W O R L D W I D E
wwPDB
P R O T E I N   D A T A   B A N K

wwPDB X-ray Structure Validation Summary Report (i)

May 16, 2020 – 04:28 pm BST

|		
PDB ID	:	2HPV
Title	:	Crystal structure of FMN-Dependent azoreductase from Enterococcus faecalis
Authors	:	Liu, Z.J.; Chen, L.; Chen, H.; Rose, J.; Wang, B.C.; Southeast Collaboratory for Structural Genomics (SECSG)
Deposited on	:	2006-07-17
Resolution	:	2.00 Å(reported)

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.
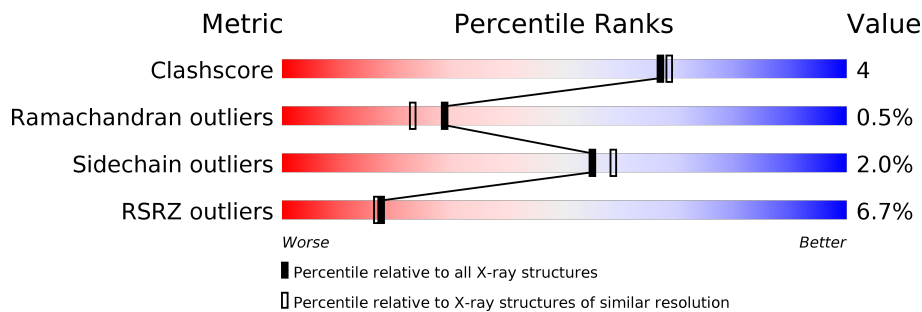
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=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 <=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	208	
1	B	208	
1	C	208	
1	D	208	

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 7142 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 FMN-dependent NADH-azoreductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	207	Total 1631	C 1039	N 270	O 319	Se 3	0	0	0
1	B	207	Total 1631	C 1039	N 270	O 319	Se 3	0	0	0
1	C	207	Total 1631	C 1039	N 270	O 319	Se 3	0	0	0
1	D	207	Total 1631	C 1039	N 270	O 319	Se 3	0	0	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
A	51	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
A	104	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
A	198	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
B	1	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
B	51	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
B	104	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
B	198	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
C	1	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
C	51	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
C	104	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
C	198	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
D	1	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
D	51	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
D	104	MSE	MET	MODIFIED RESIDUE	UNP Q831B2
D	198	MSE	MET	MODIFIED RESIDUE	UNP Q831B2

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: $C_{17}H_{21}N_4O_9P$).

FMN

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
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		

- Molecule 3 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	116	Total	O	0	0
			116	116		
3	B	127	Total	O	0	0
			127	127		
3	C	124	Total	O	0	0
			124	124		
3	D	127	Total	O	0	0
			127	127		

3 Residue-property plots (i)

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.
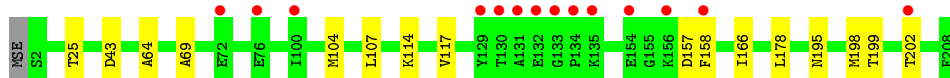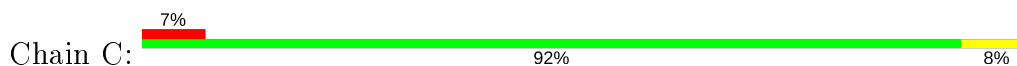
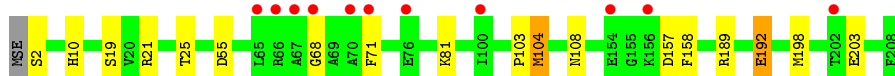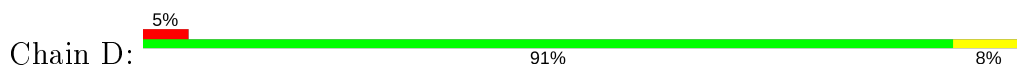- Molecule 1: FMN-dependent NADH-azoreductase
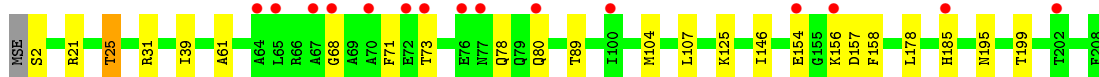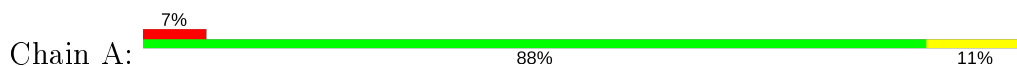
- Molecule 1: FMN-dependent NADH-azoreductase
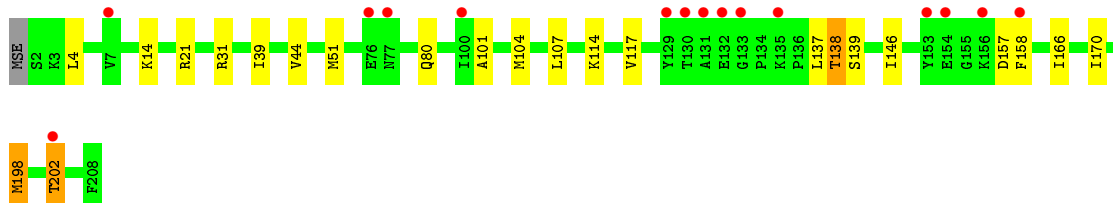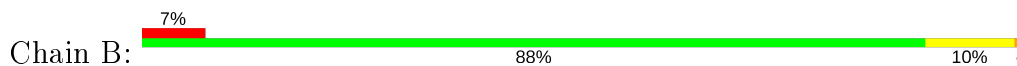
- Molecule 1: FMN-dependent NADH-azoreductase

