	5,5,5	0.82	0
3	GOL	A	403	-	5,5,5	0.66	0	5,5,5	0.77	0
3	GOL	A	402	-	5,5,5	0.58	0	5,5,5	2.18	4 (80%)
3	GOL	A	405	-	5,5,5	0.42	0	5,5,5	0.67	0
2	W7G	A	401	-	18,22,22	2.13	7 (38%)	17,30,30	3.39	6 (35%)
3	GOL	A	406	-	5,5,5	0.29	0	5,5,5	0.45	0
5	NCO	A	411	-	6,6,6	3.28	6 (100%)	-		
3	GOL	B	404	-	5,5,5	0.39	0	5,5,5	0.43	0
3	GOL	A	408	-	5,5,5	0.38	0	5,5,5	0.57	0
3	GOL	A	407[B]	-	5,5,5	0.66	0	5,5,5	0.44	0
3	GOL	A	407[A]	-	5,5,5	0.67	0	5,5,5	0.59	0
4	FMN	B	408	-	31,33,33	1.47	6 (19%)	40,50,50	2.23	9 (22%)
4	FMN	A	410	-	31,33,33	1.94	7 (22%)	40,50,50	1.94	7 (17%)
3	GOL	B	407	-	5,5,5	0.78	0	5,5,5	1.70	1 (20%)
3	GOL	B	402	-	5,5,5	0.46	0	5,5,5	0.46	0
3	GOL	A	404	-	5,5,5	0.88	0	5,5,5	0.86	0
2	W7G	B	401[B]	-	18,22,22	2.05	5 (27%)	17,30,30	3.00	7 (41%)
3	GOL	B	406	-	5,5,5	0.79	0	5,5,5	1.27	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
2	W7G	B	401[A]	-	-	0/7/11/11	0/2/2/2
3	GOL	B	403	-	-	0/4/4/4	-
3	GOL	B	405	-	-	0/4/4/4	-
3	GOL	A	409	-	-	3/4/4/4	-
3	GOL	A	403	-	-	3/4/4/4	-
3	GOL	A	402	-	-	0/4/4/4	-
3	GOL	A	405	-	-	1/4/4/4	-
2	W7G	A	401	-	-	0/7/11/11	0/2/2/2
3	GOL	A	406	-	-	2/4/4/4	-
3	GOL	B	404	-	-	2/4/4/4	-
3	GOL	A	408	-	-	0/4/4/4	-
3	GOL	A	407[B]	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	407[A]	-	-	0/4/4/4	-
4	FMN	B	408	-	-	1/18/18/18	0/3/3/3
4	FMN	A	410	-	-	1/18/18/18	0/3/3/3
3	GOL	B	407	-	-	4/4/4/4	-
3	GOL	B	402	-	-	2/4/4/4	-
3	GOL	A	404	-	-	0/4/4/4	-
2	W7G	B	401[B]	-	-	0/7/11/11	0/2/2/2
3	GOL	B	406	-	-	2/4/4/4	-

The worst 5 of 38 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	410	FMN	C4A-C10	7.06	1.45	1.38
2	A	401	W7G	CAU-NAM	5.30	1.42	1.33
2	B	401[A]	W7G	CAU-NAM	5.06	1.41	1.33
2	B	401[B]	W7G	CAJ-CAP	-4.82	1.42	1.51
2	B	401[B]	W7G	CAU-NAM	4.28	1.40	1.33

The worst 5 of 40 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	W7G	CAU-NAM-CAT	10.75	124.22	115.14
4	B	408	FMN	C4-N3-C2	8.78	122.56	115.14
4	A	410	FMN	C4-N3-C2	7.68	121.63	115.14
2	B	401[B]	W7G	CAA-OAN-CAQ	-7.30	106.52	117.53
2	B	401[A]	W7G	CAU-NAM-CAT	7.21	121.23	115.14

There are no chirality outliers.

5 of 21 torsion outliers are listed below:

