



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 23, 2022 – 04:22 PM EST

PDB ID : 7TMB
Title : Crystal Structure of N-ethylmaleimide reductase from *Klebsiella pneumoniae*
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2022-01-19
Resolution : 2.10 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

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 : 2.26
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) : 2.26

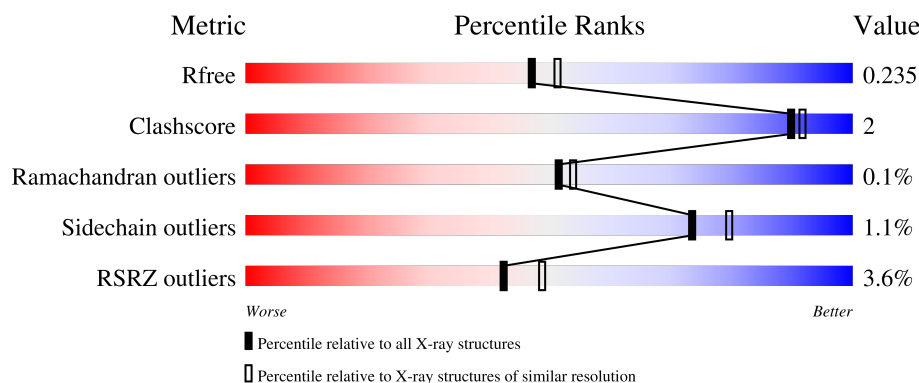
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 ≥ 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	373	<div> <div>2%</div> <div>93%</div> <div>5%</div> <div>2%</div> </div>
1	B	373	<div> <div>4%</div> <div>93%</div> <div>5%</div> <div>2%</div> </div>
1	C	373	<div> <div>3%</div> <div>91%</div> <div>5%</div> <div>2%</div> </div>
1	D	373	<div> <div>4%</div> <div>93%</div> <div>5%</div> <div>2%</div> </div>
1	E	373	<div> <div>2%</div> <div>94%</div> <div>5%</div> <div>2%</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	373	<div><div>5%</div><div>91%</div><div></div></div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 18234 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 N-ethylmaleimide reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	S	0	0	0
			2740	1719	497	520	4			
1	B	359	Total	C	N	O	S	0	1	0
			2721	1707	492	518	4			
1	C	360	Total	C	N	O	S	0	0	0
			2731	1714	495	518	4			
1	D	362	Total	C	N	O	S	0	0	0
			2740	1719	497	520	4			
1	E	362	Total	C	N	O	S	0	0	0
			2743	1720	497	522	4			
1	F	359	Total	C	N	O	S	0	0	0
			2718	1706	493	515	4			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP A0A0H3GQS4
A	2	ALA	-	expression tag	UNP A0A0H3GQS4
A	3	HIS	-	expression tag	UNP A0A0H3GQS4
A	4	HIS	-	expression tag	UNP A0A0H3GQS4
A	5	HIS	-	expression tag	UNP A0A0H3GQS4
A	6	HIS	-	expression tag	UNP A0A0H3GQS4
A	7	HIS	-	expression tag	UNP A0A0H3GQS4
A	8	HIS	-	expression tag	UNP A0A0H3GQS4
B	1	MET	-	expression tag	UNP A0A0H3GQS4
B	2	ALA	-	expression tag	UNP A0A0H3GQS4
B	3	HIS	-	expression tag	UNP A0A0H3GQS4
B	4	HIS	-	expression tag	UNP A0A0H3GQS4
B	5	HIS	-	expression tag	UNP A0A0H3GQS4
B	6	HIS	-	expression tag	UNP A0A0H3GQS4
B	7	HIS	-	expression tag	UNP A0A0H3GQS4
B	8	HIS	-	expression tag	UNP A0A0H3GQS4
C	1	MET	-	expression tag	UNP A0A0H3GQS4

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Chain	Residue	Modelled	Actual	Comment	Reference
C	2	ALA	-	expression tag	UNP A0A0H3GQS4
C	3	HIS	-	expression tag	UNP A0A0H3GQS4
C	4	HIS	-	expression tag	UNP A0A0H3GQS4
C	5	HIS	-	expression tag	UNP A0A0H3GQS4
C	6	HIS	-	expression tag	UNP A0A0H3GQS4
C	7	HIS	-	expression tag	UNP A0A0H3GQS4
C	8	HIS	-	expression tag	UNP A0A0H3GQS4
D	1	MET	-	expression tag	UNP A0A0H3GQS4
D	2	ALA	-	expression tag	UNP A0A0H3GQS4
D	3	HIS	-	expression tag	UNP A0A0H3GQS4
D	4	HIS	-	expression tag	UNP A0A0H3GQS4
D	5	HIS	-	expression tag	UNP A0A0H3GQS4
D	6	HIS	-	expression tag	UNP A0A0H3GQS4
D	7	HIS	-	expression tag	UNP A0A0H3GQS4
D	8	HIS	-	expression tag	UNP A0A0H3GQS4
E	1	MET	-	expression tag	UNP A0A0H3GQS4
E	2	ALA	-	expression tag	UNP A0A0H3GQS4
E	3	HIS	-	expression tag	UNP A0A0H3GQS4
E	4	HIS	-	expression tag	UNP A0A0H3GQS4
E	5	HIS	-	expression tag	UNP A0A0H3GQS4
E	6	HIS	-	expression tag	UNP A0A0H3GQS4
E	7	HIS	-	expression tag	UNP A0A0H3GQS4
E	8	HIS	-	expression tag	UNP A0A0H3GQS4
F	1	MET	-	expression tag	UNP A0A0H3GQS4
F	2	ALA	-	expression tag	UNP A0A0H3GQS4
F	3	HIS	-	expression tag	UNP A0A0H3GQS4
F	4	HIS	-	expression tag	UNP A0A0H3GQS4
F	5	HIS	-	expression tag	UNP A0A0H3GQS4
F	6	HIS	-	expression tag	UNP A0A0H3GQS4
F	7	HIS	-	expression tag	UNP A0A0H3GQS4
F	8	HIS	-	expression tag	UNP A0A0H3GQS4

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

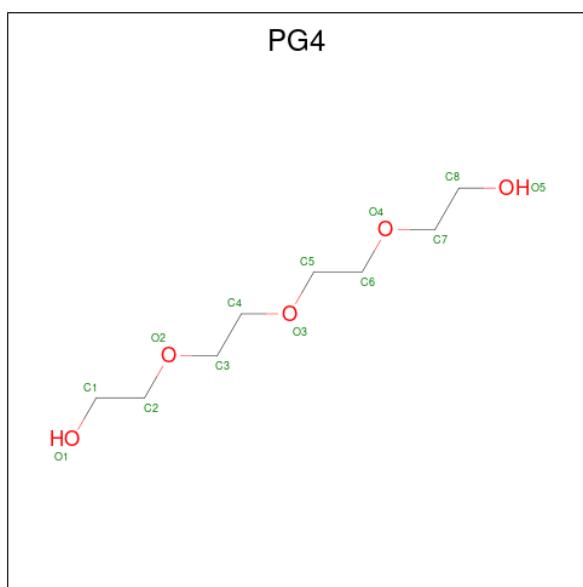
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Na 1 1	0	0
2	B	1	Total Na 1 1	0	0
2	C	1	Total Na 1 1	0	0
2	D	1	Total Na 1 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	E	1	Total 1	Na 1	0	0
2	F	1	Total 1	Na 1	0	0