- Molecule 1: FMN-dependent NADH-azoreductase

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	98.04Å 99.63Å 106.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.78 – 2.00 25.78 – 2.00	Depositor EDS
% Data completeness (in resolution range)	96.3 (25.78-2.00) 92.9 (25.78-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$<I/\sigma(I)>$ [1]	1.31 (at 1.99Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.199 , 0.250 0.200 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	25.5	Xtriage
Anisotropy	0.359	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 45.5	EDS
L-test for twinning[2]	$<|L|> = 0.49, <L^2> = 0.32$	Xtriage
Estimated twinning fraction	0.015 for k,h,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	7142	wwPDB-VP
Average B, all atoms (Å^2)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.89% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.

[2]Theoretical values of $<|L|>$, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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	0.92	0/1662	0.81	0/2248
1	B	0.93	0/1662	0.80	1/2248 (0.0%)
1	C	0.94	0/1662	0.79	1/2248 (0.0%)
1	D	0.95	3/1662 (0.2%)	0.80	3/2248 (0.1%)
All	All	0.93	3/6648 (0.0%)	0.80	5/8992 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	2	SER	CB-OG	-5.97	1.34	1.42
1	D	203	GLU	CG-CD	5.14	1.59	1.51
1	D	2	SER	N-CA	5.03	1.56	1.46

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	2	SER	N-CA-C	9.92	137.79	111.00
1	D	21	ARG	NE-CZ-NH1	6.19	123.39	120.30
1	B	198	MSE	CG-SE-CE	-6.16	85.36	98.90
1	D	21	ARG	NE-CZ-NH2	-6.08	117.26	120.30
1	C	43	ASP	CB-CG-OD1	5.45	123.20	118.30

There are no chirality outliers.

There are no planarity outliers.

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	1631	0	1604	13	0
1	B	1631	0	1604	17	0
1	C	1631	0	1604	18	0
1	D	1631	0	1604	11	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
3	A	116	0	0	0	0
3	B	127	0	0	1	0
3	C	124	0	0	3	0
3	D	127	0	0	0	0
All	All	7142	0	6492	51	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 4.

The worst 5 of 51 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:198:MSE:CE	1:B:198:MSE:SE	2.21	1.39
1:C:158:PHE:CD2	1:D:158:PHE:HE1	1.58	1.22
1:C:158:PHE:CD2	1:D:158:PHE:CE1	2.47	1.03
1:C:158:PHE:HD2	1:D:158:PHE:CE1	1.85	0.93
1:A:104:MSE:HE1	1:A:157:ASP:HB2	1.51	0.92

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	205/208 (99%)	197 (96%)	6 (3%)	2 (1%)	15	9
1	B	205/208 (99%)	202 (98%)	3 (2%)	0	100	100
1	C	205/208 (99%)	202 (98%)	3 (2%)	0	100	100
1	D	205/208 (99%)	198 (97%)	5 (2%)	2 (1%)	15	9
All	All	820/832 (99%)	799 (97%)	17 (2%)	4 (0%)	29	23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	68	GLY
1	A	71	PHE
1	D	71	PHE
1	A	68	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	174/171 (102%)	170 (98%)	4 (2%)	50	53
1	B	174/171 (102%)	168 (97%)	6 (3%)	37	36
1	C	174/171 (102%)	173 (99%)	1 (1%)	86	90
1	D	174/171 (102%)	171 (98%)	3 (2%)	60	65
All	All	696/684 (102%)	682 (98%)	14 (2%)	55	58

5 of 14 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	137	LEU
1	B	138	THR
1	D	19	SER
1	B	114	LYS
1	C	178	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	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)

4 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	B	1201	-	31,33,33	1.38	4 (12%)	40,50,50	2.30	8 (20%)
2	FMN	C	1202	-	31,33,33	1.47	6 (19%)	40,50,50	2.39	8 (20%)
2	FMN	D	1203	-	31,33,33	1.59	5 (16%)	40,50,50	2.04	10 (25%)
2	FMN	A	1200	-	31,33,33	1.74	7 (22%)	40,50,50	2.09	8 (20%)

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	B	1201	-	-	0/18/18/18	0/3/3/3
2	FMN	C	1202	-	-	0/18/18/18	0/3/3/3
2	FMN	D	1203	-	-	0/18/18/18	0/3/3/3
2	FMN	A	1200	-	-	2/18/18/18	0/3/3/3