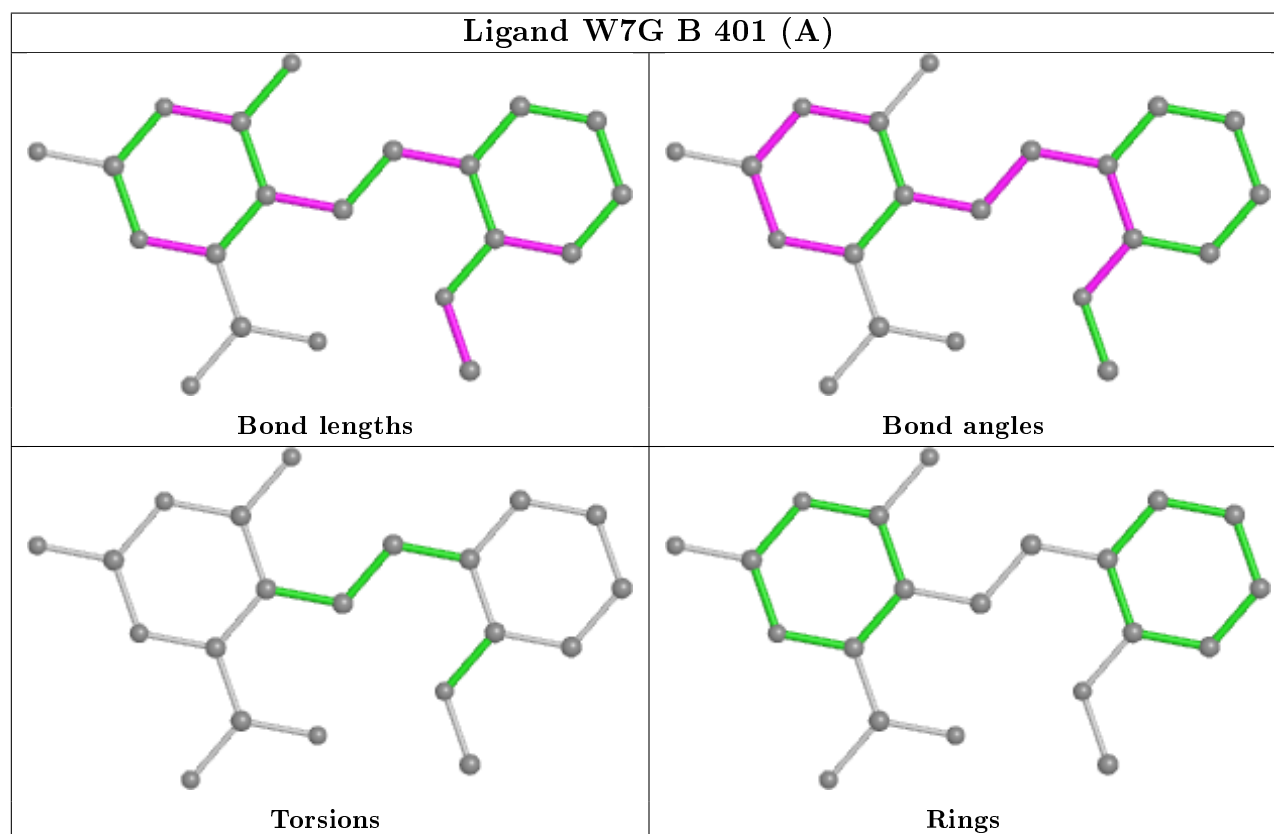
Mol	Chain	Res	Type	Atoms
3	A	409	GOL	O1-C1-C2-C3
3	A	406	GOL	O1-C1-C2-C3
3	B	404	GOL	C1-C2-C3-O3
3	A	403	GOL	O1-C1-C2-C3
3	B	407	GOL	O1-C1-C2-C3

There are no ring outliers.

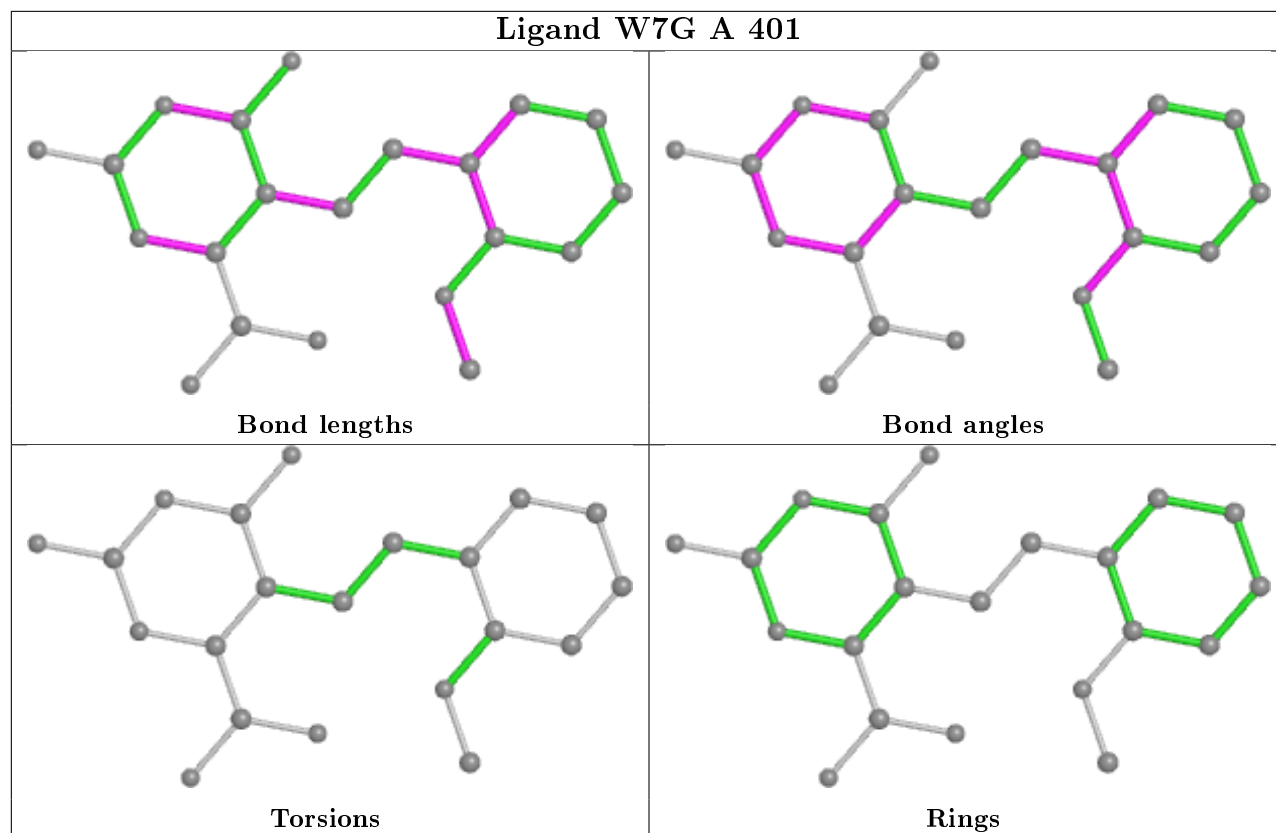
6 monomers are involved in 22 short contacts:

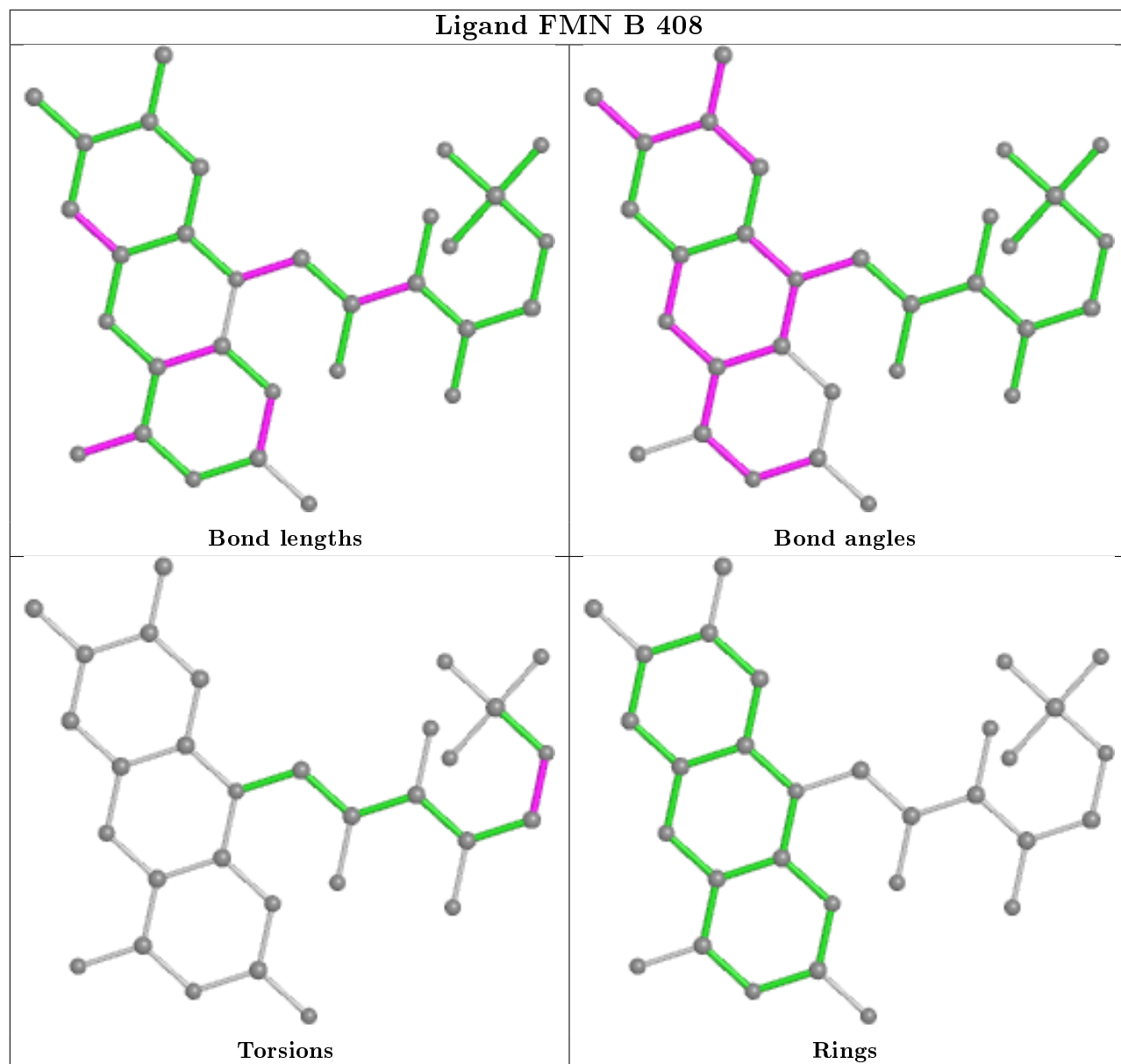
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	409	GOL	3	0
3	A	403	GOL	1	0
4	A	410	FMN	1	0
3	B	407	GOL	9	0
3	B	402	GOL	1	0
2	B	401[B]	W7G	7	0