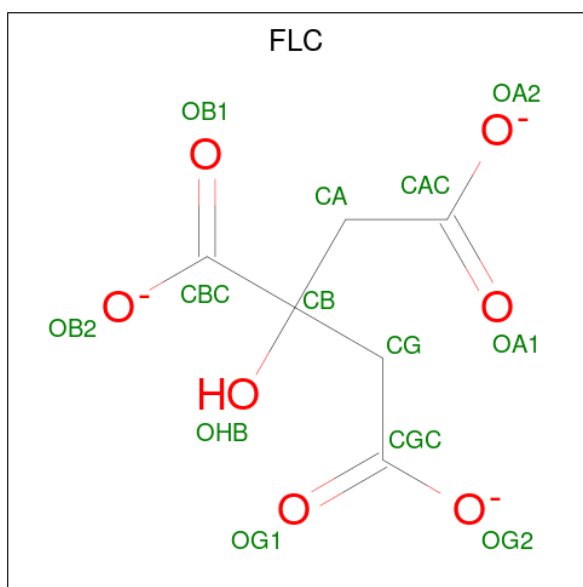
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- The image displays the chemical structure of Flavin Mononucleotide (FMN). It features an isoalloxazine ring system, which is a tricyclic aromatic heterocycle consisting of a benzene ring fused to two pyrimidine rings. The atoms in the ring are labeled: N1, N5, N10 are nitrogen atoms; C2, C4, C6, C7, C8, C9 are carbon atoms. The side chain is attached at the N10 position via a methylene group (C1'). The side chain consists of a ribityl chain: C1' is connected to C2', which is connected to C3'. C3' is a chiral center with a hydroxyl group (OH) shown with a wedge bond and a hydrogen atom (H) shown with a dashed bond. C3' is connected to C4', which is also a chiral center with a hydroxyl group (OH) shown with a wedge bond and a hydrogen atom (H) shown with a dashed bond. C4' is connected to C5', which is connected to an oxygen atom (O5'). This oxygen atom is part of a phosphate group, specifically a 5'-phosphoryl group, represented as O5'-P(=O)(OH)2-O3P, where P is the phosphorus atom, O3P is a phosphate group, and O5' is the bridging oxygen. The structure is drawn with standard chemical notation, including wedge and dash bonds to indicate stereochemistry.

- Molecule 4 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $\text{C}_8\text{H}_{18}\text{O}_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	8	5		
4	A	1	Total	C	O	0	0
			13	8	5		
4	B	1	Total	C	O	0	0
			13	8	5		
4	B	1	Total	C	O	0	0
			13	8	5		
4	C	1	Total	C	O	0	0
			13	8	5		
4	C	1	Total	C	O	0	0
			13	8	5		
4	D	1	Total	C	O	0	0
			13	8	5		
4	D	1	Total	C	O	0	0
			13	8	5		
4	E	1	Total	C	O	0	0
			13	8	5		
4	E	1	Total	C	O	0	0
			13	8	5		
4	F	1	Total	C	O	0	0
			13	8	5		
4	F	1	Total	C	O	0	0
			13	8	5		

- Molecule 5 is CITRATE ANION (three-letter code: FLC) (formula: $C_6H_5O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			13	6	7		
5	C	1	Total	C	O	0	0
			13	6	7		
5	C	1	Total	C	O	0	0
			13	6	7		
5	D	1	Total	C	O	0	0
			13	6	7		
5	E	1	Total	C	O	0	0
			13	6	7		
5	F	1	Total	C	O	0	0
			13	6	7		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	228	Total	O	0	0
			228	228		
6	B	236	Total	O	0	0
			236	236		
6	C	200	Total	O	0	0
			200	200		
6	D	264	Total	O	0	0
			264	264		
6	E	264	Total	O	0	0
			264	264		
6	F	223	Total	O	0	0
			223	223		

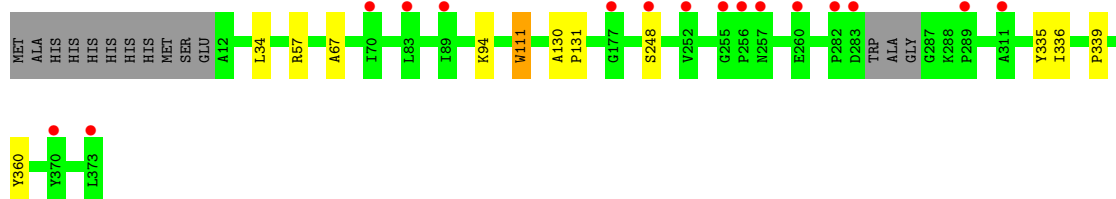
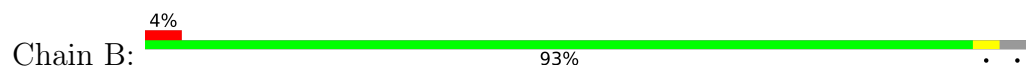
3 Residue-property plots [i](#)

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.

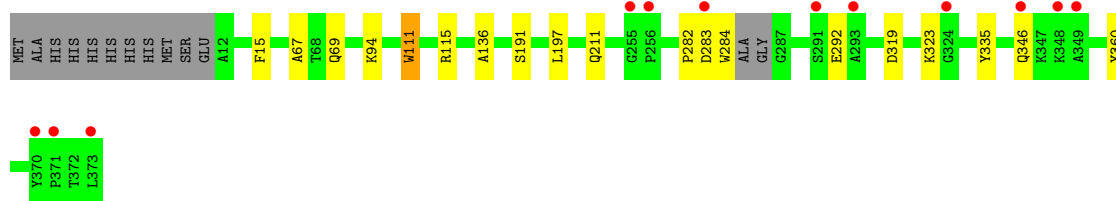
- Molecule 1: N-ethylmaleimide reductase



- Molecule 1: N-ethylmaleimide reductase



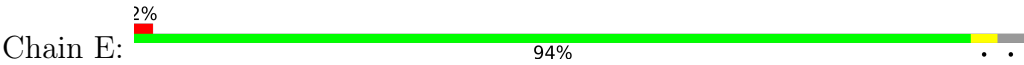
- Molecule 1: N-ethylmaleimide reductase



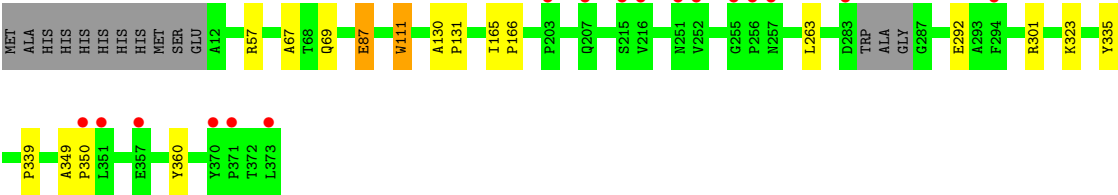
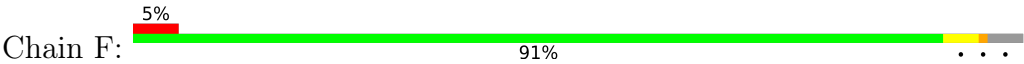
- Molecule 1: N-ethylmaleimide reductase



- Molecule 1: N-ethylmaleimide reductase



● Molecule 1: N-ethylmaleimide reductase



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	131.31Å 223.87Å 223.37Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.54 – 2.10 49.24 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.8 (25.54-2.10) 100.0 (49.24-2.10)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 2.10Å)	Xtriage
Refinement program	PHENIX 1.20_4474	Depositor
R, R_{free}	0.192 , 0.231 0.200 , 0.235	Depositor DCC
R_{free} test set	9444 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.502	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.000 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.000 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	18234	wwPDB-VP
Average B, all atoms (Å ²)	30.0	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 54.97 % 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 3.3810e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

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: FLC, NA, PG4, 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	0.41	0/2802	0.63	0/3810
1	B	0.37	0/2785	0.61	0/3788
1	C	0.37	0/2792	0.60	0/3795
1	D	0.43	0/2802	0.63	0/3810
1	E	0.41	0/2805	0.62	0/3814
1	F	0.39	0/2779	0.62	0/3779
All	All	0.40	0/16765	0.62	0/22796

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	2740	0	2685	11	0
1	B	2721	0	2657	6	0
1	C	2731	0	2676	11	0
1	D	2740	0	2685	10	0
1	E	2743	0	2687	9	0
1	F	2718	0	2659	10	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	31	0	19	2	0
3	B	31	0	19	1	0
3	C	31	0	19	2	0
3	D	31	0	19	2	0
3	E	31	0	19	2	0
3	F	31	0	19	2	0
4	A	26	0	35	1	0
4	B	26	0	35	1	0
4	C	26	0	35	1	0
4	D	26	0	36	2	0
4	E	26	0	35	0	0
4	F	26	0	34	0	0
5	B	13	0	5	0	0
5	C	26	0	10	2	0
5	D	13	0	5	2	0
5	E	13	0	5	3	0
5	F	13	0	5	0	0
6	A	228	0	0	0	0
6	B	236	0	0	1	0
6	C	200	0	0	1	0
6	D	264	0	0	1	0
6	E	264	0	0	1	0
6	F	223	0	0	1	0
All	All	18234	0	16403	58	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 2.