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1200	FMN	C10-N1	4.08	1.38	1.33
2	D	1203	FMN	C4A-C10	3.98	1.42	1.38
2	C	1202	FMN	C4-N3	3.85	1.39	1.33
2	A	1200	FMN	C4A-C10	3.84	1.42	1.38
2	B	1201	FMN	C4-N3	3.75	1.39	1.33

The worst 5 of 34 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	1202	FMN	C1'-N10-C9A	9.37	125.67	118.29
2	B	1201	FMN	C4-N3-C2	8.89	122.65	115.14
2	A	1200	FMN	C4-N3-C2	8.45	122.28	115.14
2	C	1202	FMN	C4-N3-C2	7.95	121.85	115.14
2	D	1203	FMN	C4-N3-C2	6.06	120.26	115.14

There are no chirality outliers.

All (2) torsion outliers are listed below:

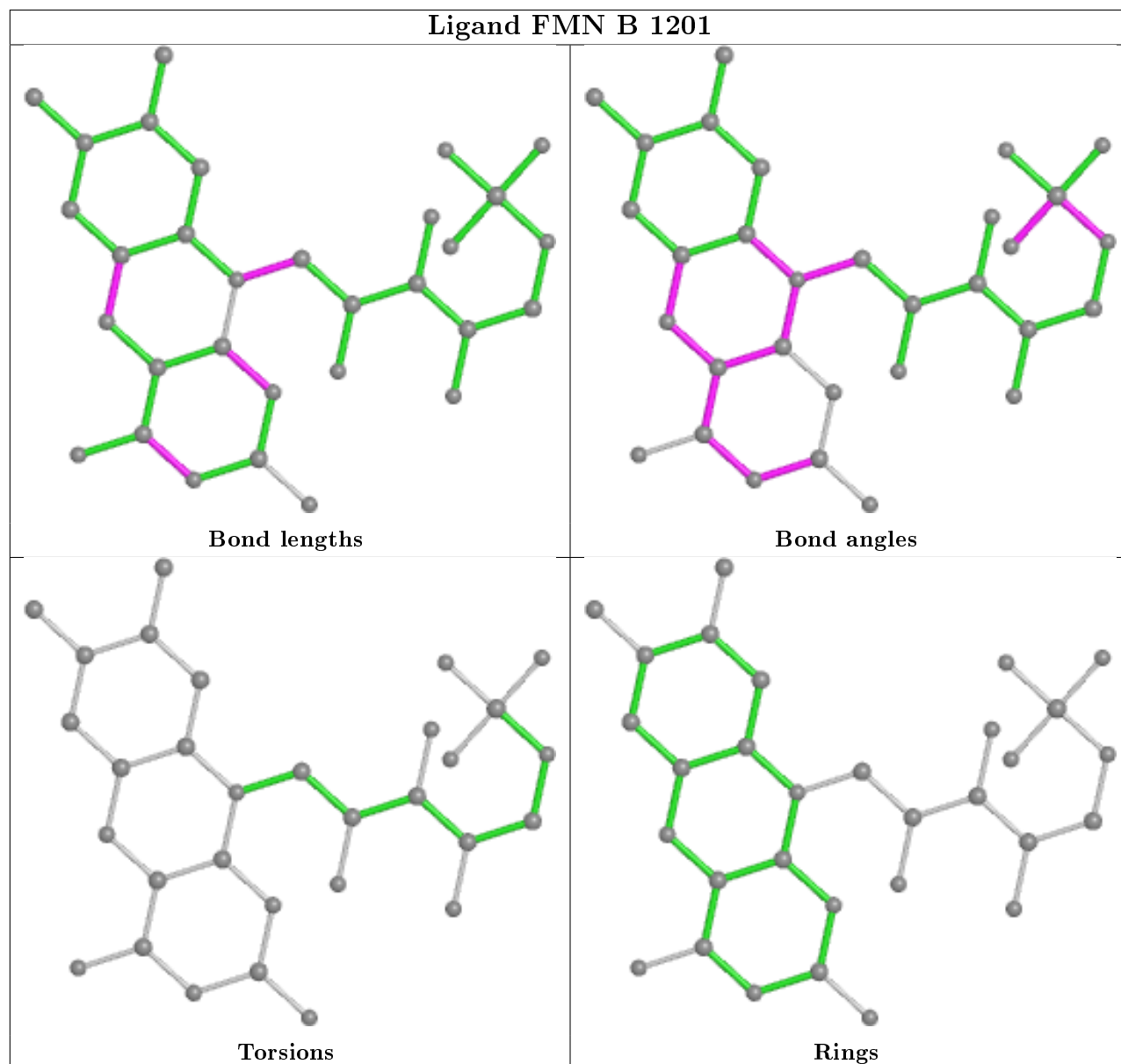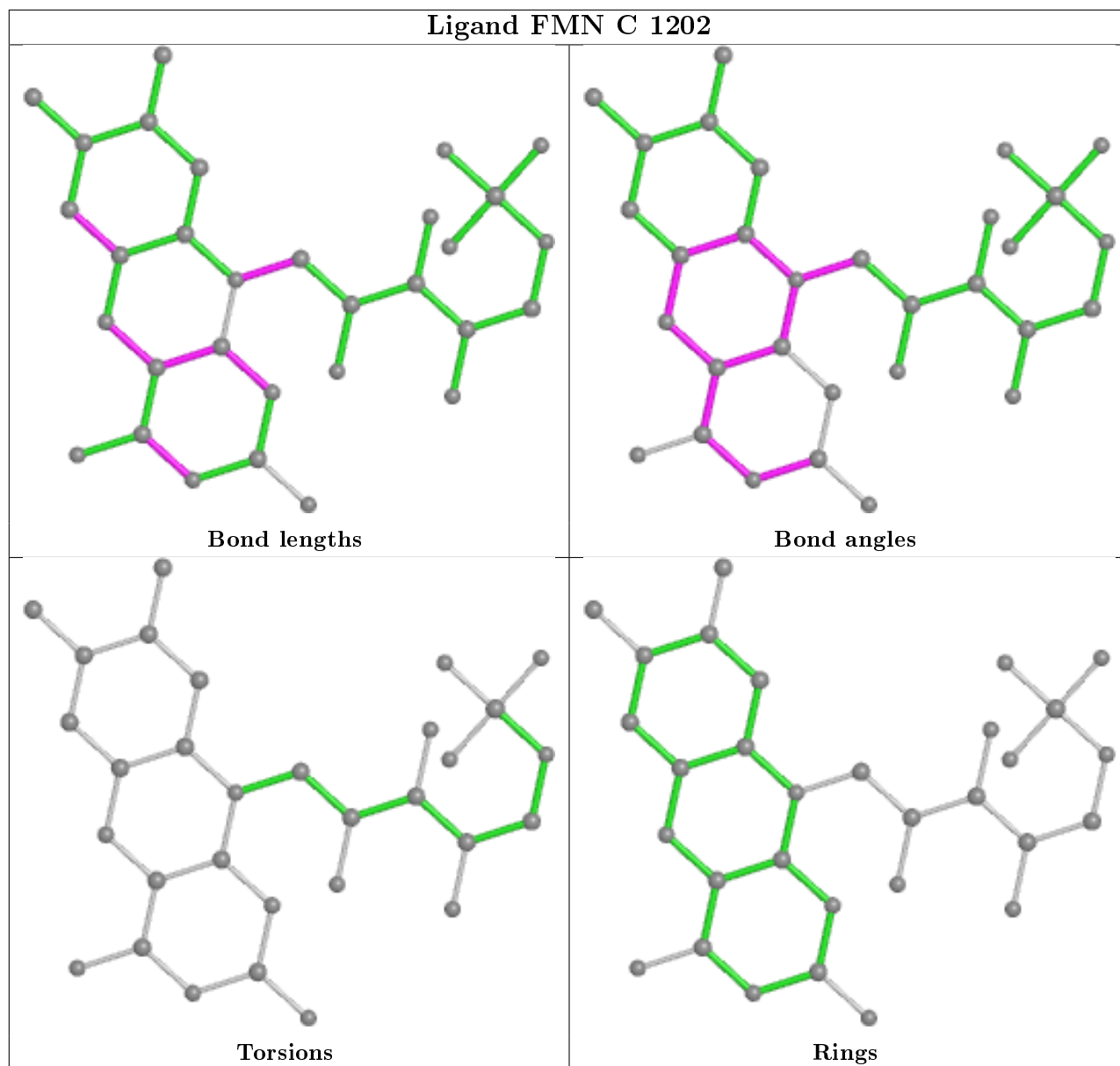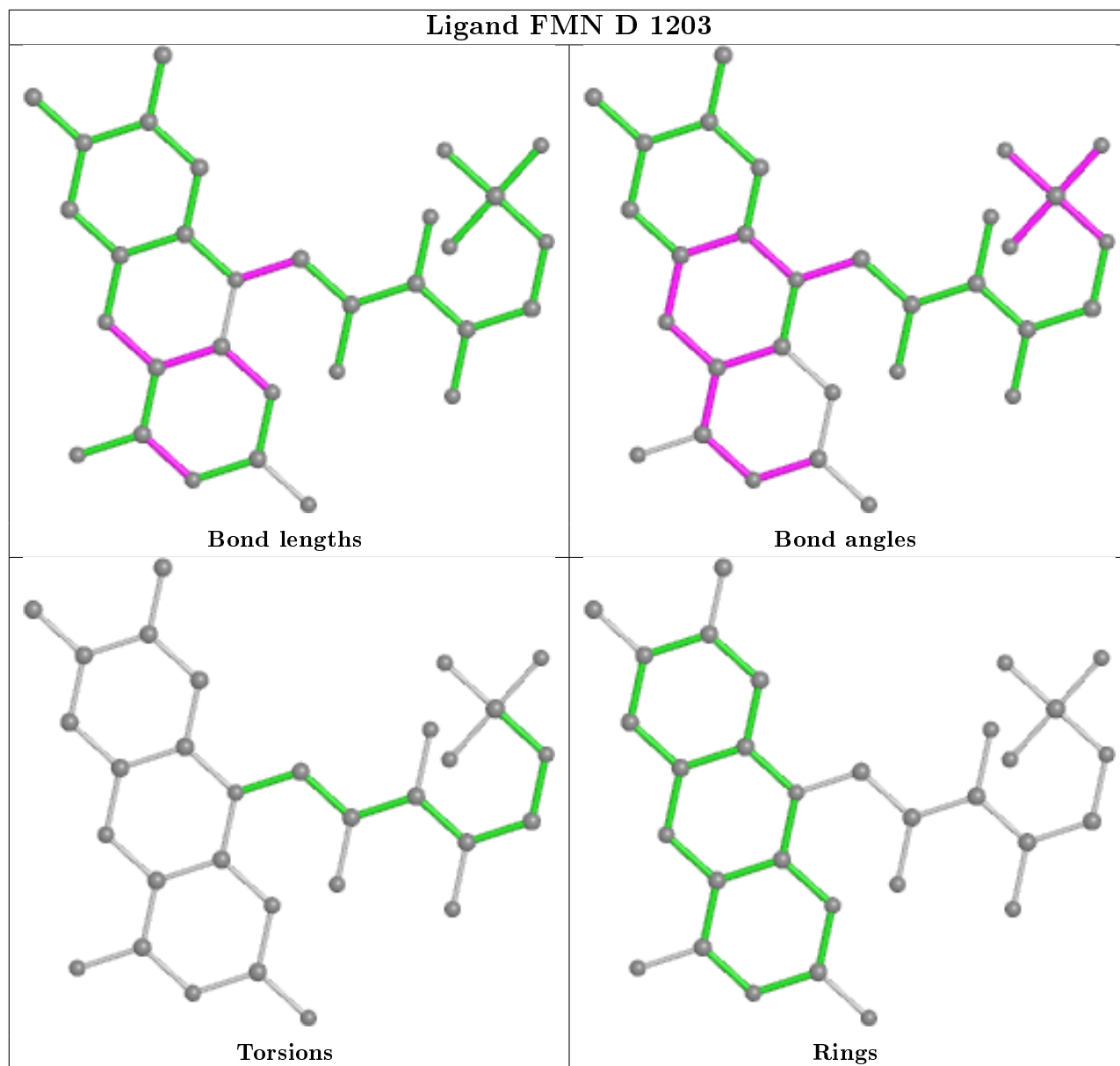| Mol	Chain	Res	Type	Atoms
2	A	1200	FMN	O3'-C3'-C4'-O4'
2	A	1200	FMN	C2'-C3'-C4'-O4'

