Full wwPDB X-ray Structure Validation Report (i)

May 13, 2020 – 08:01 pm BST

PDB ID	:	5XDD
Title	:	Crystal structure of tertiary complex of TdsC from Paenibacillus sp. A11-2 with FMN and Indole
Authors	:	Hino, T.; Hamamoto, H.; Ohshiro, T.; Nagano, S.
Deposited on	:	2017-03-28
Resolution	:	1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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 (i) symbol.

The following versions of software and data (see references (i)) 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 (i)

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

The reported resolution of this entry is 1.90 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.
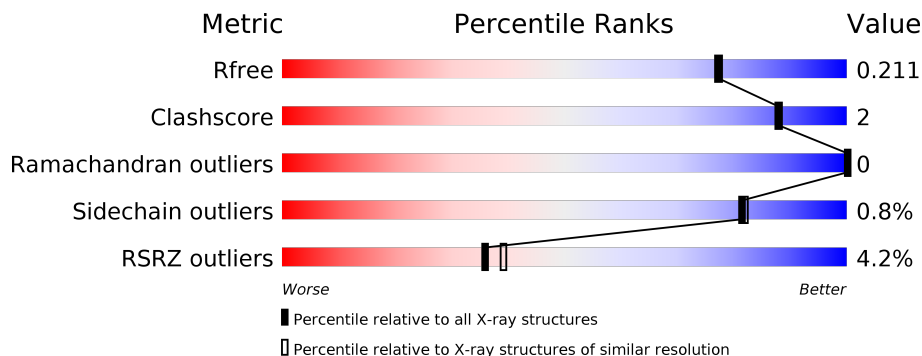
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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.

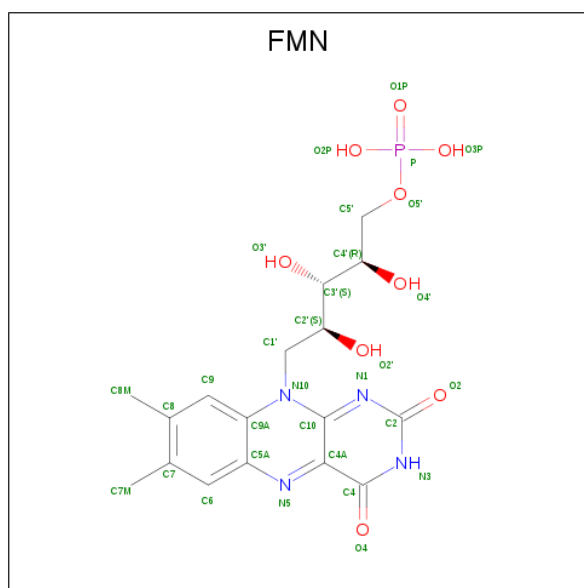| Mol	Chain	Length	Quality of chain
1	A	414	6% / 92% 5% ·
1	B	414	5% / 90% 5% 5%
1	C	414	3% / 92% 5% ·
1	D	414	2% / 94% · ·

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 26204 atoms, of which 12227 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 Thermophilic dibenzothiophene desulfurization enzyme C.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	400	Total	C	H	N	O	S	0	5	0
			6189	1984	3047	568	586	4			
1	B	394	Total	C	H	N	O	S	0	0	0
			6061	1945	2987	552	573	4			
1	C	401	Total	C	H	N	O	S	0	0	0
			6179	1980	3045	566	584	4			
1	D	400	Total	C	H	N	O	S	0	3	0
			6190	1981	3053	567	584	5			

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

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	H	N	O	P	0	0
			49	17	18	4	9	1		
2	B	1	Total	C	H	N	O	P	0	0
			49	17	18	4	9	1		

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Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
			Total	C	H	N	O	P		
2	C	1	49	17	18	4	9	1	0	0
			Total	C	H	N	O	P		
2	D	1	50	17	19	4	9	1	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).

