



# wwPDB X-ray Structure Validation Summary Report

Jan 23, 2021 – 04:19 PM EST


PDB ID : 1YRH  
Title : Crystal Structure Of Trp Repressor Binding Protein Wrba in complex with FMN  
Authors : Gorman, J.; Shapiro, L.; Burley, S.K.; New York SGX Research Center for Structural Genomics (NYSGXRC)  
Deposited on : 2005-02-03  
Resolution : 3.11 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 : 2.16  
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) : 2.16

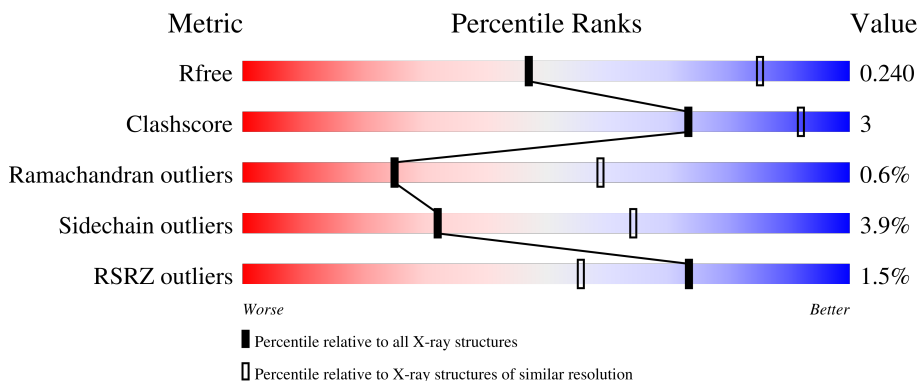
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.11 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.



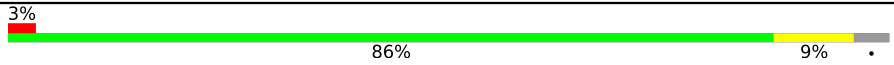


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	1292 (3.14-3.10)
Clashscore	141614	1389 (3.14-3.10)
Ramachandran outliers	138981	1337 (3.14-3.10)
Sidechain outliers	138945	1337 (3.14-3.10)
RSRZ outliers	127900	1260 (3.14-3.10)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	211	 89% 6% 5%
1	B	211	 86% 8% 5%
1	C	211	 84% 10% 5%
1	D	211	 86% 9% 5%
1	E	211	 83% 9% 6%

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Mol	Chain	Length	Quality of chain
1	F	211	
1	G	211	
1	H	211	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 12386 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 trp repressor binding protein WrbA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	200	1495	935	260	294	6	0	0	0
1	B	200	1495	935	260	294	6	0	0	0
1	C	201	1500	938	261	295	6	0	0	0
1	D	201	1500	938	261	295	6	0	0	0
1	E	198	1485	930	258	291	6	0	0	0
1	F	202	1507	942	262	297	6	0	0	0
1	G	200	1495	935	260	294	6	0	0	0
1	H	201	1500	938	261	295	6	0	0	0

There are 144 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	cloning artifact	UNP Q9RYU4
A	0	SER	-	cloning artifact	UNP Q9RYU4
A	1	LEU	-	cloning artifact	UNP Q9RYU4
A	21	MSE	MET	modified residue	UNP Q9RYU4
A	63	MSE	MET	modified residue	UNP Q9RYU4
A	95	MSE	MET	modified residue	UNP Q9RYU4
A	119	MSE	MET	modified residue	UNP Q9RYU4
A	138	MSE	MET	modified residue	UNP Q9RYU4
A	141	MSE	MET	modified residue	UNP Q9RYU4
A	200	GLU	-	cloning artifact	UNP Q9RYU4
A	202	GLY	-	cloning artifact	UNP Q9RYU4
A	203	SER	-	cloning artifact	UNP Q9RYU4
A	204	HIS	-	cloning artifact	UNP Q9RYU4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	205	HIS	-	cloning artifact	UNP Q9RYU4
A	206	HIS	-	cloning artifact	UNP Q9RYU4
A	207	HIS	-	cloning artifact	UNP Q9RYU4
A	208	HIS	-	cloning artifact	UNP Q9RYU4
A	209	HIS	-	cloning artifact	UNP Q9RYU4
B	-1	MET	-	cloning artifact	UNP Q9RYU4
B	0	SER	-	cloning artifact	UNP Q9RYU4
B	1	LEU	-	cloning artifact	UNP Q9RYU4
B	21	MSE	MET	modified residue	UNP Q9RYU4
B	63	MSE	MET	modified residue	UNP Q9RYU4
B	95	MSE	MET	modified residue	UNP Q9RYU4
B	119	MSE	MET	modified residue	UNP Q9RYU4
B	138	MSE	MET	modified residue	UNP Q9RYU4
B	141	MSE	MET	modified residue	UNP Q9RYU4
B	200	GLU	-	cloning artifact	UNP Q9RYU4
B	202	GLY	-	cloning artifact	UNP Q9RYU4
B	203	SER	-	cloning artifact	UNP Q9RYU4
B	204	HIS	-	cloning artifact	UNP Q9RYU4
B	205	HIS	-	cloning artifact	UNP Q9RYU4
B	206	HIS	-	cloning artifact	UNP Q9RYU4
B	207	HIS	-	cloning artifact	UNP Q9RYU4
B	208	HIS	-	cloning artifact	UNP Q9RYU4
B	209	HIS	-	cloning artifact	UNP Q9RYU4
C	-1	MET	-	cloning artifact	UNP Q9RYU4
C	0	SER	-	cloning artifact	UNP Q9RYU4
C	1	LEU	-	cloning artifact	UNP Q9RYU4
C	21	MSE	MET	modified residue	UNP Q9RYU4
C	63	MSE	MET	modified residue	UNP Q9RYU4
C	95	MSE	MET	modified residue	UNP Q9RYU4
C	119	MSE	MET	modified residue	UNP Q9RYU4
C	138	MSE	MET	modified residue	UNP Q9RYU4
C	141	MSE	MET	modified residue	UNP Q9RYU4
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C	203	SER	-	cloning artifact	UNP Q9RYU4
C	204	HIS	-	cloning artifact	UNP Q9RYU4
C	205	HIS	-	cloning artifact	UNP Q9RYU4
C	206	HIS	-	cloning artifact	UNP Q9RYU4
C	207	HIS	-	cloning artifact	UNP Q9RYU4
C	208	HIS	-	cloning artifact	UNP Q9RYU4
C	209	HIS	-	cloning artifact	UNP Q9RYU4
D	-1	MET	-	cloning artifact	UNP Q9RYU4

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Chain	Residue	Modelled	Actual	Comment	Reference
D	0	SER	-	cloning artifact	UNP Q9RYU4
D	1	LEU	-	cloning artifact	UNP Q9RYU4
D	21	MSE	MET	modified residue	UNP Q9RYU4
D	63	MSE	MET	modified residue	UNP Q9RYU4
D	95	MSE	MET	modified residue	UNP Q9RYU4
D	119	MSE	MET	modified residue	UNP Q9RYU4
D	138	MSE	MET	modified residue	UNP Q9RYU4
D	141	MSE	MET	modified residue	UNP Q9RYU4
D	200	GLU	-	cloning artifact	UNP Q9RYU4
D	202	GLY	-	cloning artifact	UNP Q9RYU4
D	203	SER	-	cloning artifact	UNP Q9RYU4
D	204	HIS	-	cloning artifact	UNP Q9RYU4
D	205	HIS	-	cloning artifact	UNP Q9RYU4
D	206	HIS	-	cloning artifact	UNP Q9RYU4
D	207	HIS	-	cloning artifact	UNP Q9RYU4
D	208	HIS	-	cloning artifact	UNP Q9RYU4
D	209	HIS	-	cloning artifact	UNP Q9RYU4
E	-1	MET	-	cloning artifact	UNP Q9RYU4
E	0	SER	-	cloning artifact	UNP Q9RYU4
E	1	LEU	-	cloning artifact	UNP Q9RYU4
E	21	MSE	MET	modified residue	UNP Q9RYU4
E	63	MSE	MET	modified residue	UNP Q9RYU4
E	95	MSE	MET	modified residue	UNP Q9RYU4
E	119	MSE	MET	modified residue	UNP Q9RYU4
E	138	MSE	MET	modified residue	UNP Q9RYU4
E	141	MSE	MET	modified residue	UNP Q9RYU4
E	200	GLU	-	cloning artifact	UNP Q9RYU4
E	202	GLY	-	cloning artifact	UNP Q9RYU4
E	203	SER	-	cloning artifact	UNP Q9RYU4
E	204	HIS	-	cloning artifact	UNP Q9RYU4
E	205	HIS	-	cloning artifact	UNP Q9RYU4
E	206	HIS	-	cloning artifact	UNP Q9RYU4
E	207	HIS	-	cloning artifact	UNP Q9RYU4
E	208	HIS	-	cloning artifact	UNP Q9RYU4
E	209	HIS	-	cloning artifact	UNP Q9RYU4
F	-1	MET	-	cloning artifact	UNP Q9RYU4
F	0	SER	-	cloning artifact	UNP Q9RYU4
F	1	LEU	-	cloning artifact	UNP Q9RYU4
F	21	MSE	MET	modified residue	UNP Q9RYU4
F	63	MSE	MET	modified residue	UNP Q9RYU4
F	95	MSE	MET	modified residue	UNP Q9RYU4
F	119	MSE	MET	modified residue	UNP Q9RYU4