There are no ring outliers.

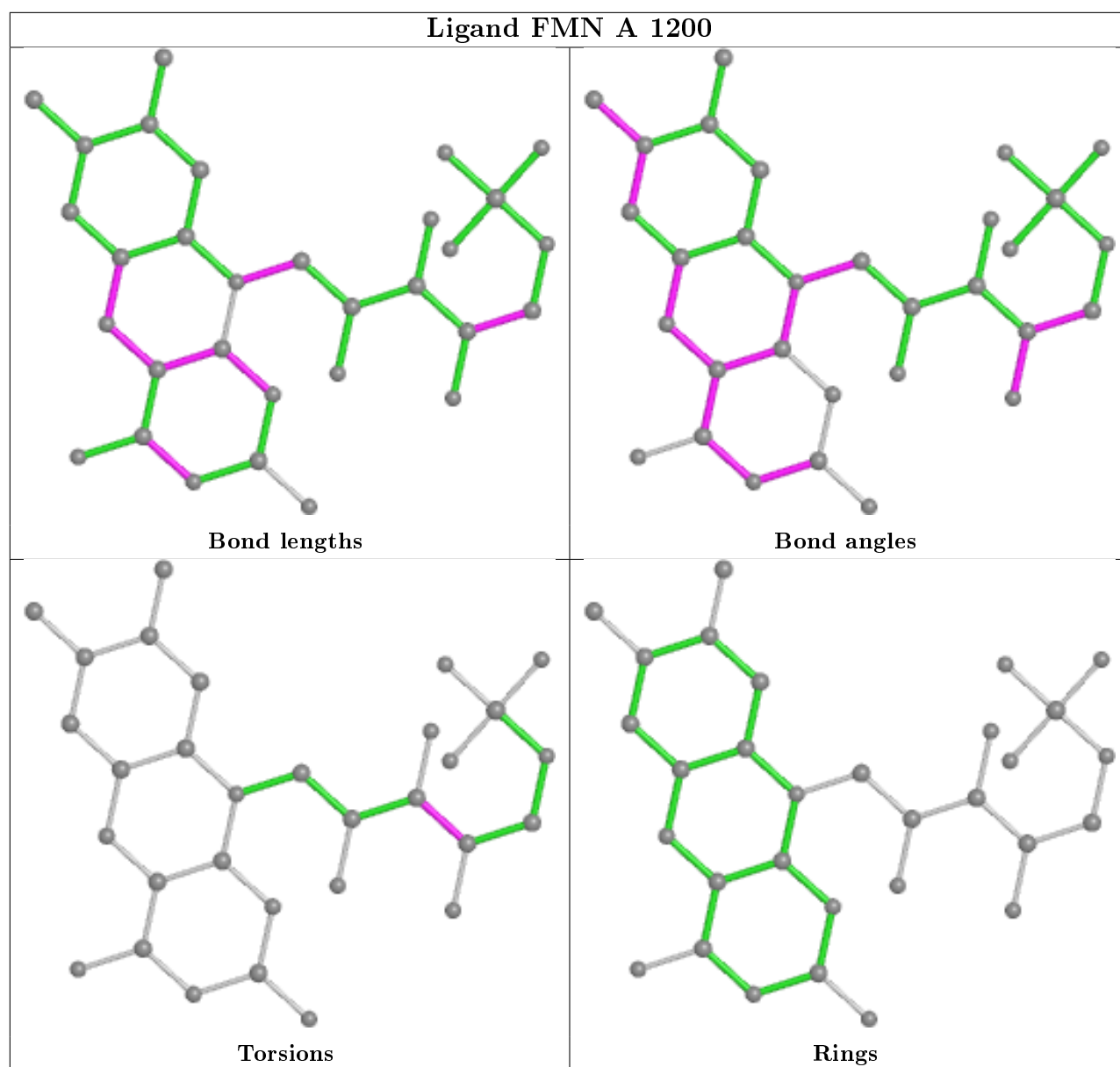No monomer is involved in short contacts.