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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total 5	O 4	S 1	0	0
3	D	1	Total 5	O 4	S 1	0	0
3	D	1	Total 5	O 4	S 1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).
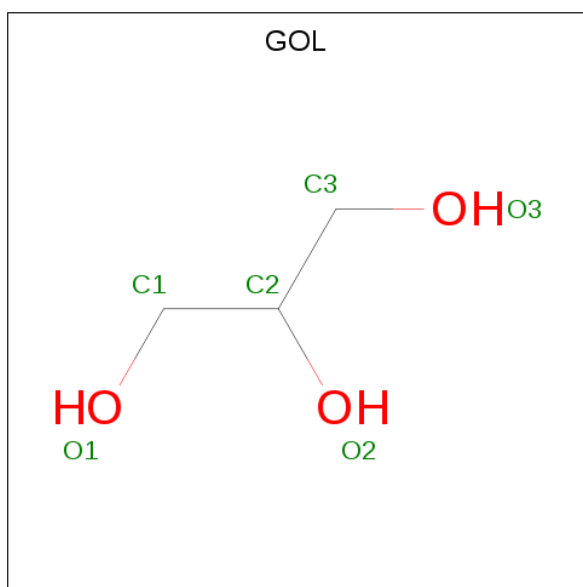
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total 14	C 3	H 8	O 3	0	0

- Molecule 5 is INDOLE (three-letter code: IND) (formula: C_8H_7N).
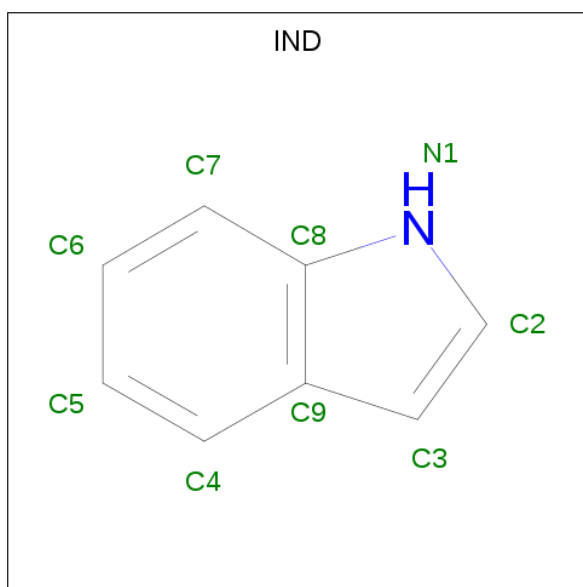
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	C	1	Total 16	C 8	H 7	N 1	0	0
5	D	1	Total 16	C 8	H 7	N 1	0	0

- Molecule 6 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	297	Total 298	O 298	0	1
6	B	333	Total 334	O 334	0	1
6	C	305	Total 306	O 306	0	1
6	D	343	Total 344	O 344	0	1