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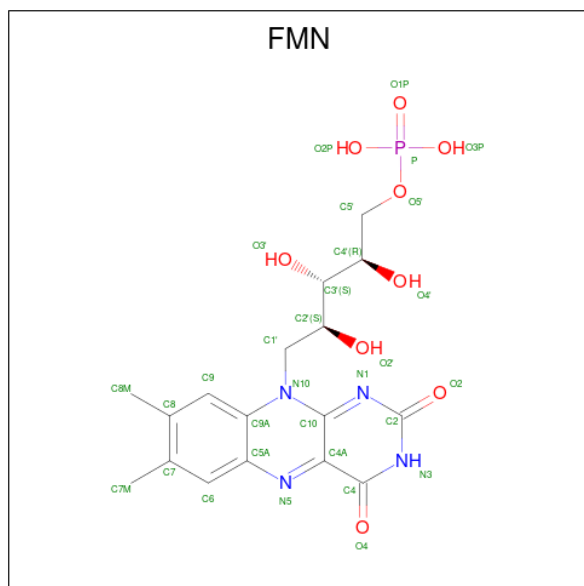
Chain	Residue	Modelled	Actual	Comment	Reference
F	138	MSE	MET	modified residue	UNP Q9RYU4
F	141	MSE	MET	modified residue	UNP Q9RYU4
F	200	GLU	-	cloning artifact	UNP Q9RYU4
F	202	GLY	-	cloning artifact	UNP Q9RYU4
F	203	SER	-	cloning artifact	UNP Q9RYU4
F	204	HIS	-	cloning artifact	UNP Q9RYU4
F	205	HIS	-	cloning artifact	UNP Q9RYU4
F	206	HIS	-	cloning artifact	UNP Q9RYU4
F	207	HIS	-	cloning artifact	UNP Q9RYU4
F	208	HIS	-	cloning artifact	UNP Q9RYU4
F	209	HIS	-	cloning artifact	UNP Q9RYU4
G	-1	MET	-	cloning artifact	UNP Q9RYU4
G	0	SER	-	cloning artifact	UNP Q9RYU4
G	1	LEU	-	cloning artifact	UNP Q9RYU4
G	21	MSE	MET	modified residue	UNP Q9RYU4
G	63	MSE	MET	modified residue	UNP Q9RYU4
G	95	MSE	MET	modified residue	UNP Q9RYU4
G	119	MSE	MET	modified residue	UNP Q9RYU4
G	138	MSE	MET	modified residue	UNP Q9RYU4
G	141	MSE	MET	modified residue	UNP Q9RYU4
G	200	GLU	-	cloning artifact	UNP Q9RYU4
G	202	GLY	-	cloning artifact	UNP Q9RYU4
G	203	SER	-	cloning artifact	UNP Q9RYU4
G	204	HIS	-	cloning artifact	UNP Q9RYU4
G	205	HIS	-	cloning artifact	UNP Q9RYU4
G	206	HIS	-	cloning artifact	UNP Q9RYU4
G	207	HIS	-	cloning artifact	UNP Q9RYU4
G	208	HIS	-	cloning artifact	UNP Q9RYU4
G	209	HIS	-	cloning artifact	UNP Q9RYU4
H	-1	MET	-	cloning artifact	UNP Q9RYU4
H	0	SER	-	cloning artifact	UNP Q9RYU4
H	1	LEU	-	cloning artifact	UNP Q9RYU4
H	21	MSE	MET	modified residue	UNP Q9RYU4
H	63	MSE	MET	modified residue	UNP Q9RYU4
H	95	MSE	MET	modified residue	UNP Q9RYU4
H	119	MSE	MET	modified residue	UNP Q9RYU4
H	138	MSE	MET	modified residue	UNP Q9RYU4
H	141	MSE	MET	modified residue	UNP Q9RYU4
H	200	GLU	-	cloning artifact	UNP Q9RYU4
H	202	GLY	-	cloning artifact	UNP Q9RYU4
H	203	SER	-	cloning artifact	UNP Q9RYU4
H	204	HIS	-	cloning artifact	UNP Q9RYU4

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Chain	Residue	Modelled	Actual	Comment	Reference
H	205	HIS	-	cloning artifact	UNP Q9RYU4
H	206	HIS	-	cloning artifact	UNP Q9RYU4
H	207	HIS	-	cloning artifact	UNP Q9RYU4
H	208	HIS	-	cloning artifact	UNP Q9RYU4
H	209	HIS	-	cloning artifact	UNP Q9RYU4

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: C<sub>17</sub>H<sub>21</sub>N<sub>4</sub>O<sub>9</sub>P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	P		
2	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	B	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	C	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	D	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	E	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	F	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	G	1	Total	C	N	O	P	0	0
			31	17	4	9	1		
2	H	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 3 is water.



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	29	Total O 29 29	0	0
3	B	21	Total O 21 21	0	0
3	C	23	Total O 23 23	0	0
3	D	24	Total O 24 24	0	0
3	E	12	Total O 12 12	0	0
3	F	21	Total O 21 21	0	0
3	G	12	Total O 12 12	0	0
3	H	19	Total O 19 19	0	0

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


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: trp repressor binding protein WrbA

Chain A: 




- Molecule 1: trp repressor binding protein WrbA

Chain B: 




- Molecule 1: trp repressor binding protein WrbA

Chain C: 




- Molecule 1: trp repressor binding protein WrbA

Chain D: 



- Molecule 1: trp repressor binding protein WrbA

Chain E: 



- Molecule 1: trp repressor binding protein WrbA



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	122.33Å 122.33Å 208.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 3.11 20.02 – 3.11	Depositor EDS
% Data completeness (in resolution range)	97.0 (20.00-3.11) 97.1 (20.02-3.11)	Depositor EDS
$R_{merge}$	0.17	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.65 (at 3.09Å)	Xtrriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.198 , 0.233 0.203 , 0.240	Depositor DCC
$R_{free}$ test set	1632 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	36.2	Xtrriage
Anisotropy	0.058	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 26.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.032 for -h,-k,l	Xtrriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	12386	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.31% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality [i](#)

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

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

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	0.56	0/1518	0.60	0/2051
1	B	0.55	0/1518	0.61	0/2051
1	C	0.55	0/1523	0.57	0/2059
1	D	0.55	0/1523	0.59	0/2059
1	E	0.53	0/1508	0.59	0/2038
1	F	0.51	0/1530	0.55	0/2069
1	G	0.53	0/1518	0.56	0/2051
1	H	0.52	0/1523	0.62	1/2059 (0.0%)
All	All	0.54	0/12161	0.59	1/16437 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	30	ARG	NE-CZ-NH1	5.35	122.97	120.30

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	1495	0	1466	6	0
1	B	1495	0	1466	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1500	0	1470	11	0
1	D	1500	0	1470	14	0
1	E	1485	0	1458	15	0
1	F	1507	0	1477	14	0
1	G	1495	0	1466	11	0
1	H	1500	0	1470	14	0
2	A	31	0	19	0	0
2	B	31	0	19	0	0
2	C	31	0	19	0	0
2	D	31	0	19	0	0
2	E	31	0	19	0	0
2	F	31	0	19	1	0
2	G	31	0	19	0	0
2	H	31	0	19	2	0
3	A	29	0	0	0	0
3	B	21	0	0	0	0
3	C	23	0	0	1	0
3	D	24	0	0	1	0
3	E	12	0	0	2	0
3	F	21	0	0	1	0
3	G	12	0	0	1	0
3	H	19	0	0	1	0
All	All	12386	0	11895	80	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 3.

The worst 5 of 80 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:134:GLN:HE22	1:D:134:GLN:HE22	1.26	0.81
1:F:134:GLN:HE22	1:H:134:GLN:HE22	1.34	0.75
1:F:134:GLN:O	1:F:138:MSE:HG2	1.88	0.73
1:E:134:GLN:HE21	1:F:134:GLN:NE2	1.87	0.71
1:G:115:THR:HG21	1:G:195:THR:OG1	1.93	0.69

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	198/211 (94%)	191 (96%)	7 (4%)	0	100	100
1	B	198/211 (94%)	192 (97%)	6 (3%)	0	100	100
1	C	199/211 (94%)	189 (95%)	8 (4%)	2 (1%)	15	48
1	D	199/211 (94%)	192 (96%)	7 (4%)	0	100	100
1	E	196/211 (93%)	181 (92%)	12 (6%)	3 (2%)	10	38
1	F	200/211 (95%)	178 (89%)	21 (10%)	1 (0%)	29	63
1	G	198/211 (94%)	183 (92%)	13 (7%)	2 (1%)	15	48
1	H	199/211 (94%)	192 (96%)	5 (2%)	2 (1%)	15	48
All	All	1587/1688 (94%)	1498 (94%)	79 (5%)	10 (1%)	25	59

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	109	GLY
1	E	88	PHE
1	F	100	ASP
1	H	109	GLY
1	C	155	GLU

### 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	152/156 (97%)	147 (97%)	5 (3%)	38	68

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	152/156 (97%)	143 (94%)	9 (6%)	19	50
1	C	152/156 (97%)	146 (96%)	6 (4%)	32	64
1	D	152/156 (97%)	146 (96%)	6 (4%)	32	64
1	E	151/156 (97%)	148 (98%)	3 (2%)	55	79
1	F	153/156 (98%)	149 (97%)	4 (3%)	46	74
1	G	152/156 (97%)	142 (93%)	10 (7%)	16	46
1	H	152/156 (97%)	147 (97%)	5 (3%)	38	68
All	All	1216/1248 (97%)	1168 (96%)	48 (4%)	32	64