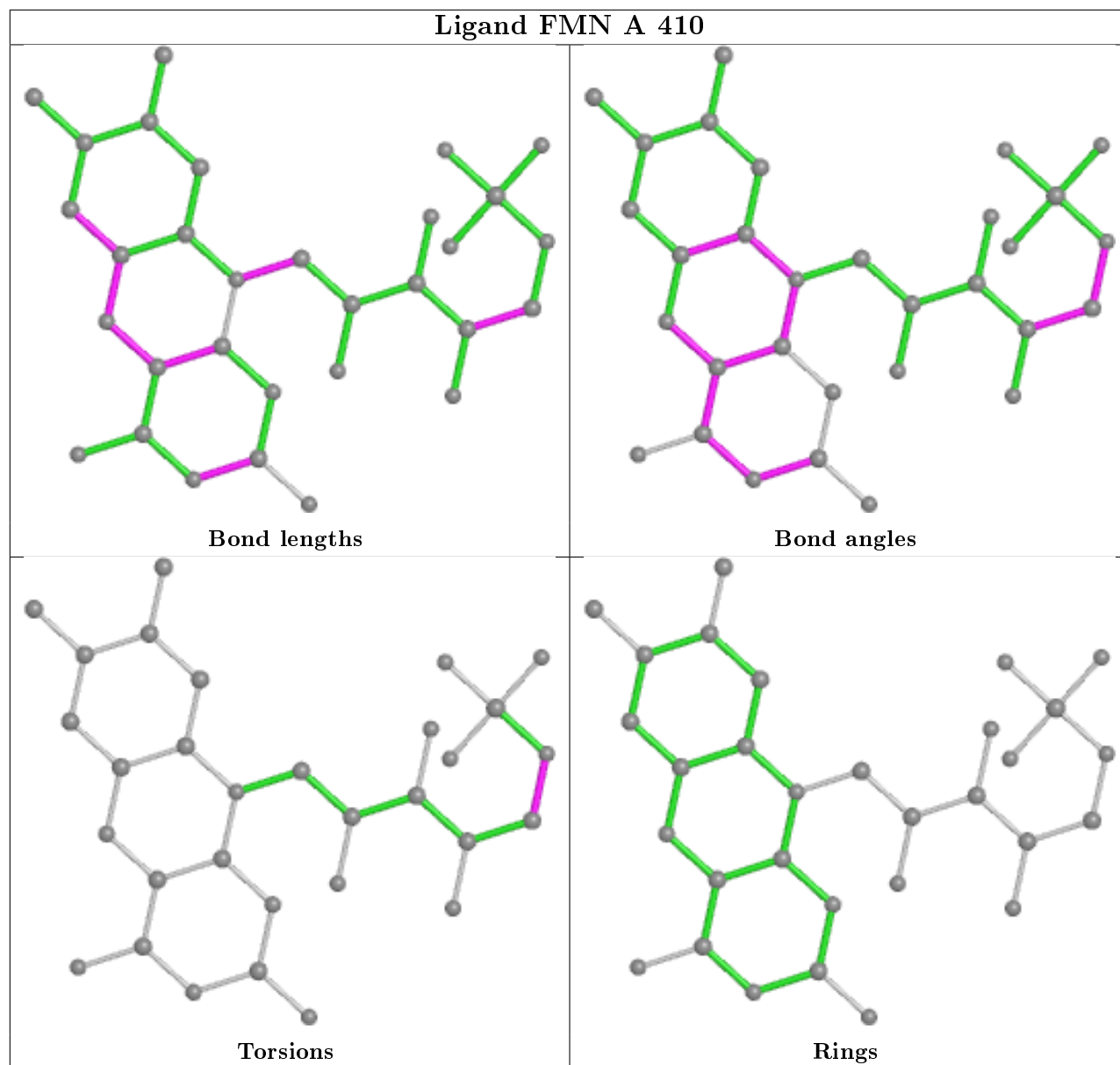
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