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.
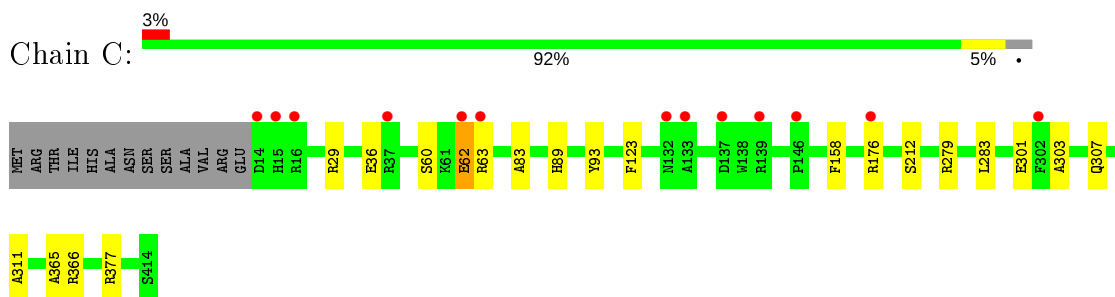
- Molecule 1: Thermophilic dibenzothiophene desulfurization enzyme C
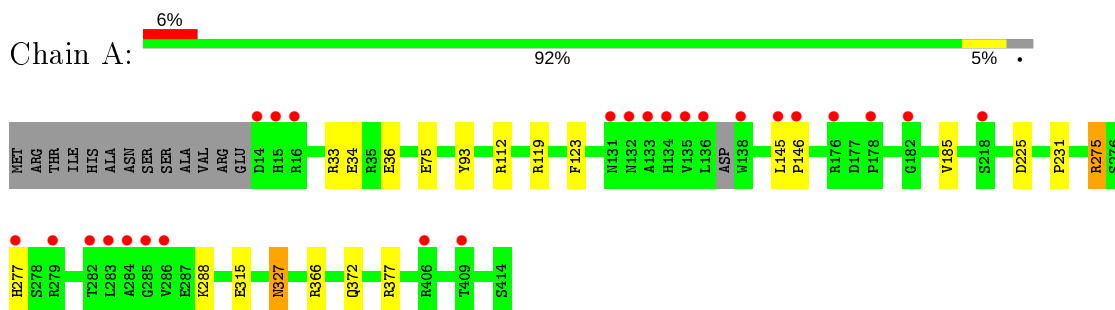
- Molecule 1: Thermophilic dibenzothiophene desulfurization enzyme C
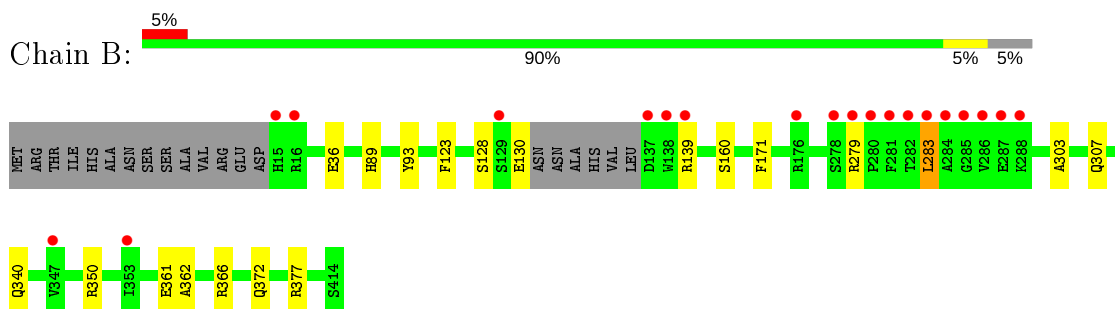
- Molecule 1: Thermophilic dibenzothiophene desulfurization enzyme C

- Molecule 1: Thermophilic dibenzothiophene desulfurization enzyme C

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α, β, γ	100.79Å 100.79Å 424.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.47 – 1.90 47.47 – 1.81	Depositor EDS
% Data completeness (in resolution range)	92.6 (47.47-1.90) 89.4 (47.47-1.81)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	1.11 (at 1.81Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155)	Depositor
R, R_{free}	0.178 , 0.211 0.179 , 0.211	Depositor DCC
R_{free} test set	8952 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	18.6	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.43 , 50.3	EDS
L-test for twinning[2]	$<|L|> = 0.48$, $<L^2> = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	26204	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.11% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section:
FMN, GOL, IND, SO4

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	0.54	0/3238	0.66	1/4401 (0.0%)
1	B	0.55	0/3150	0.64	1/4281 (0.0%)
1	C	0.52	0/3212	0.61	0/4367
1	D	0.52	0/3228	0.66	1/4386 (0.0%)
All	All	0.53	0/12828	0.64	3/17435 (0.0%)

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	C	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	275	ARG	NE-CZ-NH2	-6.15	117.22	120.30
1	D	355	LEU	CA-CB-CG	5.65	128.30	115.30
1	B	283	LEU	CA-CB-CG	5.52	128.00	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	62	GLU	Peptide

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	3142	3047	3030	12	0
1	B	3074	2987	2985	12	0
1	C	3134	3045	3045	17	0
1	D	3137	3053	3038	7	0
2	A	31	18	19	0	0
2	B	31	18	19	1	0
2	C	31	18	19	0	0
2	D	31	19	19	1	0
3	A	15	0	0	1	0
3	B	10	0	0	0	0
3	C	15	0	0	0	0
3	D	20	0	0	1	0
4	B	6	8	8	0	0
5	C	9	7	7	0	0
5	D	9	7	7	1	0
6	A	298	0	0	4	1
6	B	334	0	0	7	1
6	C	306	0	0	3	0
6	D	344	0	0	7	0
All	All	13977	12227	12196	50	1

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.