5 of 48 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	64	LYS
1	E	17	THR
1	H	54	ASP
1	D	87	ARG
1	D	155	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 21 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	191	GLN
1	E	191	GLN
1	H	53	GLN
1	D	134	GLN
1	H	123	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

8 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	FMN	F	407	-	31,33,33	1.51	5 (16%)	40,50,50	1.59	8 (20%)
2	FMN	D	401	-	31,33,33	1.47	5 (16%)	40,50,50	1.79	7 (17%)
2	FMN	A	404	-	31,33,33	1.54	5 (16%)	40,50,50	1.50	4 (10%)
2	FMN	G	406	-	31,33,33	1.49	4 (12%)	40,50,50	1.69	7 (17%)
2	FMN	E	408	-	31,33,33	1.63	5 (16%)	40,50,50	1.97	7 (17%)
2	FMN	H	405	-	31,33,33	1.58	4 (12%)	40,50,50	1.62	5 (12%)
2	FMN	C	402	-	31,33,33	1.40	4 (12%)	40,50,50	1.67	6 (15%)
2	FMN	B	403	-	31,33,33	1.44	4 (12%)	40,50,50	1.76	5 (12%)

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	FMN	F	407	-	-	7/18/18/18	0/3/3/3
2	FMN	D	401	-	-	4/18/18/18	0/3/3/3
2	FMN	A	404	-	-	5/18/18/18	0/3/3/3
2	FMN	G	406	-	-	7/18/18/18	0/3/3/3
2	FMN	E	408	-	-	0/18/18/18	0/3/3/3
2	FMN	H	405	-	-	5/18/18/18	0/3/3/3
2	FMN	C	402	-	-	9/18/18/18	0/3/3/3
2	FMN	B	403	-	-	4/18/18/18	0/3/3/3

The worst 5 of 36 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	405	FMN	C10-N1	4.79	1.39	1.33
2	G	406	FMN	C10-N1	4.52	1.39	1.33
2	B	403	FMN	C10-N1	4.30	1.38	1.33
2	E	408	FMN	C10-N1	4.26	1.38	1.33
2	F	407	FMN	C10-N1	4.25	1.38	1.33

The worst 5 of 49 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	403	FMN	C4-N3-C2	6.43	120.57	115.14
2	E	408	FMN	C4-N3-C2	6.39	120.54	115.14
2	D	401	FMN	C4-N3-C2	6.31	120.47	115.14
2	H	405	FMN	C4-N3-C2	6.24	120.41	115.14
2	G	406	FMN	C4-N3-C2	5.73	119.98	115.14

There are no chirality outliers.

5 of 41 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	F	407	FMN	C2'-C3'-C4'-O4'
2	F	407	FMN	C2'-C3'-C4'-C5'
2	F	407	FMN	O3'-C3'-C4'-C5'
2	F	407	FMN	C5'-O5'-P-O3P
2	D	401	FMN	C2'-C3'-C4'-O4'

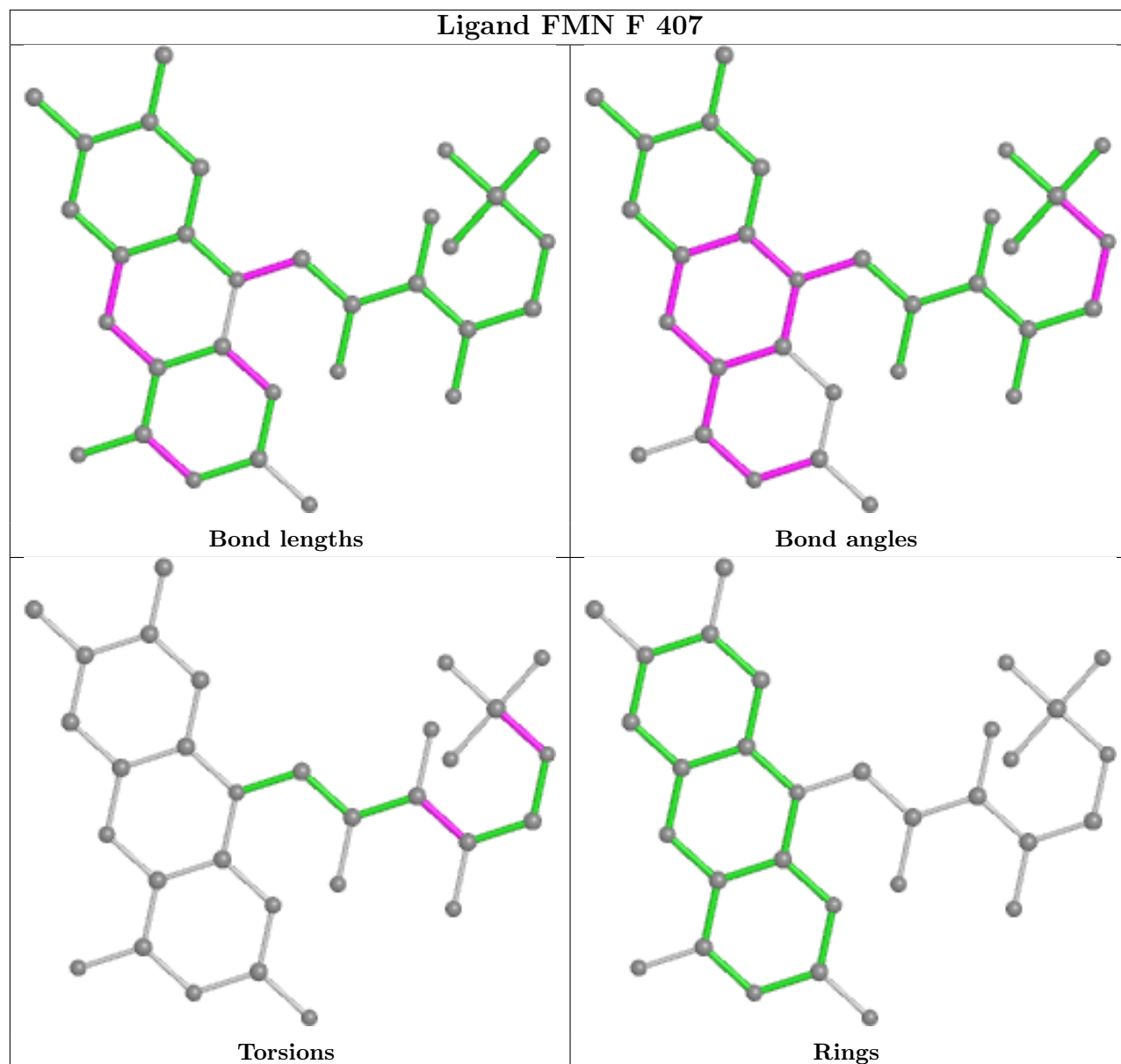
There are no ring outliers.

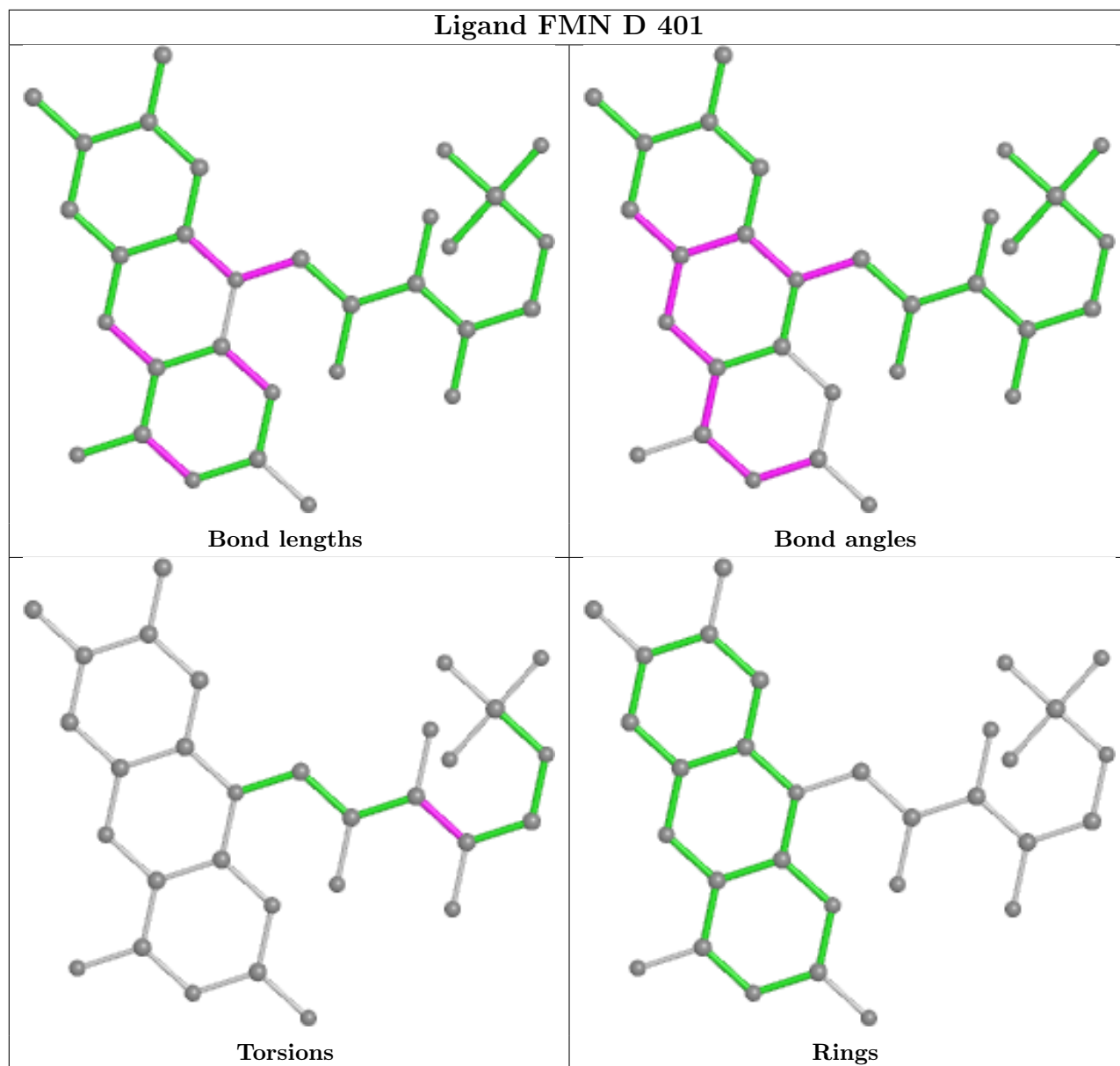
2 monomers are involved in 3 short contacts:

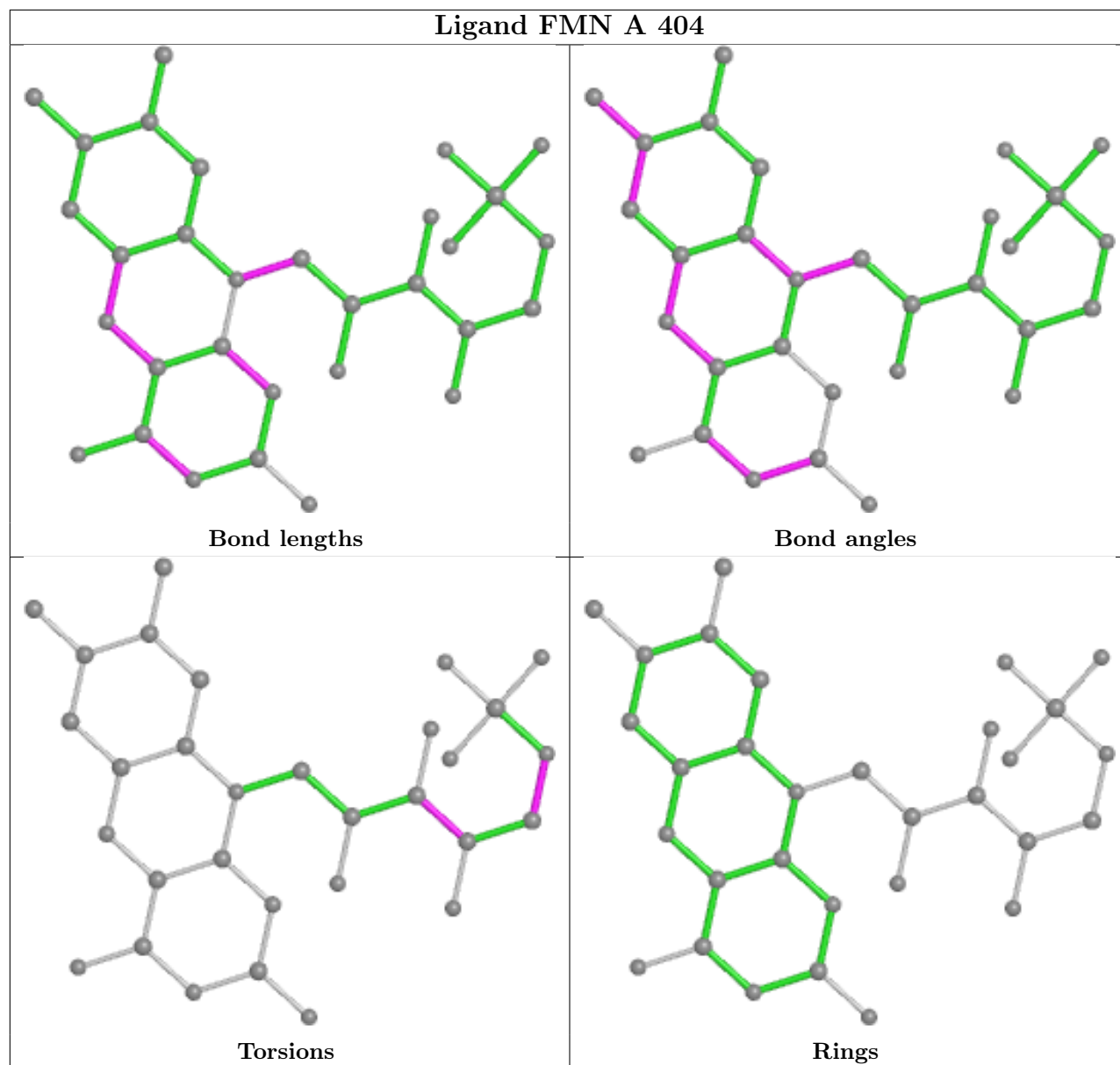
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	407	FMN	1	0
2	H	405	FMN	2	0

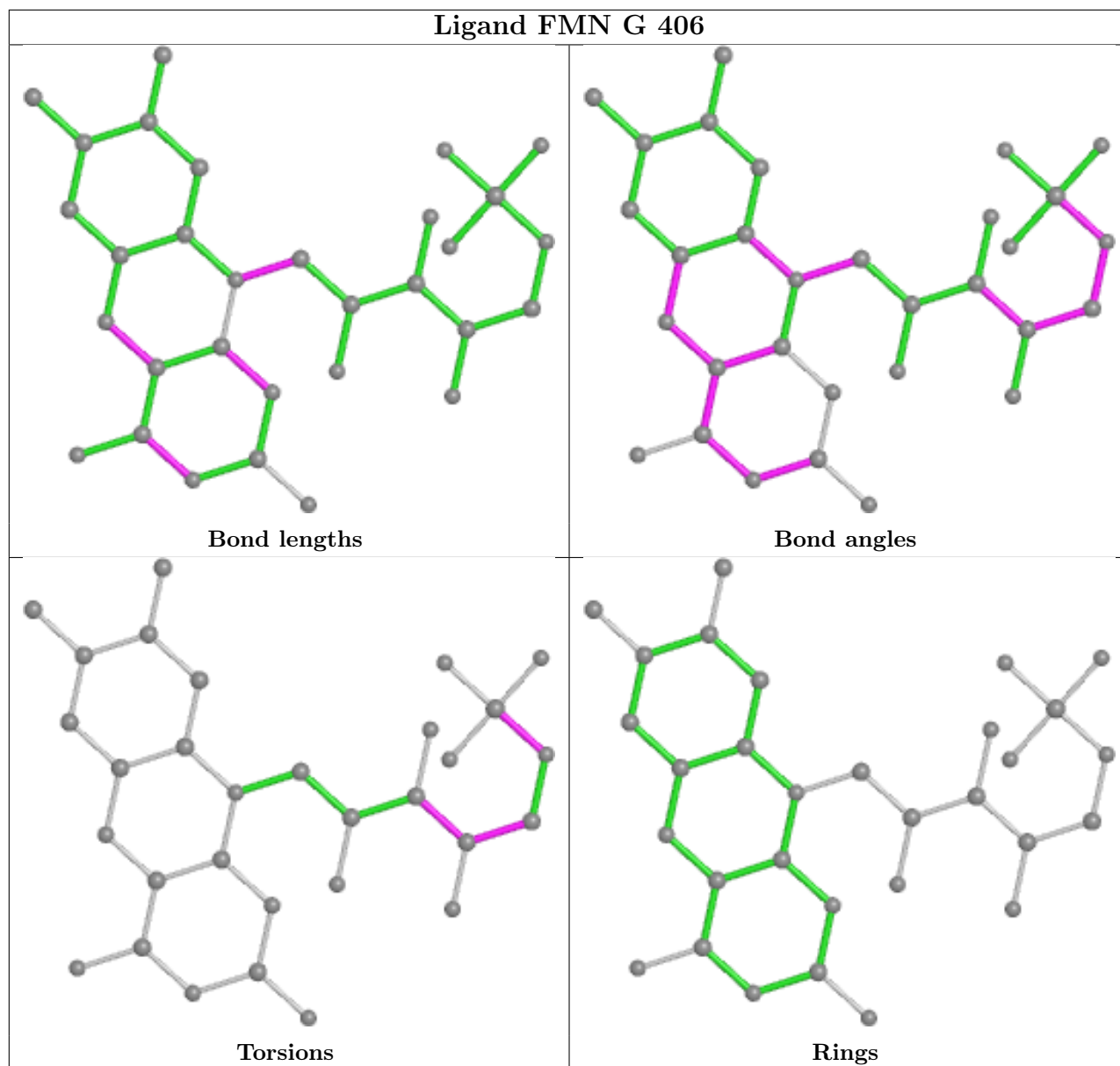
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