The worst 5 of 58 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:360:TYR:CZ	3:A:402:FMN:HM72	2.34	0.62
1:E:292:GLU:OE1	1:E:323:LYS:NZ	2.34	0.60
1:B:67:ALA:HB1	1:B:111:TRP:CD2	2.37	0.59
1:F:360:TYR:CZ	3:F:402:FMN:HM72	2.37	0.58
1:A:235:SER:HB2	5:C:404:FLC:CA	2.33	0.58

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	360/373 (96%)	348 (97%)	11 (3%)	1 (0%)	41	41
1	B	356/373 (95%)	342 (96%)	14 (4%)	0	100	100
1	C	356/373 (95%)	342 (96%)	13 (4%)	1 (0%)	41	41
1	D	360/373 (96%)	347 (96%)	12 (3%)	1 (0%)	41	41
1	E	360/373 (96%)	348 (97%)	12 (3%)	0	100	100
1	F	355/373 (95%)	341 (96%)	14 (4%)	0	100	100
All	All	2147/2238 (96%)	2068 (96%)	76 (4%)	3 (0%)	51	54

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	283	ASP
1	D	284	TRP
1	A	286	GLY

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	279/293 (95%)	277 (99%)	2 (1%)	84	88
1	B	278/293 (95%)	275 (99%)	3 (1%)	73	79

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	279/293 (95%)	276 (99%)	3 (1%)	73	79
1	D	279/293 (95%)	276 (99%)	3 (1%)	73	79
1	E	280/293 (96%)	277 (99%)	3 (1%)	73	79
1	F	277/293 (94%)	273 (99%)	4 (1%)	67	73
All	All	1672/1758 (95%)	1654 (99%)	18 (1%)	73	79

5 of 18 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	F	69	GLN
1	F	335	TYR
1	F	111	TRP
1	D	111	TRP
1	E	335	TYR

Sometimes 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 30 ligands modelled in this entry, 6 are monoatomic - leaving 24 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
5	FLC	C	403	-	3,12,12	0.56	0	3,17,17	0.86	0
3	FMN	E	402	-	31,33,33	1.49	5 (16%)	40,50,50	2.01	8 (20%)
4	PG4	A	403	-	12,12,12	0.22	0	11,11,11	0.35	0
4	PG4	C	406	2	12,12,12	0.14	0	11,11,11	0.61	0
4	PG4	F	404	-	12,12,12	0.25	0	11,11,11	0.41	0
3	FMN	B	402	-	31,33,33	1.43	4 (12%)	40,50,50	2.10	7 (17%)
3	FMN	F	402	-	31,33,33	1.53	4 (12%)	40,50,50	1.99	6 (15%)
5	FLC	E	403	-	3,12,12	0.38	0	3,17,17	2.75	2 (66%)
4	PG4	F	405	2	12,12,12	0.12	0	11,11,11	0.52	0
3	FMN	A	402	-	31,33,33	1.49	6 (19%)	40,50,50	2.19	9 (22%)
5	FLC	F	403	-	3,12,12	0.38	0	3,17,17	0.99	0
5	FLC	D	403	-	3,12,12	0.58	0	3,17,17	0.24	0
4	PG4	E	404	-	12,12,12	0.15	0	11,11,11	0.51	0
3	FMN	C	402	-	31,33,33	1.50	5 (16%)	40,50,50	2.01	9 (22%)
5	FLC	B	403	-	3,12,12	0.16	0	3,17,17	1.50	1 (33%)
4	PG4	C	405	-	12,12,12	0.21	0	11,11,11	0.42	0
4	PG4	A	404	2	12,12,12	0.11	0	11,11,11	0.46	0
4	PG4	E	405	2	12,12,12	0.14	0	11,11,11	0.46	0
4	PG4	B	404	-	12,12,12	0.18	0	11,11,11	0.59	0
4	PG4	B	405	2	12,12,12	0.08	0	11,11,11	0.70	0
5	FLC	C	404	-	3,12,12	0.37	0	3,17,17	2.64	1 (33%)
4	PG4	D	404	-	12,12,12	0.18	0	11,11,11	0.52	0
4	PG4	D	405	2	12,12,12	0.24	0	11,11,11	0.34	0
3	FMN	D	402	-	31,33,33	1.44	4 (12%)	40,50,50	2.20	9 (22%)

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
5	FLC	C	403	-	-	3/6/16/16	-
3	FMN	E	402	-	-	2/18/18/18	0/3/3/3
4	PG4	A	403	-	-	5/10/10/10	-
4	PG4	C	406	2	-	4/10/10/10	-
4	PG4	F	404	-	-	5/10/10/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	FMN	B	402	-	-	1/18/18/18	0/3/3/3
3	FMN	F	402	-	-	1/18/18/18	0/3/3/3
5	FLC	E	403	-	-	1/6/16/16	-
4	PG4	F	405	2	-	4/10/10/10	-
3	FMN	A	402	-	-	2/18/18/18	0/3/3/3
5	FLC	F	403	-	-	0/6/16/16	-
5	FLC	D	403	-	-	0/6/16/16	-
4	PG4	E	404	-	-	5/10/10/10	-
3	FMN	C	402	-	-	1/18/18/18	0/3/3/3
5	FLC	B	403	-	-	0/6/16/16	-
4	PG4	C	405	-	-	4/10/10/10	-
4	PG4	A	404	2	-	4/10/10/10	-
4	PG4	E	405	2	-	5/10/10/10	-
4	PG4	B	404	-	-	4/10/10/10	-
4	PG4	B	405	2	-	4/10/10/10	-
5	FLC	C	404	-	-	2/6/16/16	-
4	PG4	D	404	-	-	4/10/10/10	-
4	PG4	D	405	2	-	3/10/10/10	-
3	FMN	D	402	-	-	1/18/18/18	0/3/3/3

The worst 5 of 28 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	402	FMN	C10-N1	4.39	1.38	1.33
3	F	402	FMN	C10-N1	4.24	1.38	1.33
3	A	402	FMN	C10-N1	4.05	1.38	1.33
3	E	402	FMN	C4-N3	3.91	1.39	1.33
3	B	402	FMN	C10-N1	3.86	1.38	1.33

The worst 5 of 52 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	FMN	C4A-N5-C5A	7.69	124.46	116.77
3	D	402	FMN	C4A-N5-C5A	7.69	124.45	116.77
3	B	402	FMN	C4A-N5-C5A	7.04	123.81	116.77
3	D	402	FMN	C10-C4A-N5	-6.79	116.56	121.26
3	C	402	FMN	C4A-N5-C5A	6.70	123.46	116.77

There are no chirality outliers.