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 then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

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

Ligand FMN B 1201

Bond lengths

Bond angles

Torsions

Rings

Ligand FMN C 1202

Bond lengths

Bond angles

Torsions

Rings

Ligand FMN D 1203

Bond lengths

Bond angles

Torsions

Rings

Ligand FMN A 1200

Bond lengths

Bond angles

Torsions

Rings

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(Å²)	Q<0.9
1	A	204/208 (98%)	0.20	15 (7%)	14	13	17, 27, 49, 57	0
1	B	204/208 (98%)	0.14	15 (7%)	14	13	15, 25, 45, 50	0
1	C	204/208 (98%)	0.14	14 (6%)	16	16	14, 25, 46, 52	0
1	D	204/208 (98%)	0.16	11 (5%)	25	24	14, 25, 47, 55	0
All	All	816/832 (98%)	0.16	55 (6%)	17	17	14, 25, 47, 57	0

The worst 5 of 55 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	131	ALA	5.0
1	B	132	GLU	5.0
1	B	131	ALA	4.8
1	A	67	ALA	4.8
1	A	156	LYS	4.3

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

There are no non-standard protein/DNA/RNA residues in this entry.
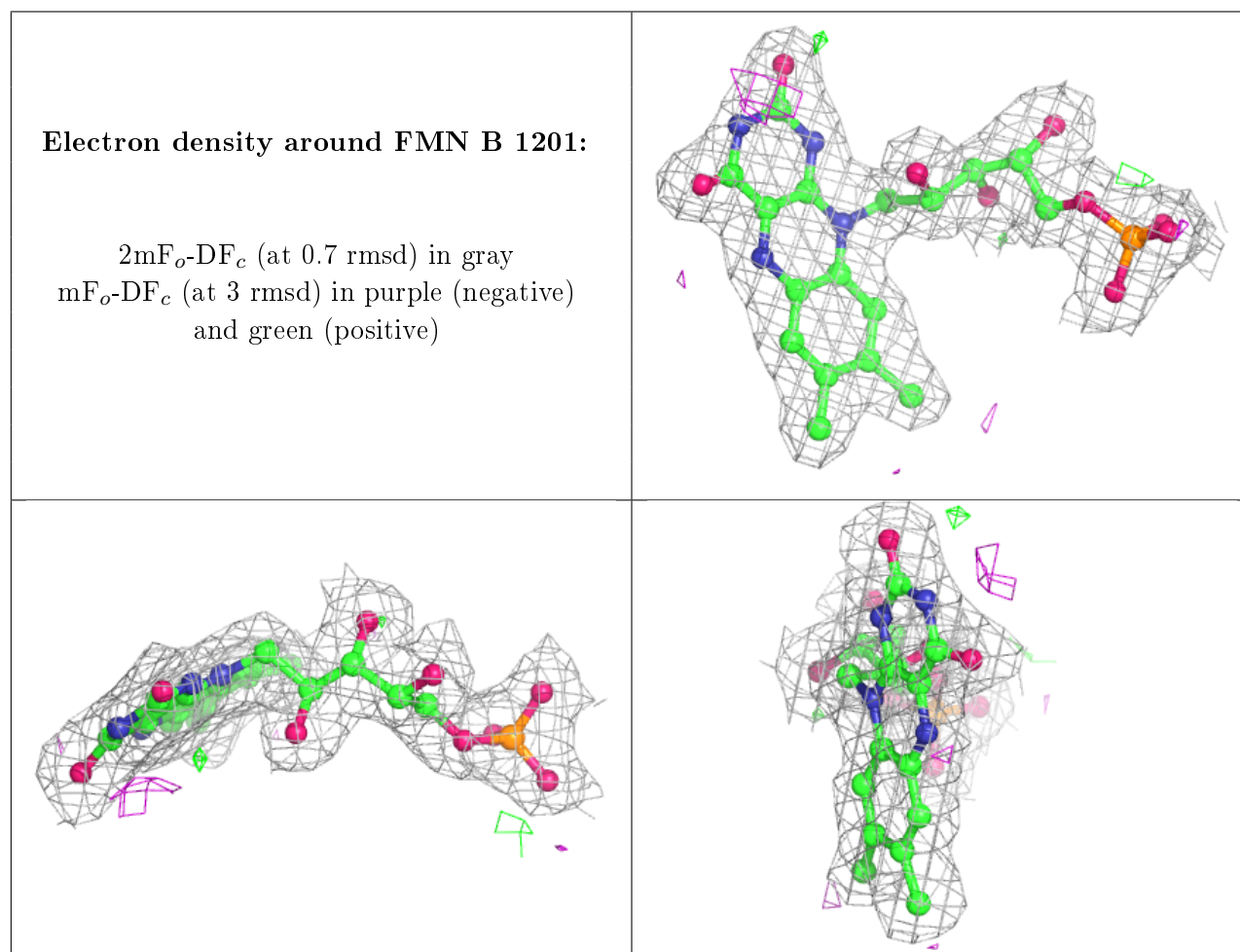
6.3 Carbohydrates (i)