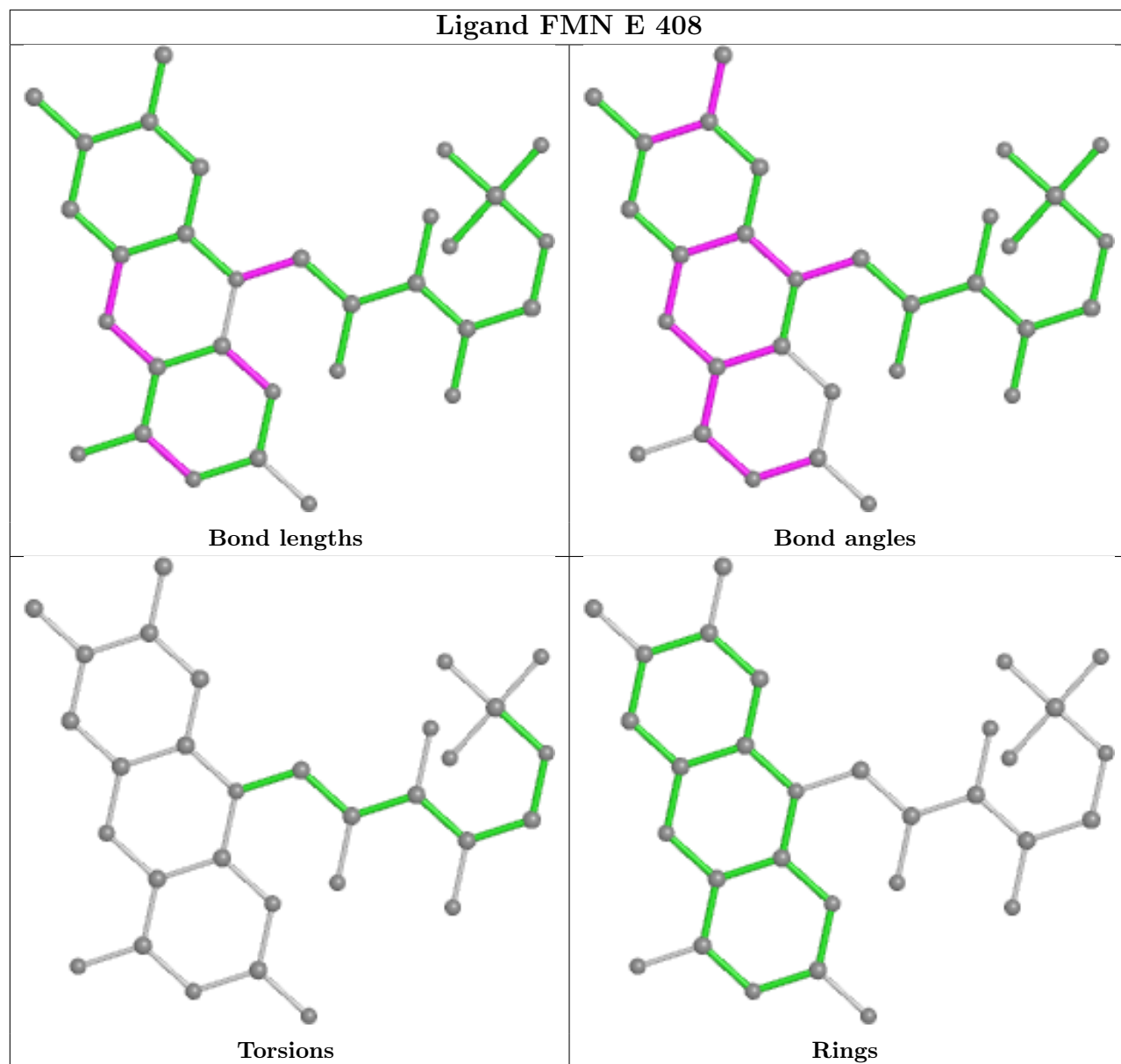
any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