All (50) 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:D:211:ASP:OD1	6:D:601:HOH:O	2.00	0.78
1:B:361:GLU:OE2	6:B:601:HOH:O	2.03	0.76
1:A:327[A]:ASN:ND2	3:A:503:SO4:O4	2.23	0.71
1:B:130:GLU:OE1	1:B:139:ARG:N	2.21	0.69
1:C:301:GLU:OE2	6:C:601:HOH:O	2.13	0.66
3:D:503:SO4:O4	6:D:602:HOH:O	2.15	0.64
1:C:303:ALA:O	1:C:307:GLN:HG2	2.01	0.60
1:C:60:SER:C	1:C:62:GLU:H	2.07	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:SER:HB3	1:B:171:PHE:CD1	2.39	0.57
6:B:634:HOH:O	1:C:307:GLN:HG3	2.04	0.57
1:B:350:ARG:NH2	6:B:604:HOH:O	2.28	0.57
1:D:89:HIS:CE1	1:D:93:TYR:CE1	2.93	0.56
1:C:89:HIS:CE1	1:C:93:TYR:CE1	2.94	0.56
1:B:89:HIS:CE1	1:B:93:TYR:CE1	2.95	0.55
1:C:365:ALA:HB3	2:D:501:FMN:H5'2	1.88	0.55
1:D:361:GLU:OE2	6:D:603:HOH:O	2.19	0.53
1:C:279:ARG:HG3	1:C:366:ARG:NH2	2.22	0.53
1:A:119[A]:ARG:HD3	6:A:601:HOH:O	2.09	0.52
1:A:36:GLU:HG2	1:A:377:ARG:HD3	1.91	0.52
1:B:340:GLN:HA	6:B:634:HOH:O	2.08	0.52
1:C:62:GLU:O	1:C:63:ARG:CB	2.58	0.51
1:C:62:GLU:O	1:C:63:ARG:HG3	2.11	0.51
1:B:372:GLN:OE1	6:B:602:HOH:O	2.19	0.51
1:C:29:ARG:HG3	1:C:83:ALA:HB1	1.95	0.48
1:D:37:ARG:NH1	6:D:616:HOH:O	2.44	0.48
1:A:275:ARG:NH2	6:A:606:HOH:O	2.37	0.48
5:D:506:IND:H2	6:D:635:HOH:O	2.14	0.48
1:A:112:ARG:NH2	1:A:225:ASP:HB2	2.30	0.47
1:A:277:HIS:O	1:A:366:ARG:NH2	2.48	0.47
1:B:36:GLU:HG2	1:B:377:ARG:HD3	1.97	0.46
1:D:121:ARG:NH2	6:D:608:HOH:O	2.32	0.46
1:D:288:LYS:NZ	6:D:622:HOH:O	2.49	0.46
1:C:62:GLU:O	1:C:63:ARG:CG	2.64	0.45
1:C:36:GLU:HG2	1:C:377:ARG:HD3	1.99	0.45
1:B:279:ARG:HD2	1:B:366:ARG:CZ	2.47	0.45
1:A:75:GLU:OE2	6:A:602:HOH:O	2.21	0.44
6:B:634:HOH:O	1:C:307:GLN:CB	2.66	0.43
1:A:315[A]:GLU:OE2	6:A:603:HOH:O	2.21	0.43
1:A:33:ARG:HD2	1:A:372:GLN:O	2.19	0.43
1:A:33:ARG:NH1	1:A:34:GLU:OE2	2.52	0.43
1:C:176:ARG:NH2	6:C:614:HOH:O	2.52	0.43
1:A:145:LEU:HB3	1:A:146:PRO:HD2	2.01	0.42
1:B:362:ALA:O	6:B:603:HOH:O	2.21	0.42
1:B:160:SER:HB2	2:B:501:FMN:O4	2.19	0.42
1:A:185:VAL:HG12	1:A:231:PRO:HB3	2.01	0.41
1:B:303:ALA:O	1:B:307:GLN:HG2	2.19	0.41
1:C:311:ALA:O	6:C:602:HOH:O	2.21	0.41
1:C:158:PHE:HA	1:C:212:SER:O	2.21	0.41
1:D:36:GLU:HG2	1:D:377:ARG:HD3	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:29:ARG:HG3	1:C:83:ALA:CB	2.52	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:874:HOH:O	6:B:906:HOH:O[7_455]	1.84	0.36