There are no 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, 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	FMN	B	1201	31/31	0.97	0.09	14,19,22,23	0
2	FMN	C	1202	31/31	0.97	0.10	14,19,21,23	0
2	FMN	D	1203	31/31	0.97	0.11	15,24,27,28	0
2	FMN	A	1200	31/31	0.97	0.09	16,22,25,26	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.
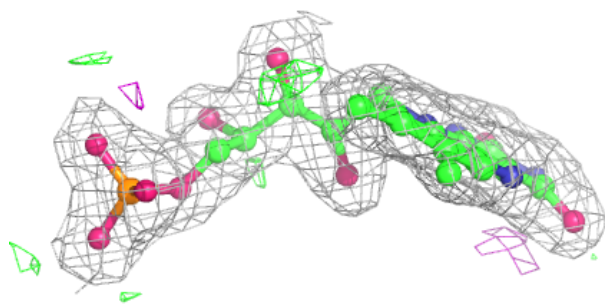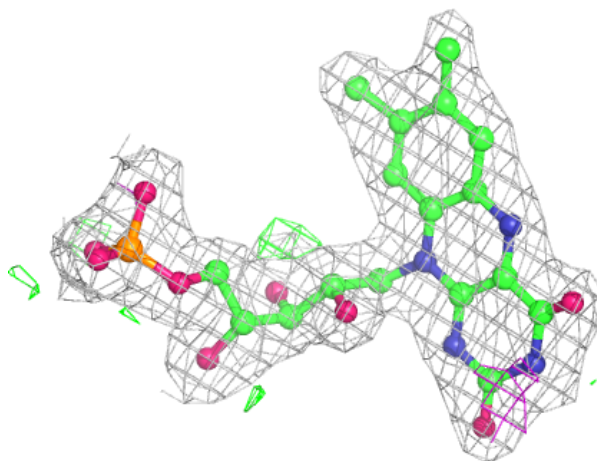
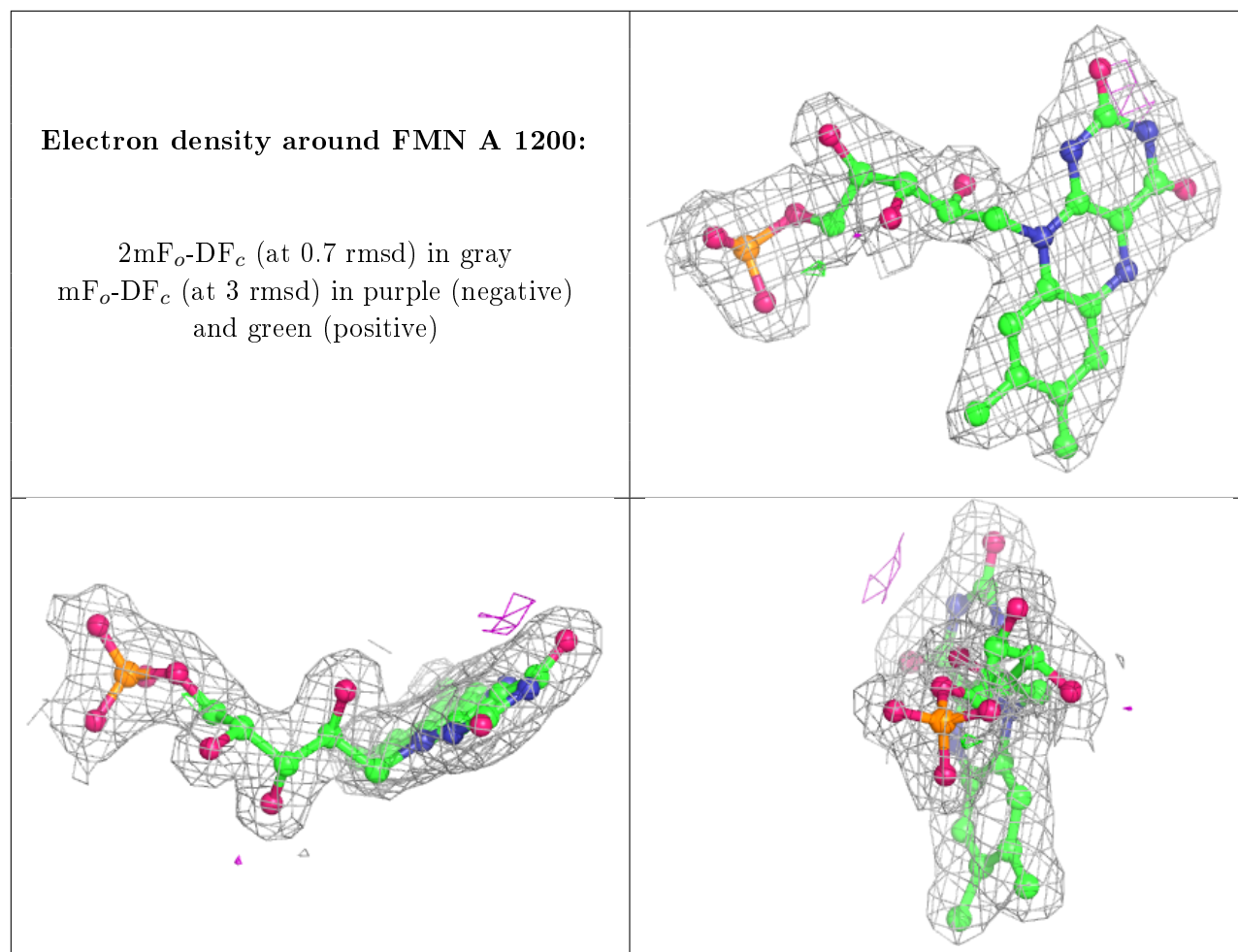