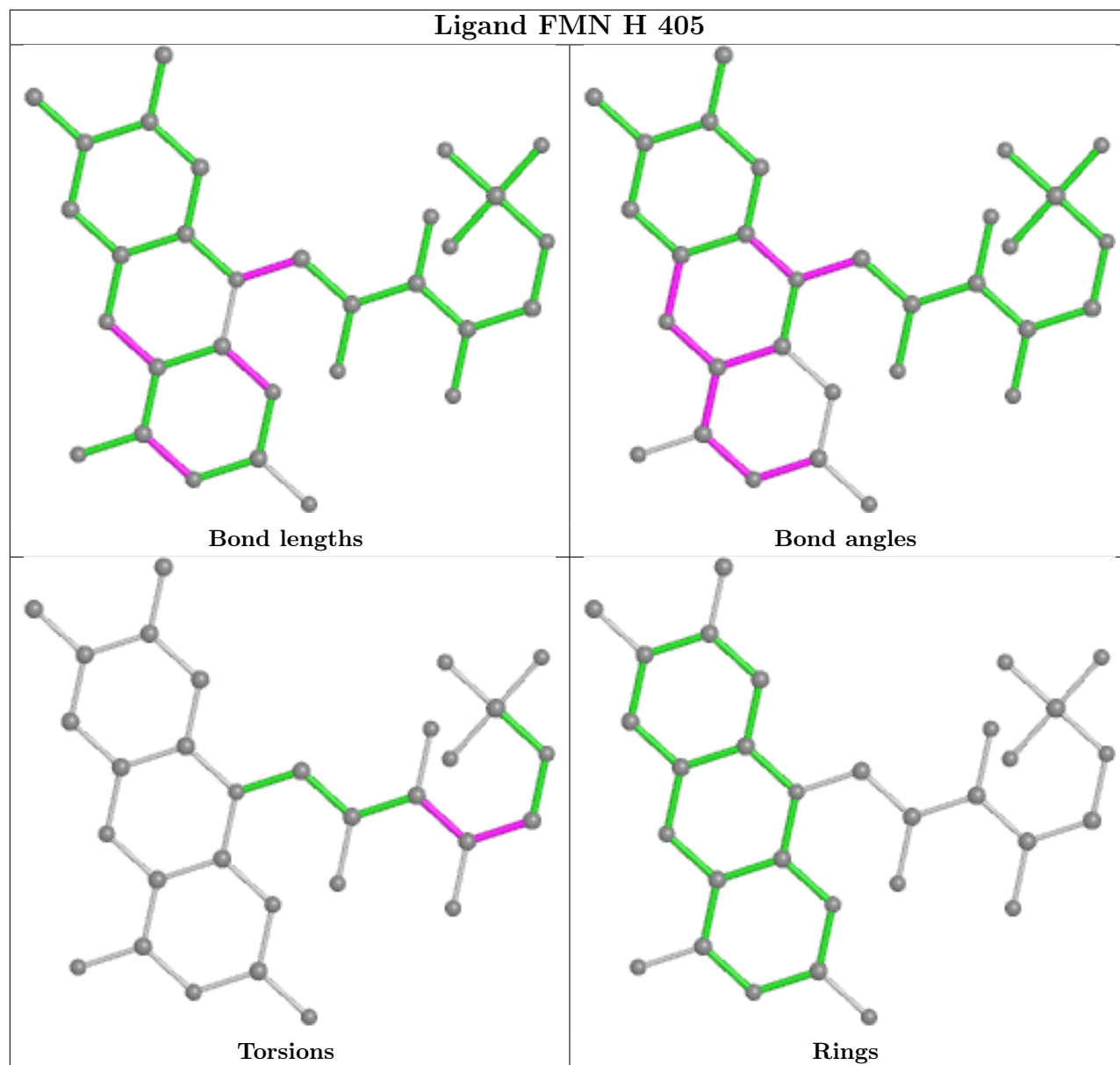




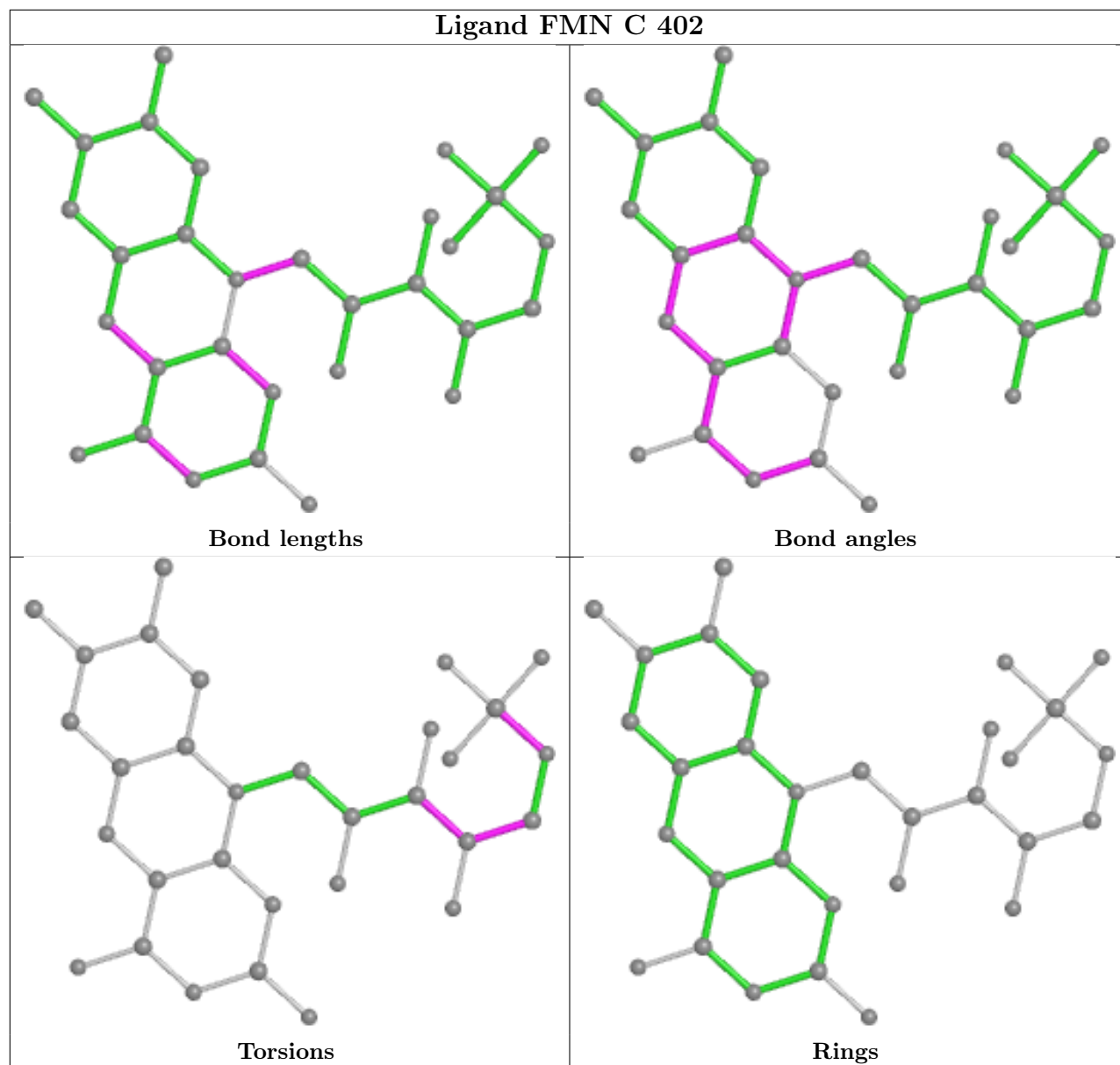


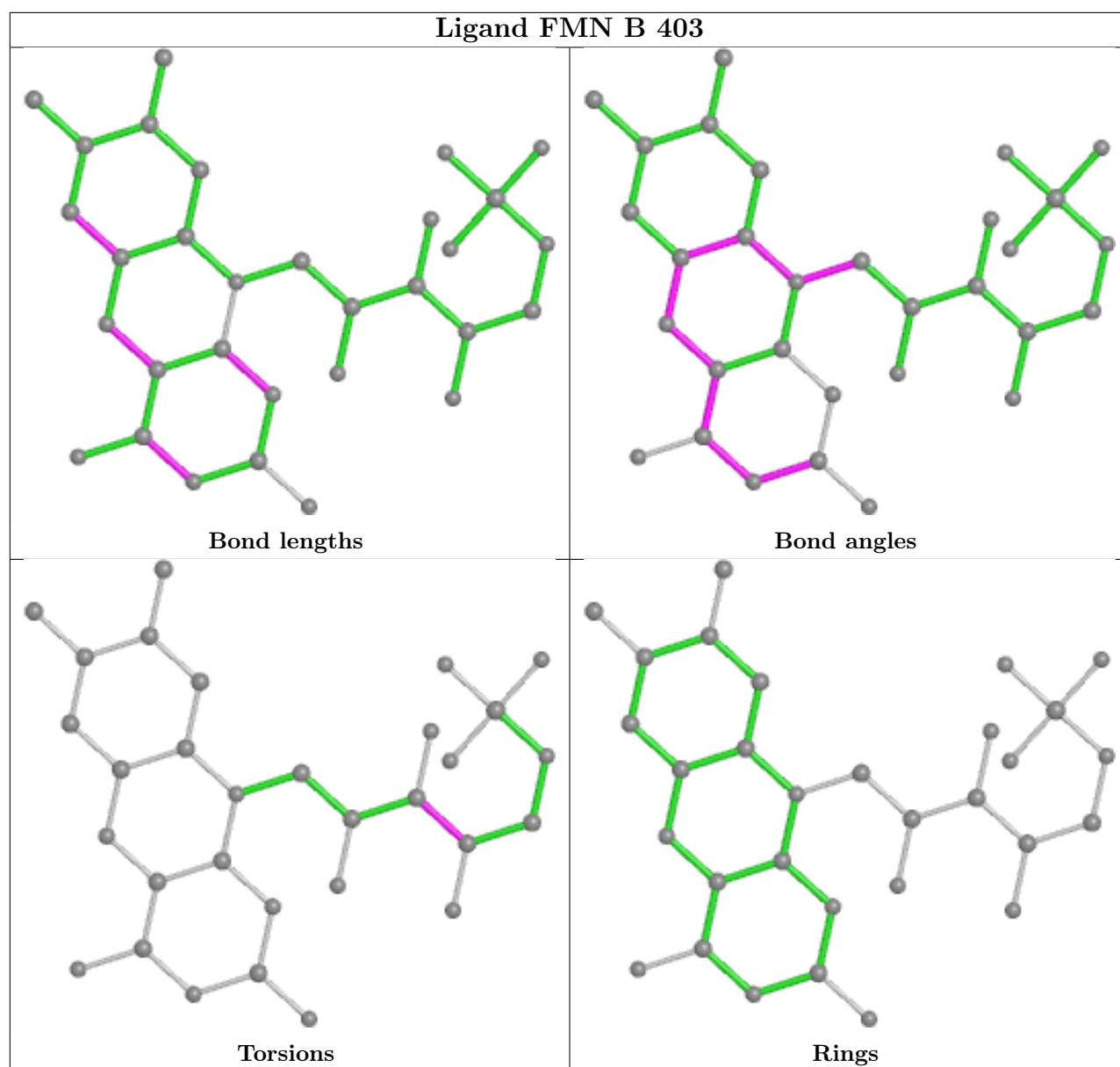












## 5.7 Other polymers [i](#)

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	194/211 (91%)	-0.41	1 (0%) 91 82	25, 26, 32, 36	0
1	B	194/211 (91%)	-0.37	1 (0%) 91 82	25, 28, 39, 51	0
1	C	195/211 (92%)	-0.36	1 (0%) 91 82	25, 28, 37, 44	0
1	D	195/211 (92%)	-0.40	2 (1%) 82 69	25, 27, 34, 42	0
1	E	192/211 (90%)	-0.24	2 (1%) 82 69	29, 37, 46, 49	0
1	F	196/211 (92%)	-0.03	7 (3%) 42 22	35, 44, 58, 61	0
1	G	194/211 (91%)	0.02	7 (3%) 42 22	33, 44, 61, 64	0
1	H	195/211 (92%)	-0.37	3 (1%) 73 56	26, 29, 38, 42	0
All	All	1555/1688 (92%)	-0.27	24 (1%) 73 56	25, 30, 50, 64	0

The worst 5 of 24 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	108	SER	5.4
1	C	203	SER	4.1
1	F	107	SER	3.6
1	G	54	ASP	3.3
1	G	51	ASP	3.3

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

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

### 6.3 Carbohydrates [i](#)

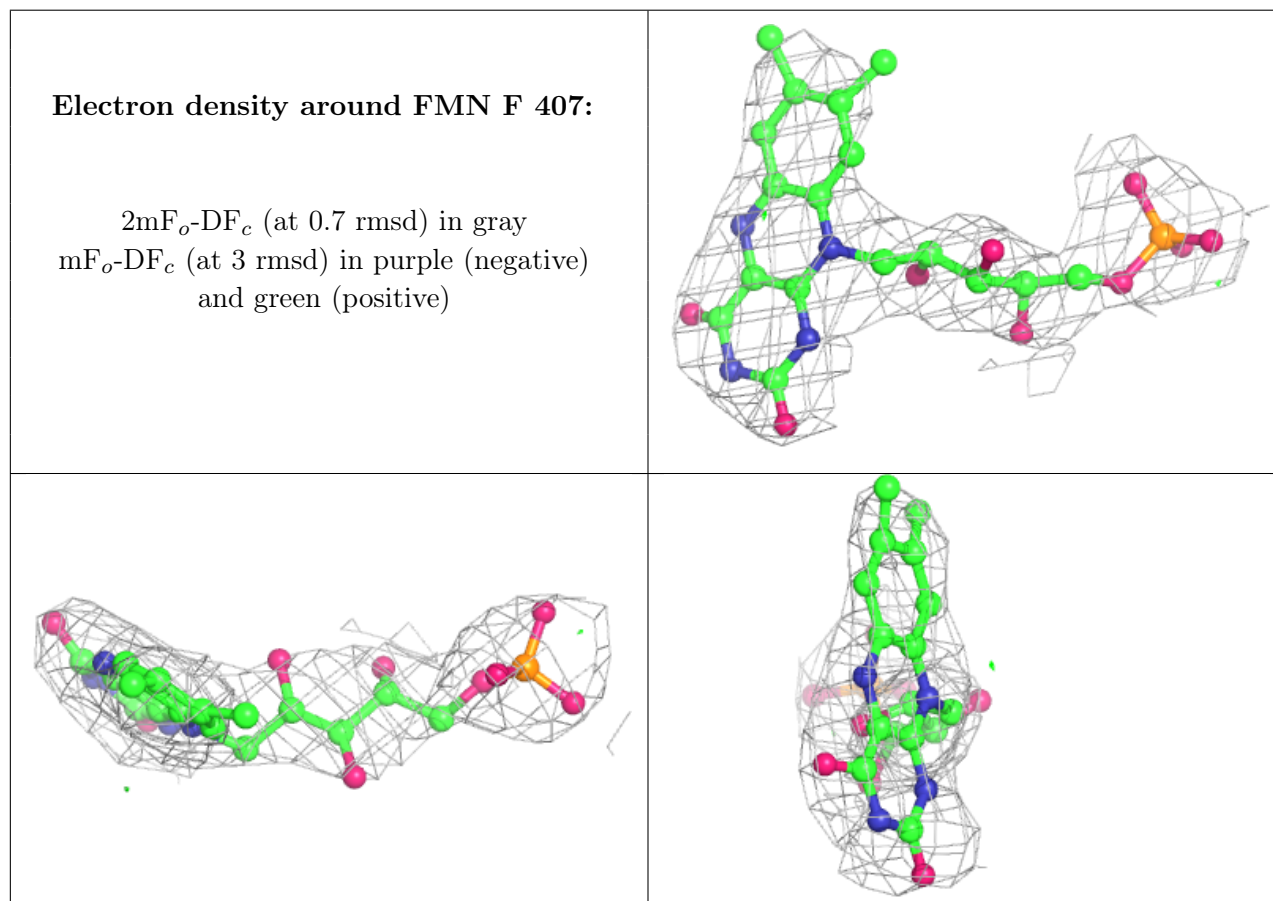
There are no monosaccharides in this entry.