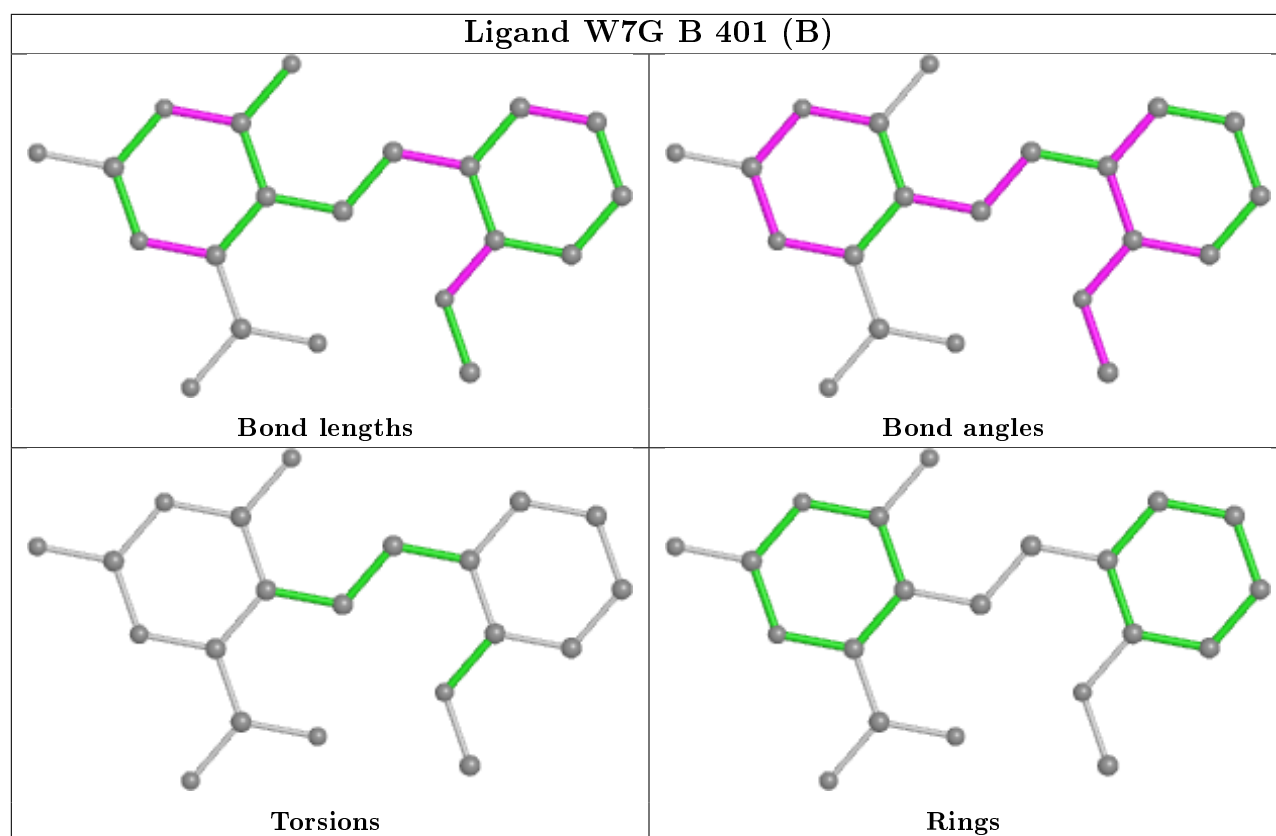












## 5.7 Other polymers [i](#)

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	313/313 (100%)	-0.45	1 (0%) 94 95	6, 10, 20, 41	0
1	B	313/313 (100%)	-0.29	5 (1%) 72 77	6, 11, 25, 46	0
All	All	626/626 (100%)	-0.37	6 (0%) 82 86	6, 11, 24, 46	0

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	131	PRO	5.0
1	B	135	GLY	3.5
1	A	312	GLU	2.8
1	B	312	GLU	2.4
1	B	130[A]	CYS	2.4