5 of 65 torsion outliers are listed below:

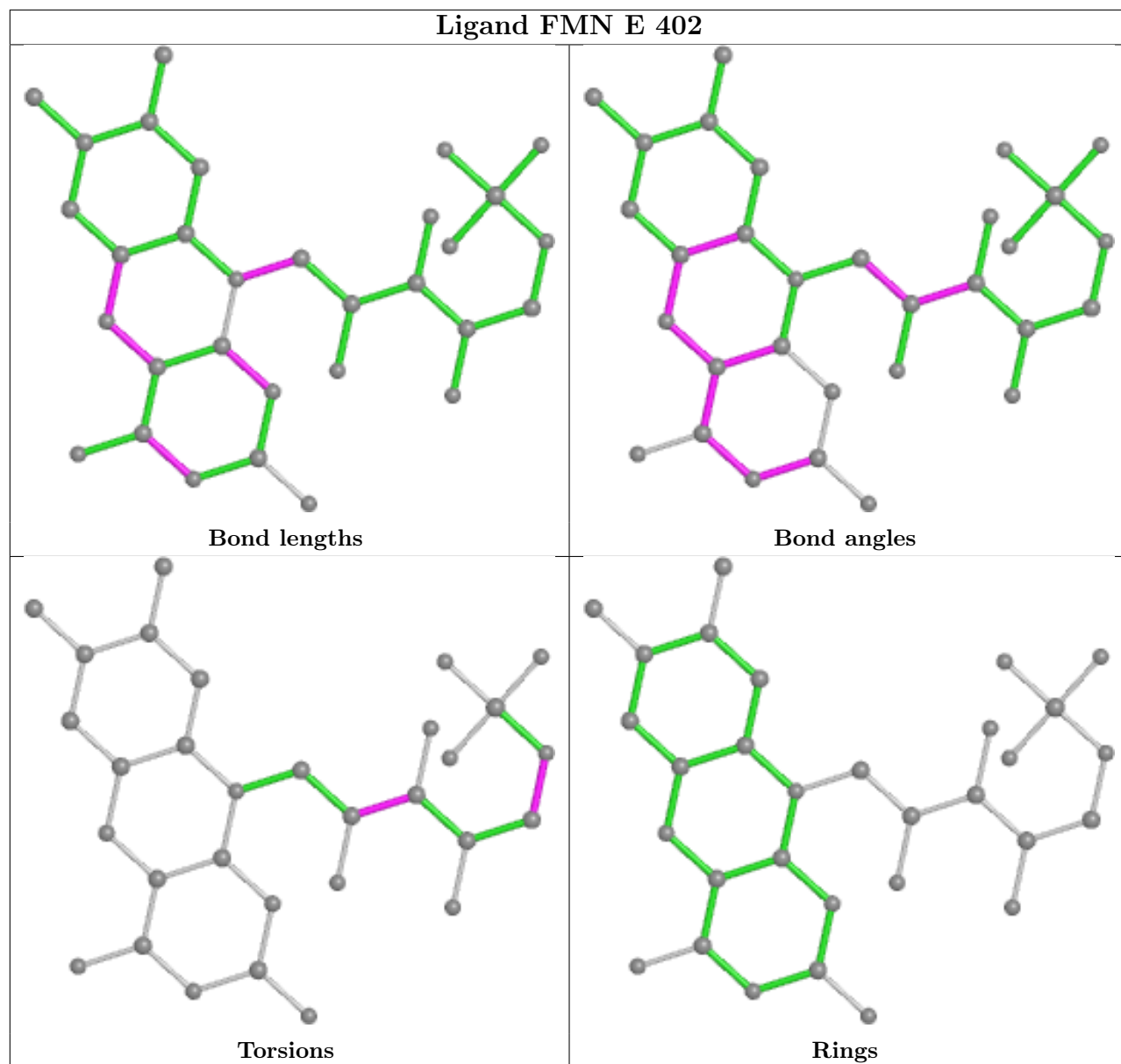
Mol	Chain	Res	Type	Atoms
5	C	403	FLC	CAC-CA-CB-CBC
5	C	403	FLC	CAC-CA-CB-CG
4	E	404	PG4	O4-C7-C8-O5
4	E	404	PG4	O3-C5-C6-O4
4	D	405	PG4	O3-C5-C6-O4

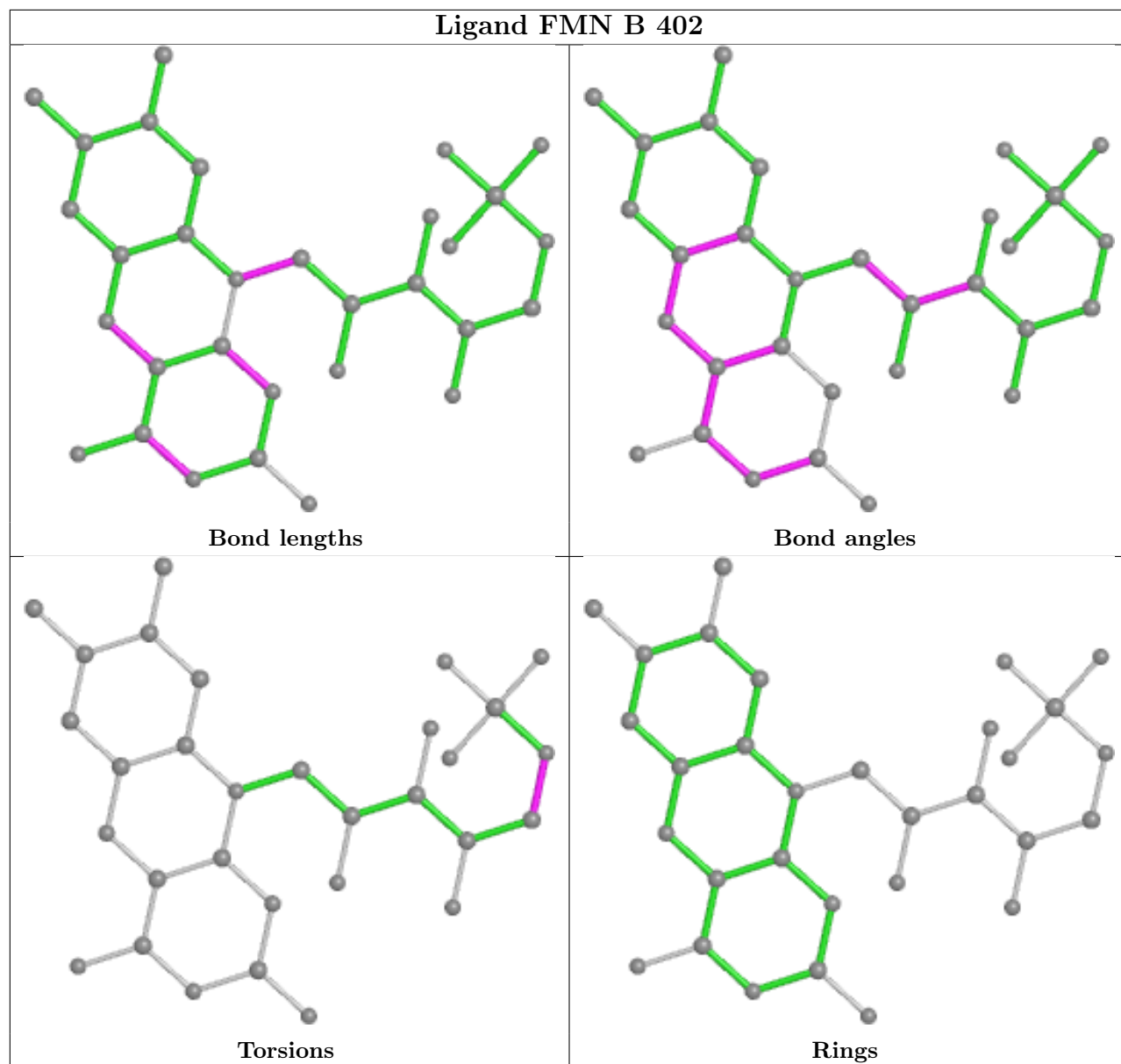
There are no ring outliers.