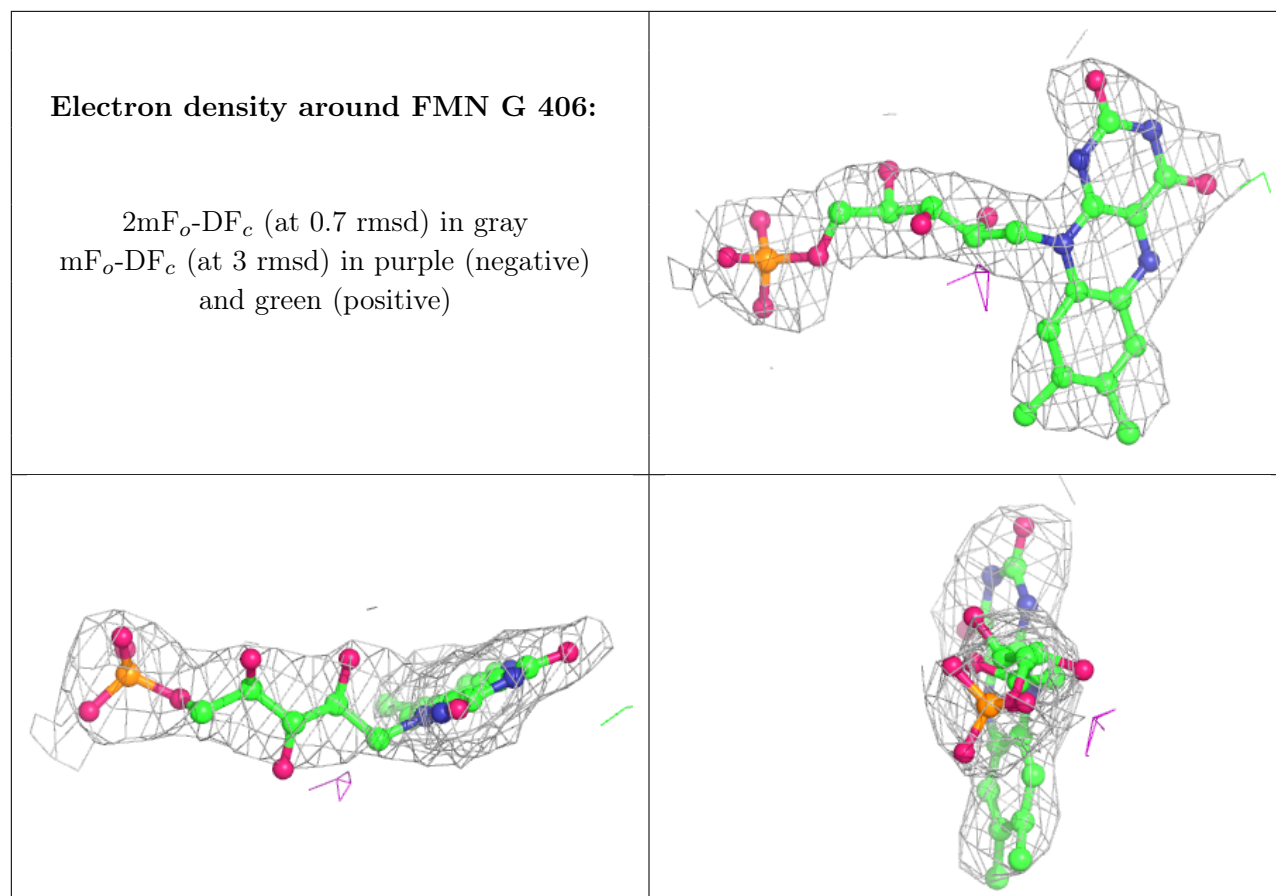
## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	FMN	F	407	31/31	0.95	0.16	35,37,38,39	0
2	FMN	G	406	31/31	0.95	0.15	33,34,45,46	0
2	FMN	E	408	31/31	0.95	0.16	30,31,33,34	0
2	FMN	C	402	31/31	0.95	0.18	25,25,27,28	0
2	FMN	A	404	31/31	0.96	0.16	25,25,26,27	0
2	FMN	H	405	31/31	0.97	0.15	27,29,30,33	0
2	FMN	D	401	31/31	0.97	0.13	25,25,25,25	0
2	FMN	B	403	31/31	0.97	0.15	25,26,27,27	0

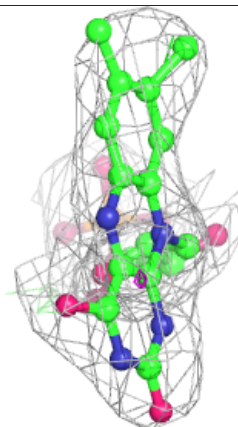
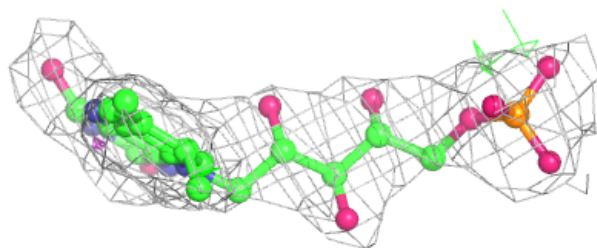
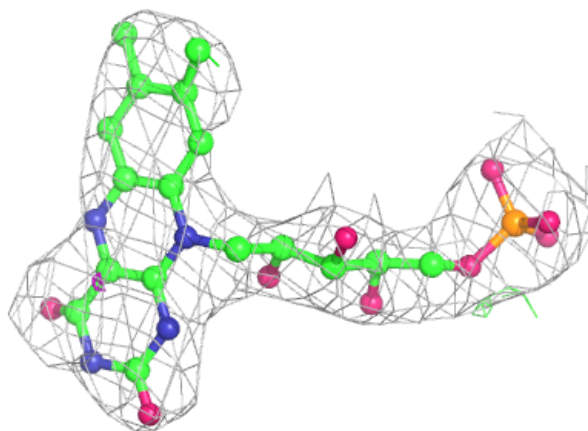
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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





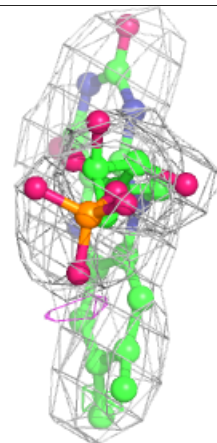
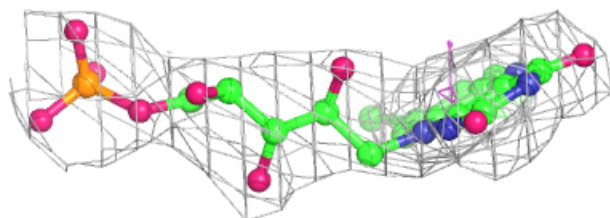
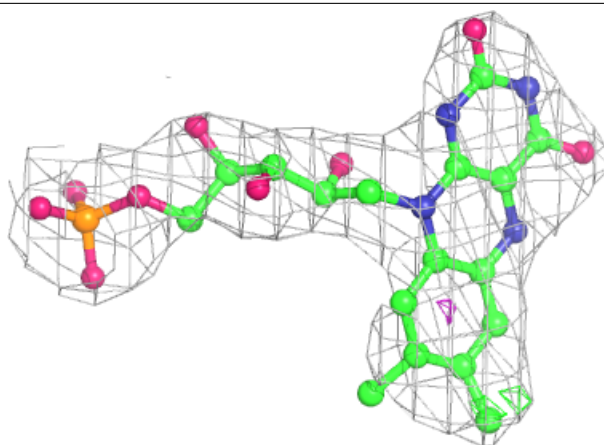
**Electron density around FMN E 408:**

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

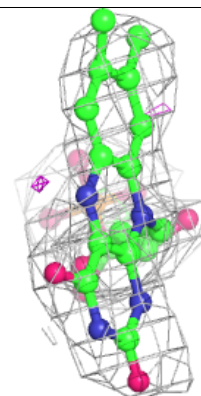
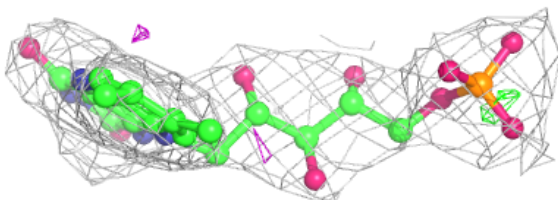
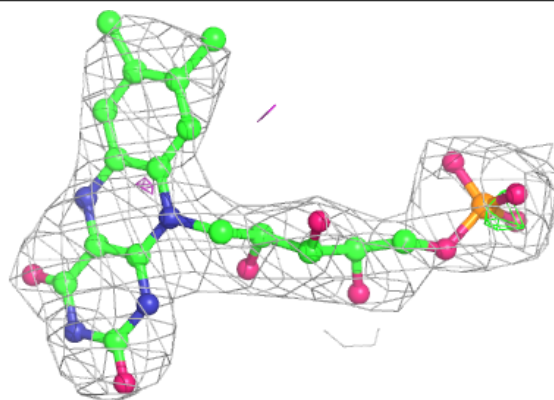