### 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 carbohydrates in this entry.

### 6.4 Ligands [i](#)

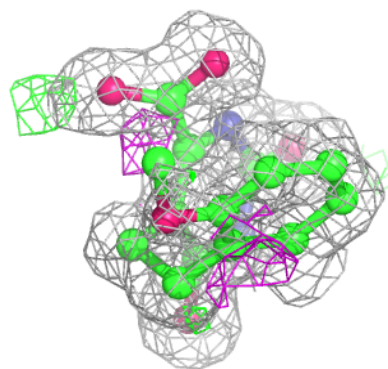
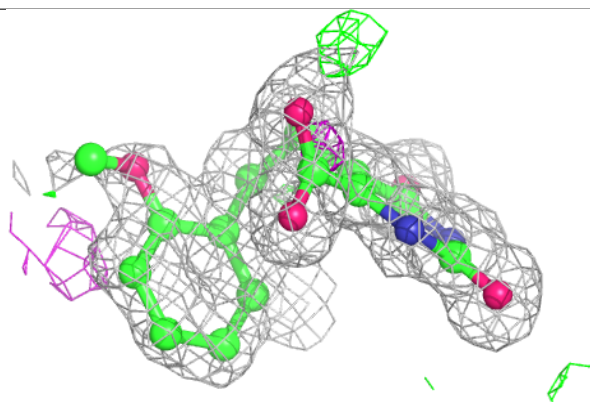
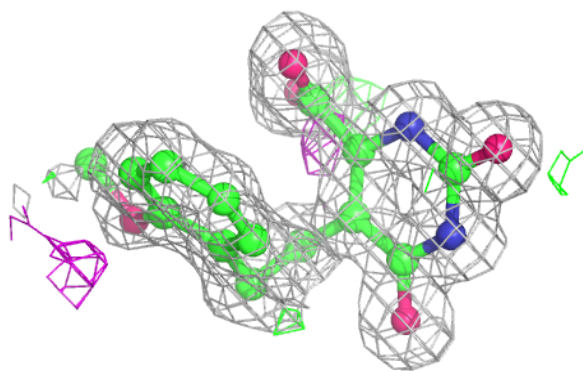
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 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
3	GOL	B	404	6/6	0.78	0.18	33,39,42,44	0
3	GOL	A	406	6/6	0.79	0.16	55,55,59,62	0
3	GOL	B	406	6/6	0.80	0.20	31,35,36,39	0
3	GOL	A	409	6/6	0.81	0.18	36,40,43,44	0
3	GOL	B	402	6/6	0.82	0.16	28,44,46,50	0
3	GOL	A	402	6/6	0.82	0.14	29,37,39,39	0
3	GOL	B	403	6/6	0.84	0.15	27,32,37,42	0
3	GOL	A	403	6/6	0.86	0.19	23,28,32,39	0
3	GOL	B	407	6/6	0.89	0.33	25,35,38,41	0
3	GOL	A	404	6/6	0.90	0.12	27,29,32,32	0
3	GOL	A	405	6/6	0.92	0.12	23,26,28,31	0
3	GOL	B	405	6/6	0.95	0.07	16,17,19,20	0
2	W7G	B	401[B]	21/21	0.96	0.08	9,10,16,20	21
2	W7G	B	401[A]	21/21	0.96	0.08	8,10,12,13	21
2	W7G	A	401	21/21	0.97	0.06	8,9,11,12	0
3	GOL	A	408	6/6	0.97	0.05	15,17,19,21	0
4	FMN	A	410	31/31	0.98	0.07	6,7,9,14	0
3	GOL	A	407[A]	6/6	0.98	0.06	12,15,16,17	6
3	GOL	A	407[B]	6/6	0.98	0.06	12,14,15,16	6
5	NCO	A	411	7/7	0.99	0.07	10,10,12,12	0
4	FMN	B	408	31/31	0.99	0.06	6,7,9,12	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.

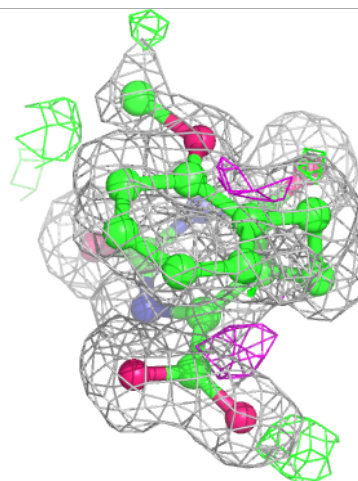
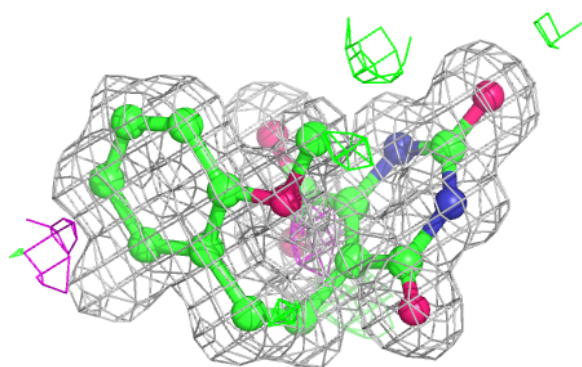
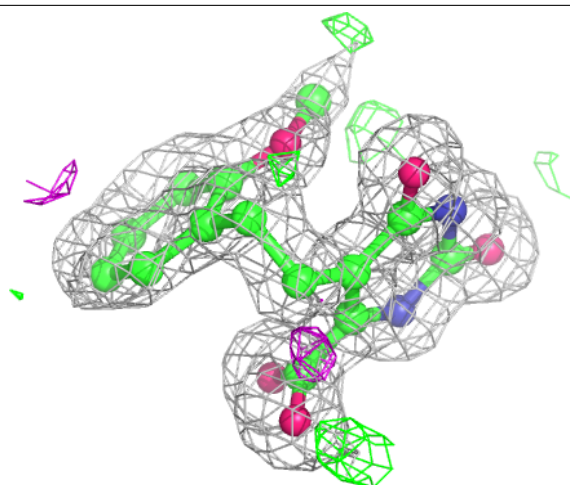
**Electron density around W7G B 401 (B):**