Electron density around FMN B 1201:

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

Electron density around FMN C 1202:

$2mF_o\text{-}DF_c$ (at 0.7 rmsd) in gray
$mF_o\text{-}DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)
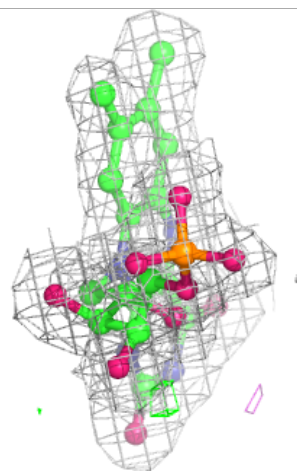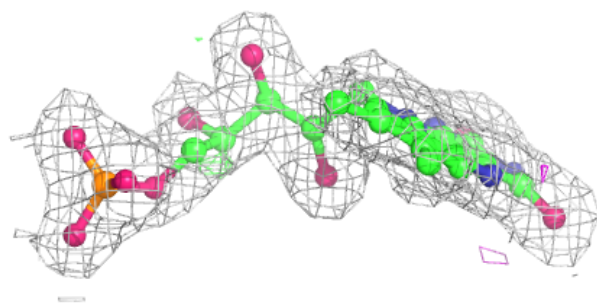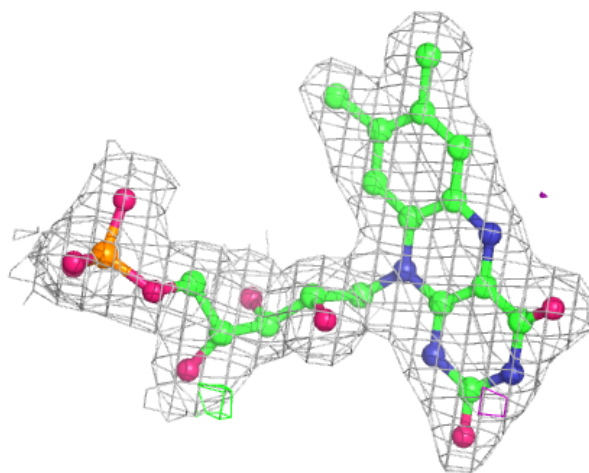
Electron density around FMN D 1203:

$2mF_o$-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 1200:

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

6.5 Other polymers ⓘ

There are no such residues in this entry.