**Electron density around FMN C 402:**

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

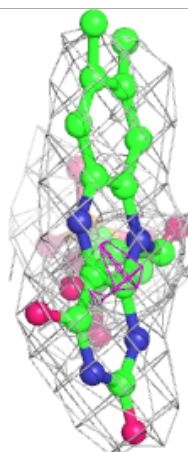
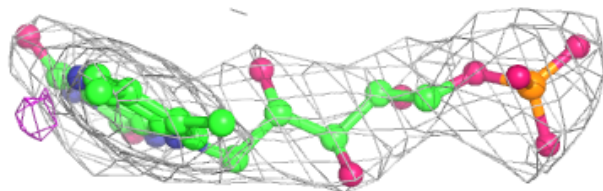
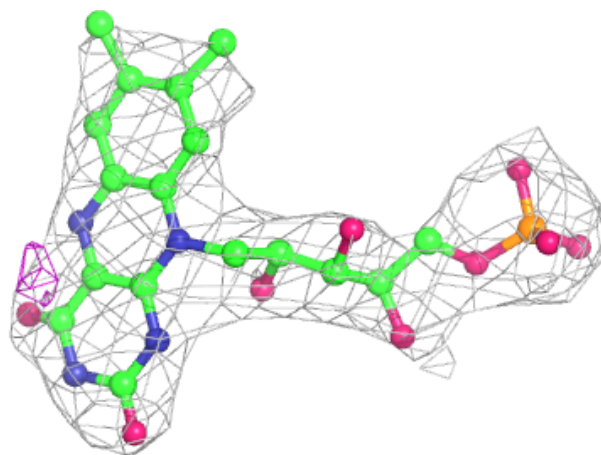
**Electron density around FMN A 404:**

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

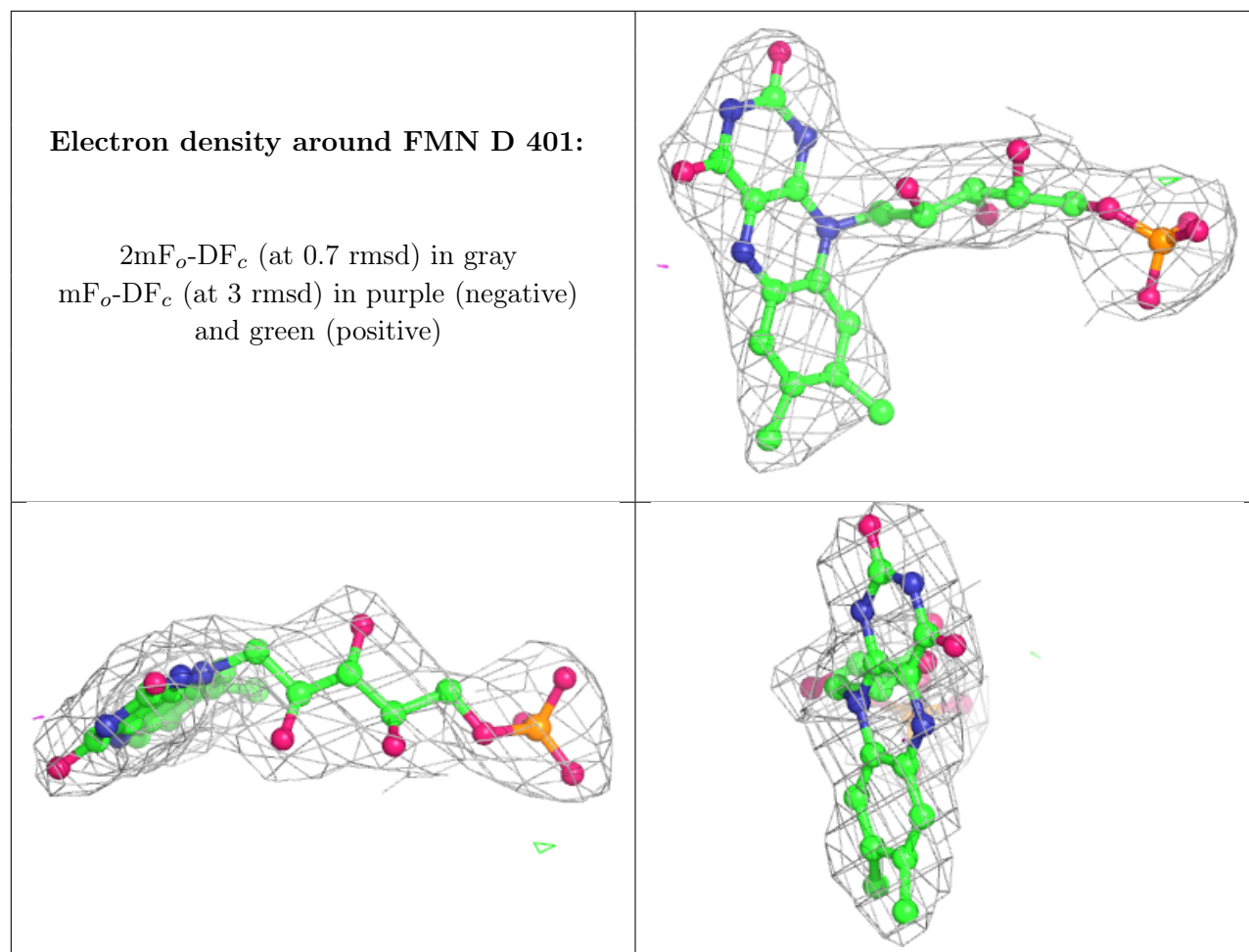


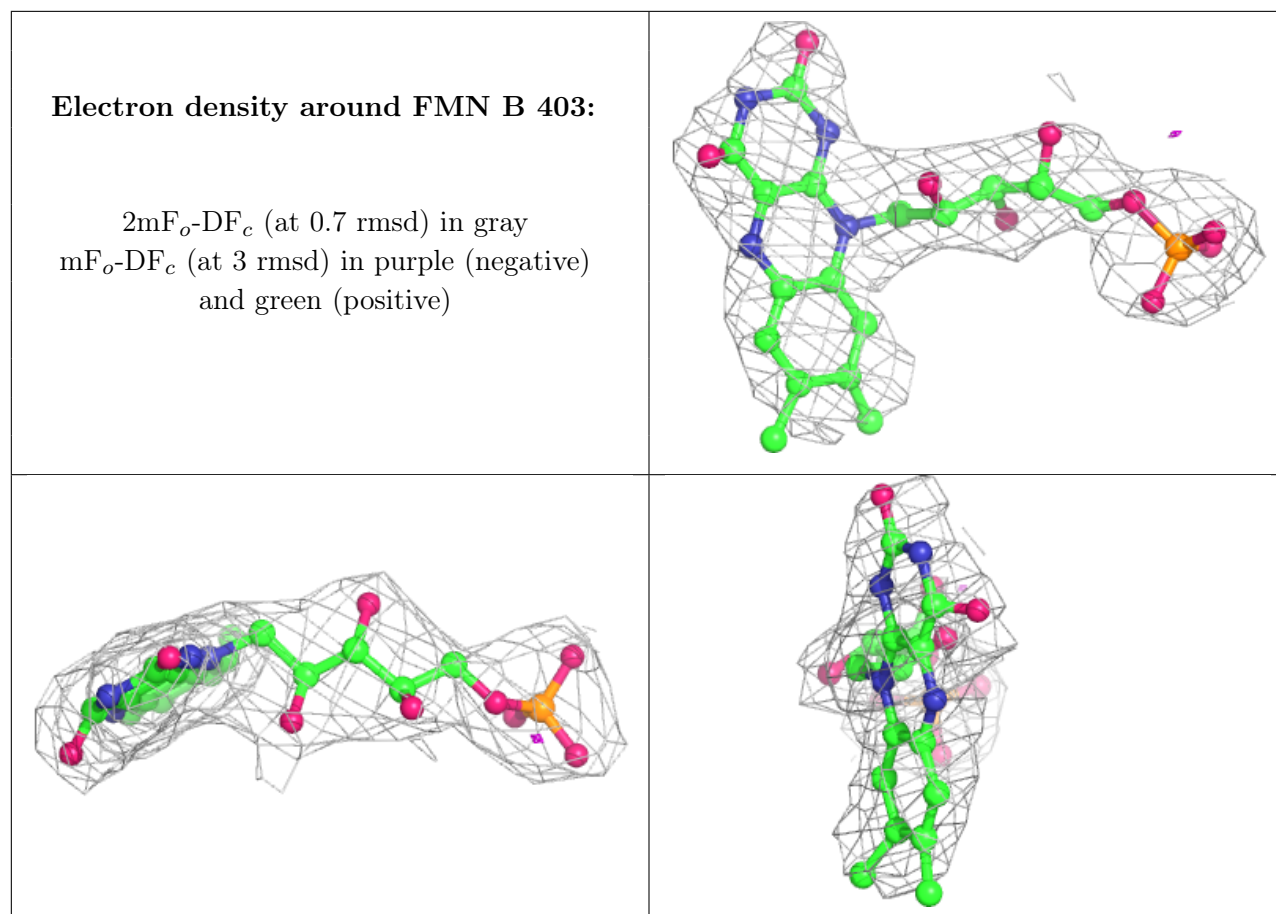
**Electron density around FMN H 405:**

$2mF_o-DF_c$  (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.