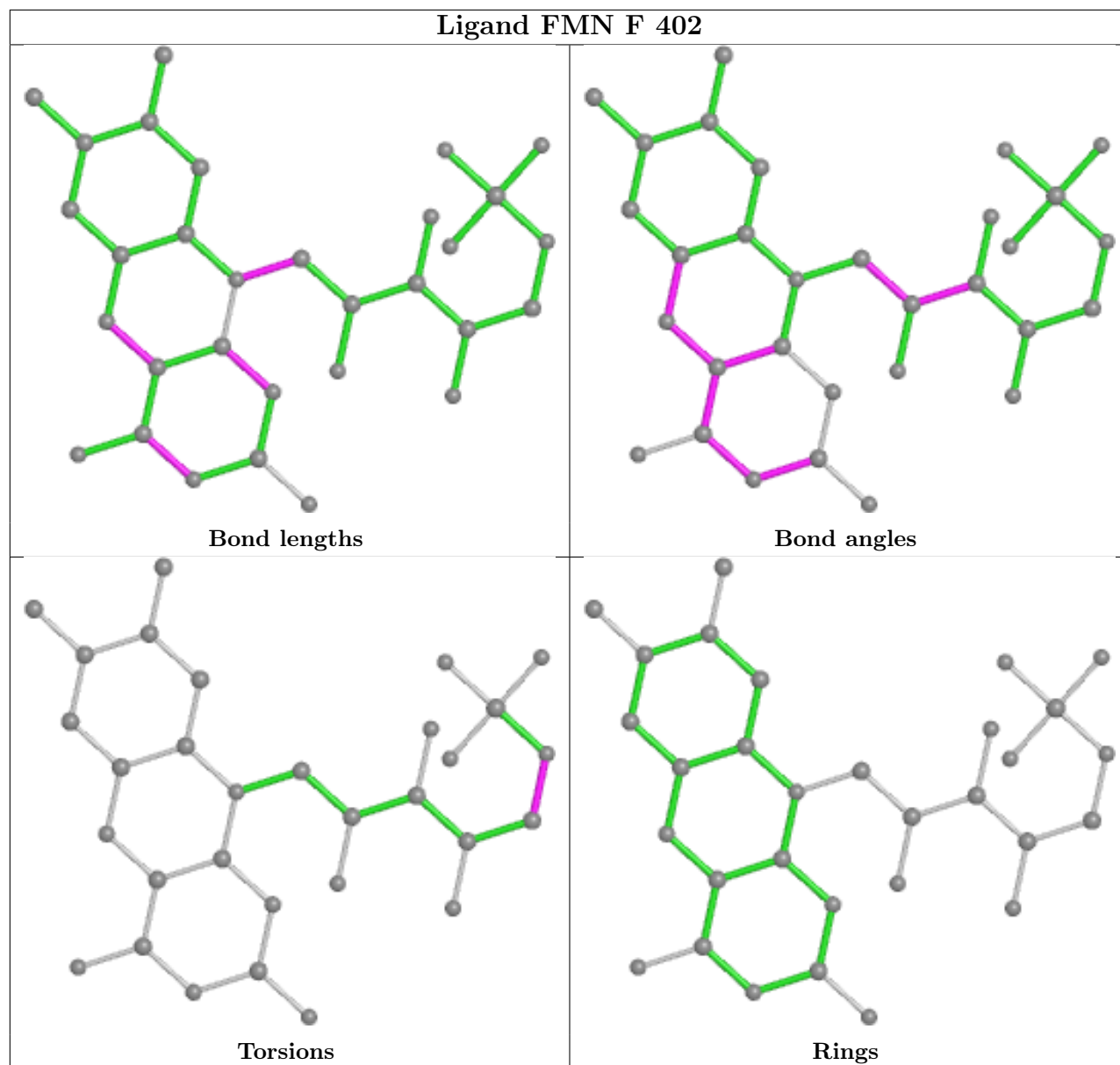
13 monomers are involved in 23 short contacts:

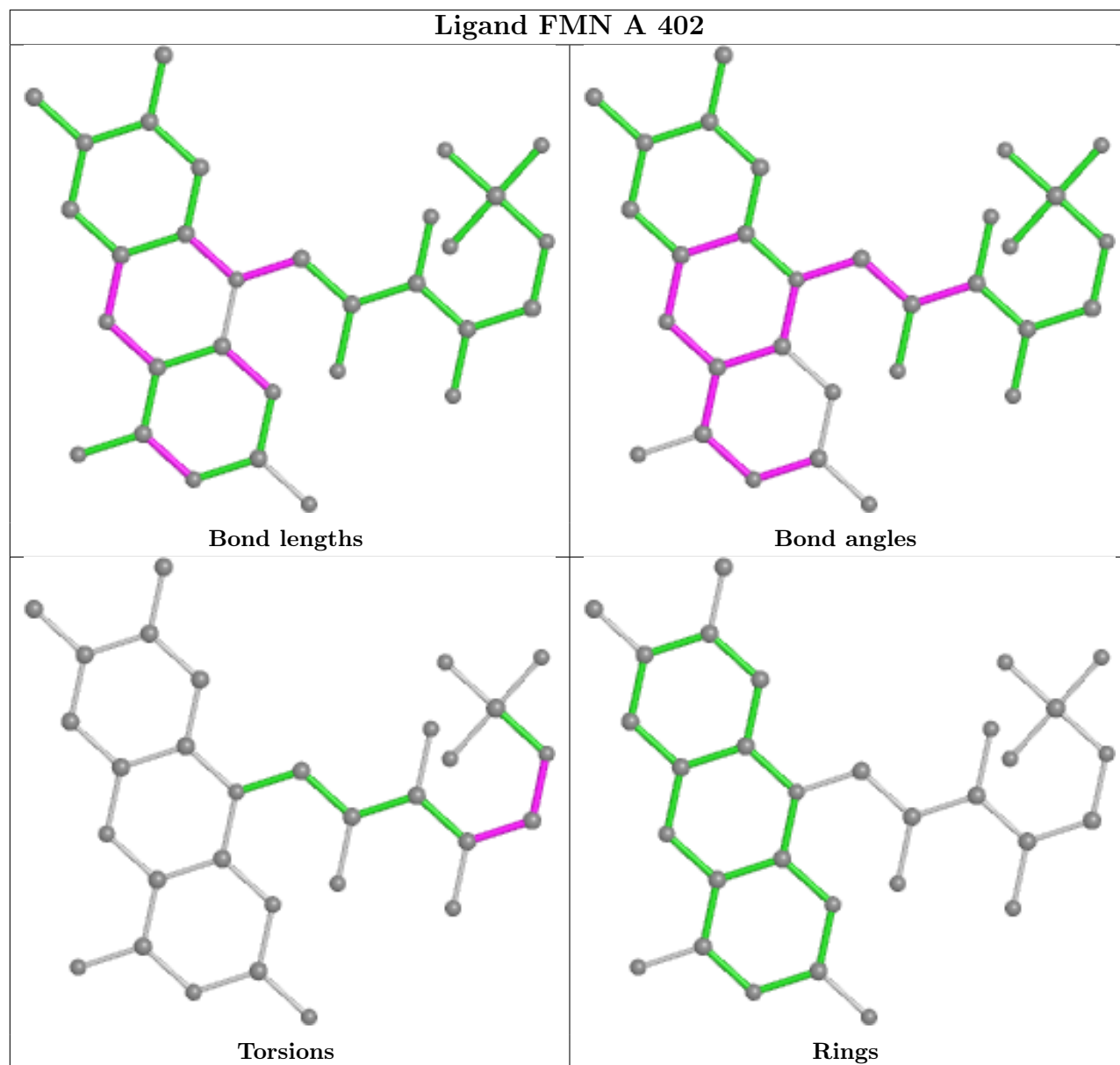
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	402	FMN	2	0
4	A	403	PG4	1	0
3	B	402	FMN	1	0
3	F	402	FMN	2	0
5	E	403	FLC	3	0
3	A	402	FMN	2	0
5	D	403	FLC	2	0
3	C	402	FMN	2	0
4	C	405	PG4	1	0
4	B	404	PG4	1	0
5	C	404	FLC	2	0
4	D	404	PG4	2	0
3	D	402	FMN	2	0