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



**Electron density around W7G B 401 (A):**

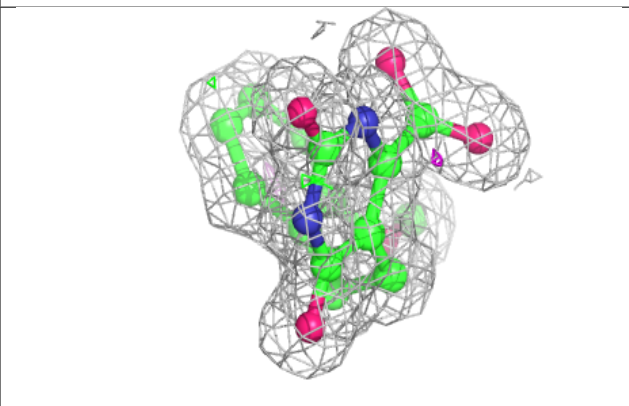
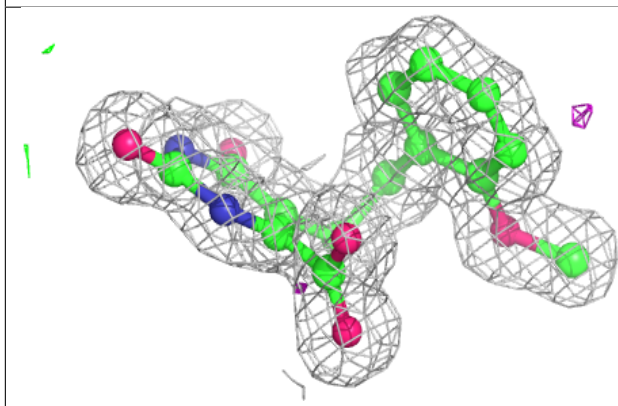
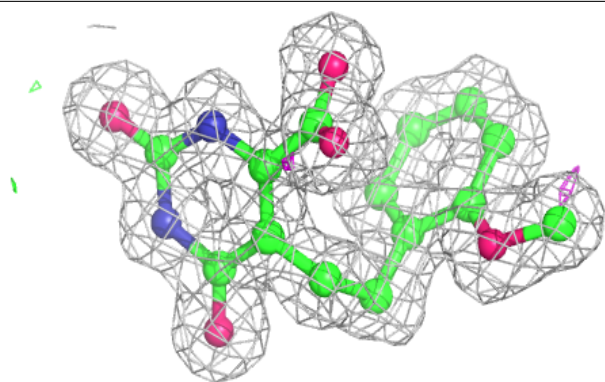
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



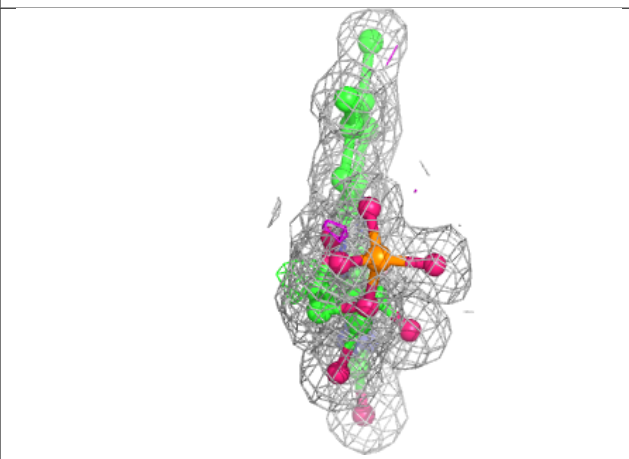
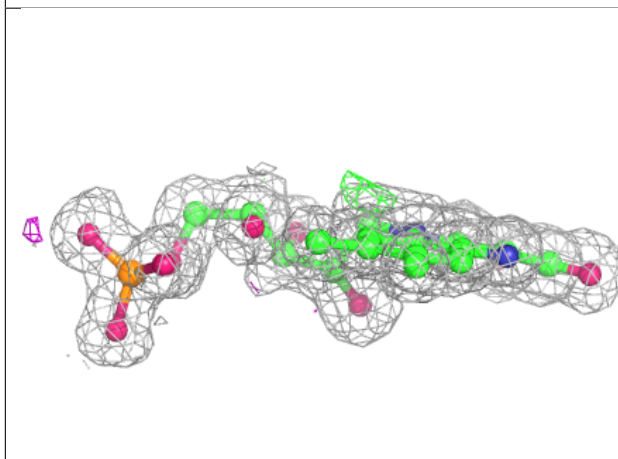
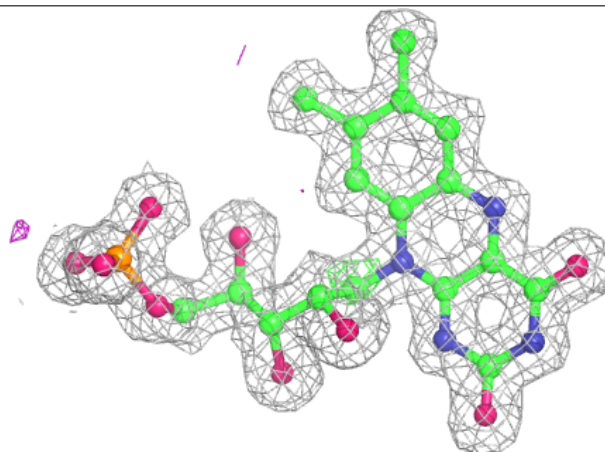