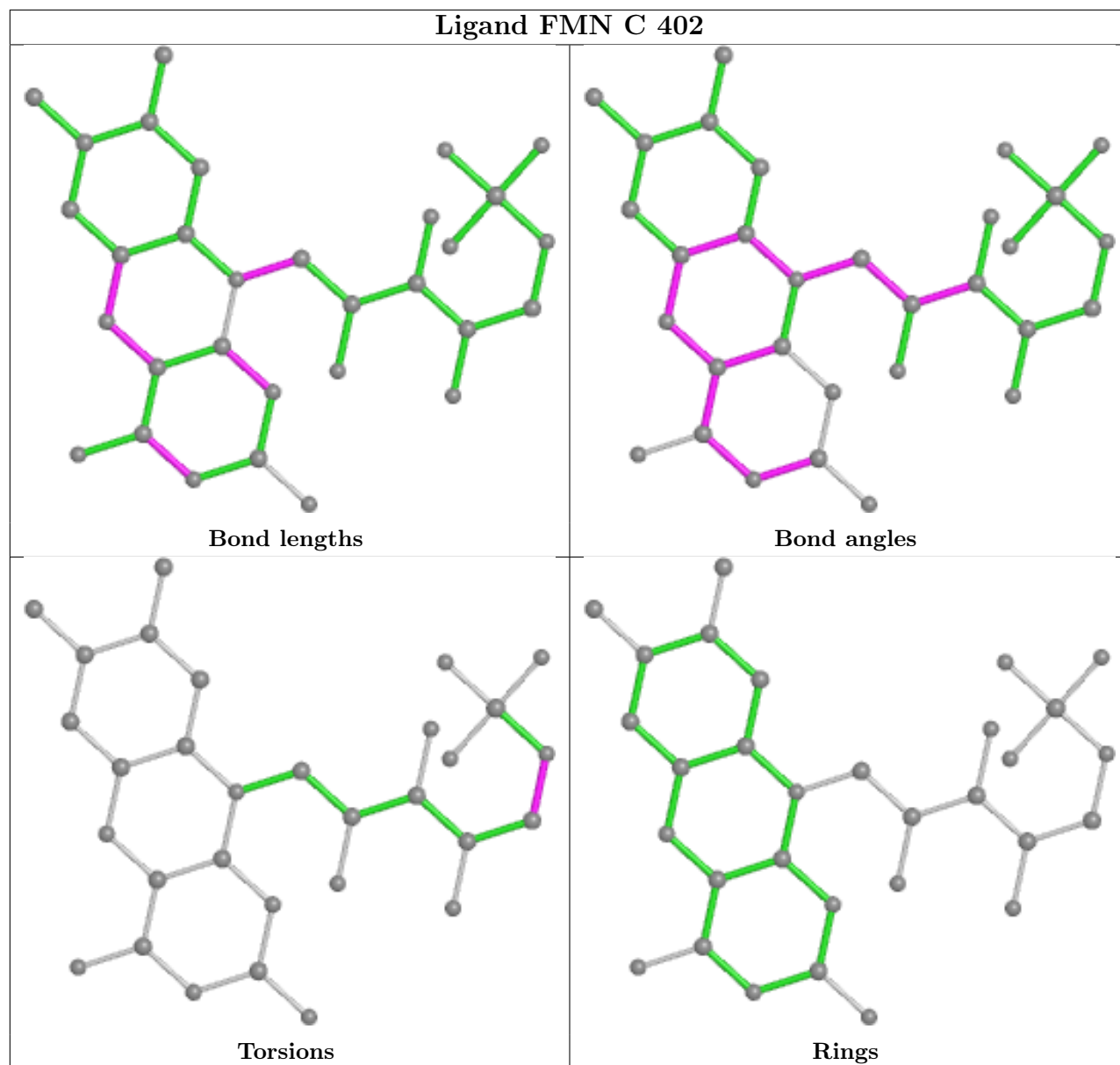
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

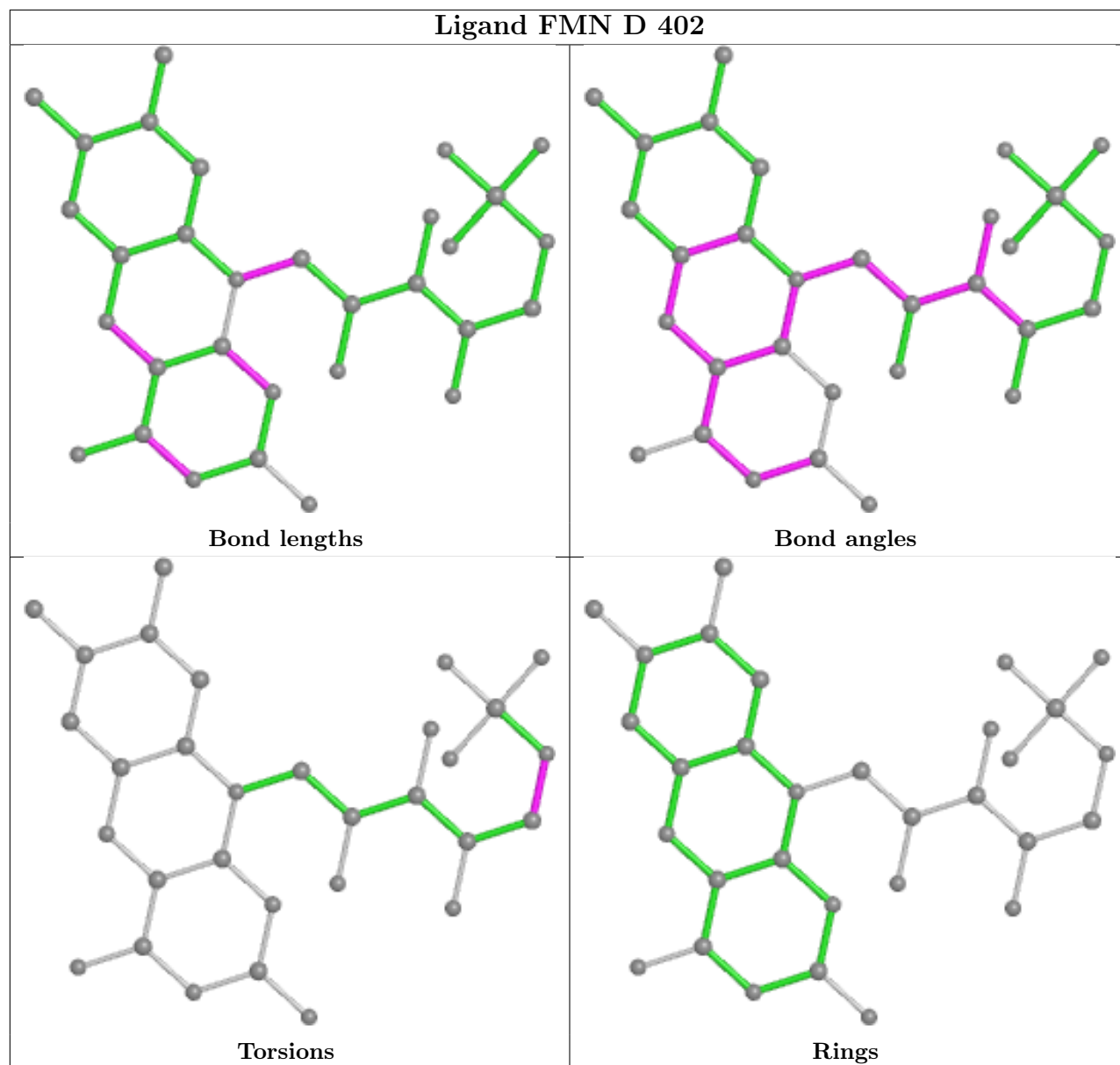












5.7 Other polymers [i](#)

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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	362/373 (97%)	0.07	9 (2%) 57 62	17, 26, 48, 74	0
1	B	359/373 (96%)	0.26	16 (4%) 33 38	17, 29, 49, 66	0
1	C	360/373 (96%)	0.16	12 (3%) 46 53	19, 31, 55, 70	0
1	D	362/373 (97%)	0.10	14 (3%) 39 45	16, 23, 42, 66	0
1	E	362/373 (97%)	0.15	9 (2%) 57 62	16, 26, 46, 74	0
1	F	359/373 (96%)	0.27	17 (4%) 31 37	18, 30, 50, 70	0
All	All	2164/2238 (96%)	0.17	77 (3%) 42 49	16, 28, 49, 74	0

The worst 5 of 77 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	283	ASP	5.5
1	F	373	LEU	4.5
1	E	284	TRP	4.0
1	F	256	PRO	3.8
1	A	351	LEU	3.7

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 ⓘ

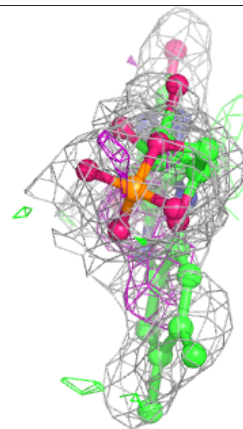
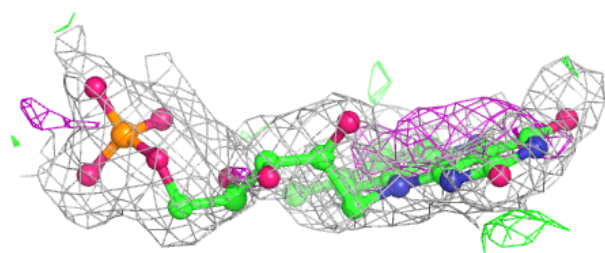
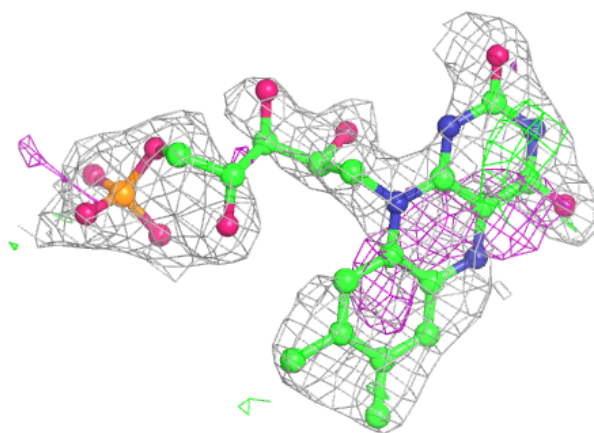
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NA	B	401	1/1	0.66	0.27	61,61,61,61	0
5	FLC	E	403	13/13	0.72	0.25	40,49,56,63	0
3	FMN	F	402	31/31	0.76	0.26	30,56,67,76	0
2	NA	F	401	1/1	0.76	0.33	56,56,56,56	0
2	NA	E	401	1/1	0.77	0.22	51,51,51,51	0
5	FLC	C	404	13/13	0.78	0.26	46,53,61,67	0
4	PG4	F	404	13/13	0.78	0.24	36,47,59,61	0
5	FLC	D	403	13/13	0.79	0.25	48,51,58,59	0
4	PG4	D	405	13/13	0.79	0.23	32,42,54,61	0
3	FMN	B	402	31/31	0.80	0.29	33,53,70,77	0
3	FMN	C	402	31/31	0.81	0.26	35,53,69,76	0
4	PG4	B	404	13/13	0.82	0.24	34,46,51,63	0
3	FMN	E	402	31/31	0.83	0.24	25,44,60,64	0
5	FLC	C	403	13/13	0.84	0.30	43,53,59,60	0
4	PG4	D	404	13/13	0.84	0.22	33,42,53,54	0
3	FMN	A	402	31/31	0.85	0.27	24,43,61,62	0
4	PG4	A	403	13/13	0.85	0.22	42,49,58,61	0
4	PG4	C	405	13/13	0.86	0.19	35,45,57,58	0
2	NA	A	401	1/1	0.86	0.20	45,45,45,45	0
4	PG4	E	404	13/13	0.87	0.20	33,42,56,57	0
4	PG4	C	406	13/13	0.87	0.21	43,50,57,62	0
4	PG4	B	405	13/13	0.88	0.21	43,47,54,56	0
4	PG4	F	405	13/13	0.88	0.26	40,50,56,57	0
3	FMN	D	402	31/31	0.88	0.24	21,39,49,56	0
5	FLC	F	403	13/13	0.88	0.19	40,46,51,53	0
4	PG4	E	405	13/13	0.90	0.21	33,46,54,55	0
2	NA	D	401	1/1	0.90	0.29	44,44,44,44	0
5	FLC	B	403	13/13	0.91	0.18	38,47,50,52	0
4	PG4	A	404	13/13	0.91	0.23	40,43,50,51	0
2	NA	C	401	1/1	0.95	0.27	53,53,53,53	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.

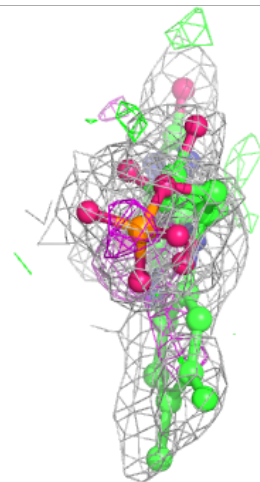
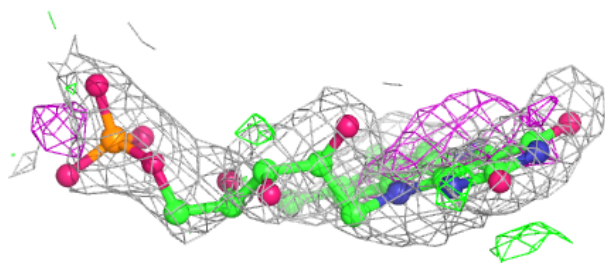
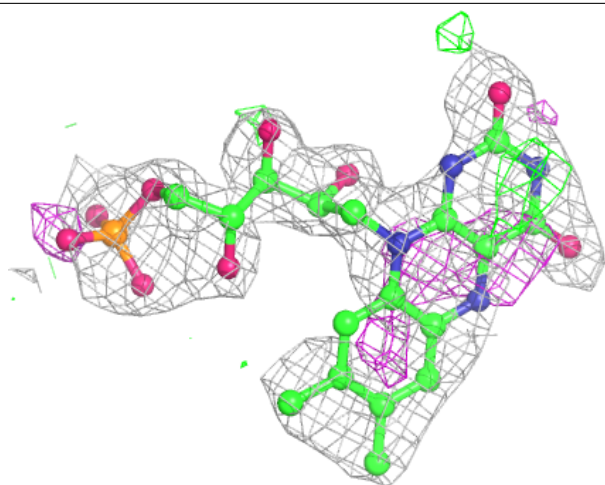
Electron density around FMN F 402:

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



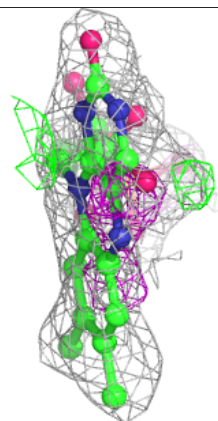
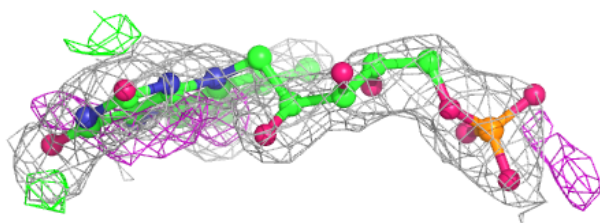
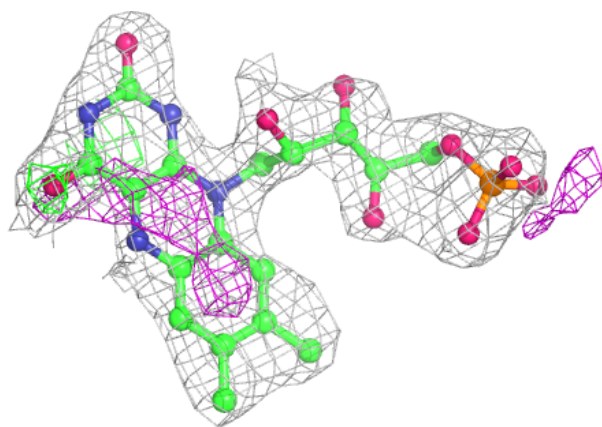
Electron density around FMN B 402:

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



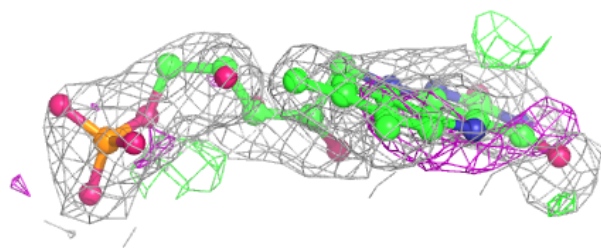
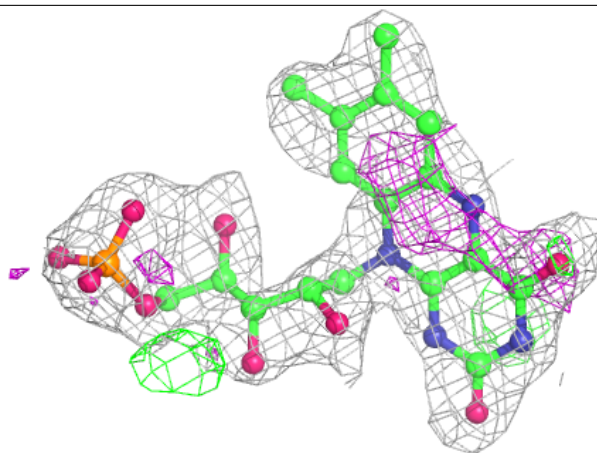
Electron density around FMN C 402:

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



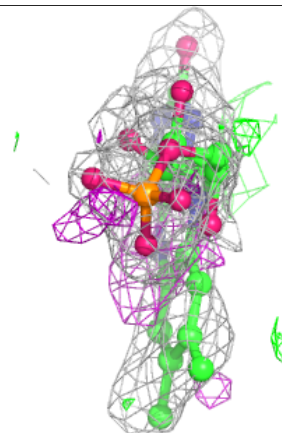
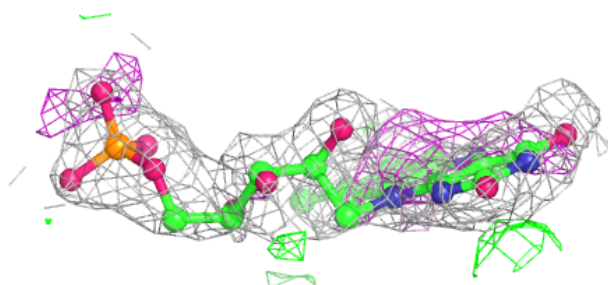
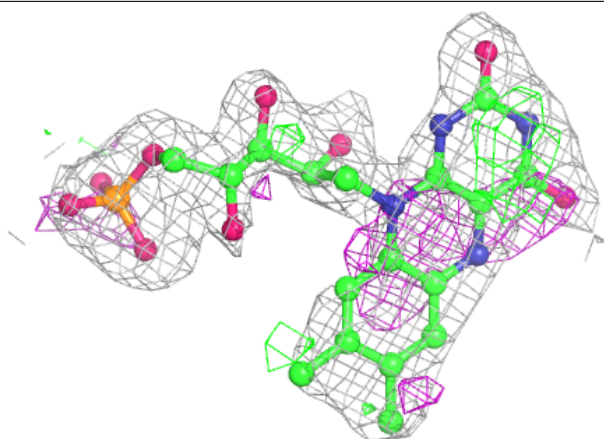
Electron density around FMN E 402:

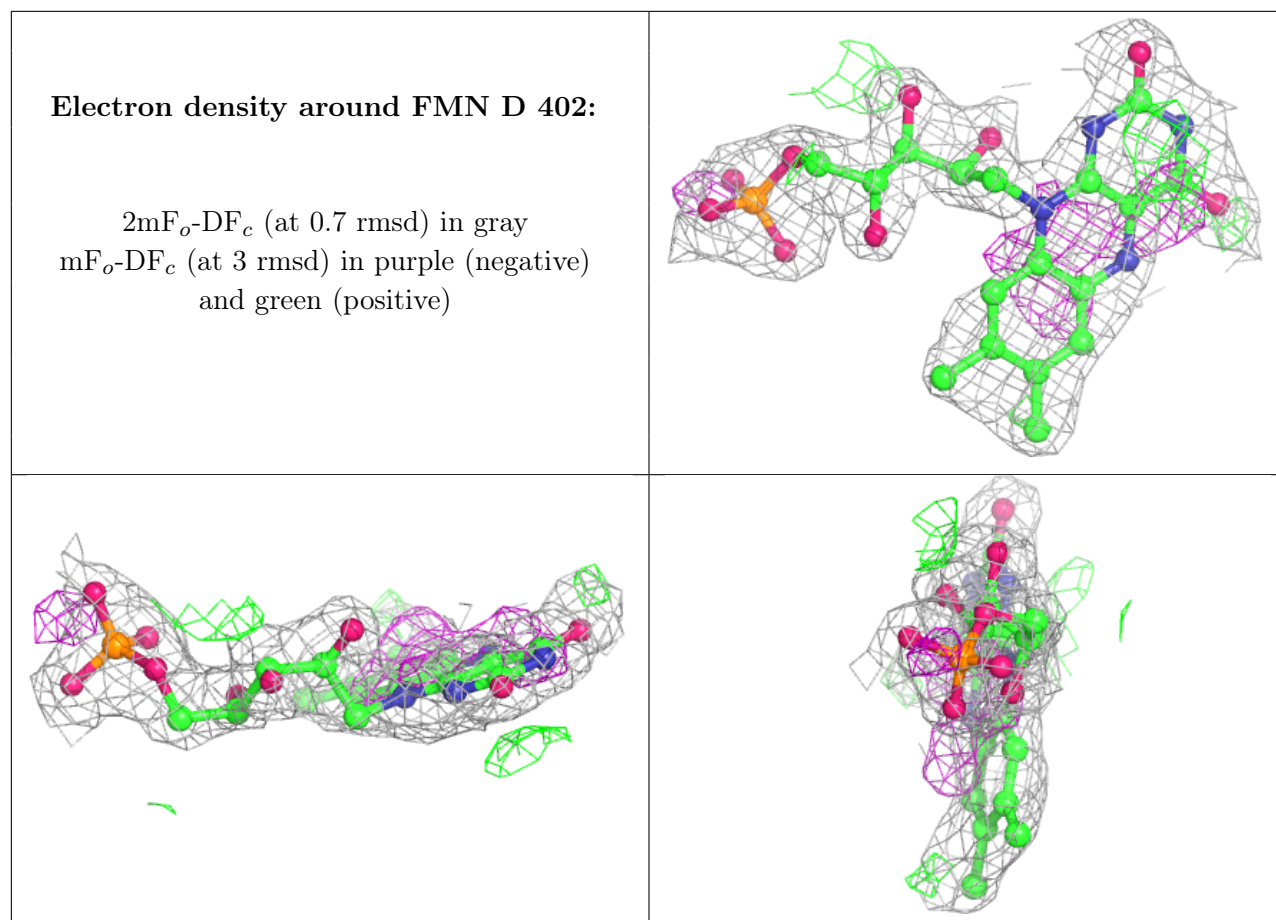
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FMN A 402:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.