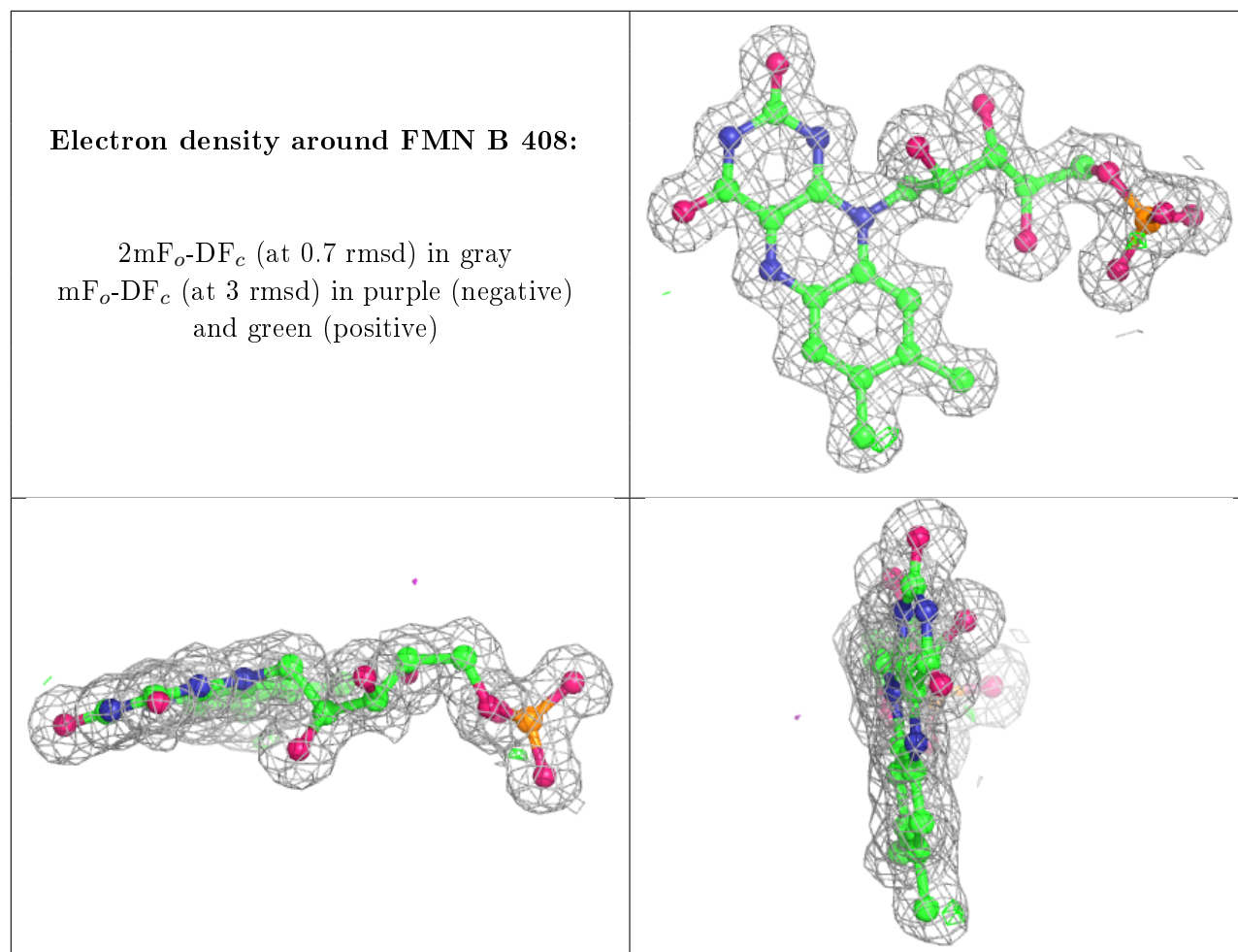
**Electron density around W7G A 401:**

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

**Electron density around FMN A 410:**

$2mF_o-DF_c$  (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.