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	401/414 (97%)	396 (99%)	5 (1%)	0	100	100
1	B	390/414 (94%)	387 (99%)	3 (1%)	0	100	100
1	C	399/414 (96%)	390 (98%)	9 (2%)	0	100	100
1	D	401/414 (97%)	393 (98%)	8 (2%)	0	100	100
All	All	1591/1656 (96%)	1566 (98%)	25 (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	321/331 (97%)	316 (98%)	5 (2%)	62	60
1	B	312/331 (94%)	310 (99%)	2 (1%)	86	87
1	C	319/331 (96%)	317 (99%)	2 (1%)	86	87
1	D	320/331 (97%)	318 (99%)	2 (1%)	86	87
All	All	1272/1324 (96%)	1261 (99%)	11 (1%)	81	79

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93	TYR
1	A	123	PHE
1	A	288	LYS
1	A	327[A]	ASN
1	A	327[B]	ASN
1	B	123	PHE
1	B	283	LEU
1	C	123	PHE
1	C	283	LEU
1	D	123	PHE
1	D	350	ARG

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

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)

19 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	501	-	31,33,33	1.85	6 (19%)	40,50,50	2.81	10 (25%)
3	SO4	C	502	-	4,4,4	0.29	0	6,6,6	0.37	0
2	FMN	D	501	-	31,33,33	1.80	6 (19%)	40,50,50	2.22	12 (30%)
3	SO4	A	502	-	4,4,4	0.18	0	6,6,6	0.45	0
3	SO4	D	504	-	4,4,4	0.18	0	6,6,6	0.62	0
3	SO4	A	503	-	4,4,4	0.24	0	6,6,6	0.43	0
3	SO4	D	503	-	4,4,4	0.16	0	6,6,6	0.41	0
3	SO4	B	503	-	4,4,4	0.11	0	6,6,6	0.24	0
5	IND	D	506	-	8,10,10	1.47	0	9,13,13	2.49	3 (33%)
3	SO4	A	504	-	4,4,4	0.19	0	6,6,6	0.28	0
3	SO4	C	503	-	4,4,4	0.15	0	6,6,6	0.20	0
3	SO4	C	504	-	4,4,4	0.11	0	6,6,6	0.41	0
2	FMN	C	501	-	31,33,33	1.80	5 (16%)	40,50,50	1.72	7 (17%)
3	SO4	B	502	-	4,4,4	0.17	0	6,6,6	0.39	0
4	GOL	B	504	-	5,5,5	0.59	0	5,5,5	0.27	0
3	SO4	D	505	-	4,4,4	0.18	0	6,6,6	0.18	0
3	SO4	D	502	-	4,4,4	0.15	0	6,6,6	0.34	0
2	FMN	A	501	-	31,33,33	1.69	6 (19%)	40,50,50	1.80	8 (20%)
5	IND	C	505	-	8,10,10	1.54	0	9,13,13	2.47	3 (33%)

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	501	-	-	4/18/18/18	0/3/3/3
5	IND	D	506	-	-	-	0/2/2/2
2	FMN	D	501	-	-	8/18/18/18	0/3/3/3
5	IND	C	505	-	-	-	0/2/2/2
2	FMN	A	501	-	-	4/18/18/18	0/3/3/3
2	FMN	C	501	-	-	6/18/18/18	0/3/3/3
4	GOL	B	504	-	-	0/4/4/4	-

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	FMN	C1'-N10	5.93	1.54	1.48
2	A	501	FMN	C1'-N10	5.31	1.53	1.48
2	C	501	FMN	C4A-N5	5.11	1.40	1.33
2	D	501	FMN	C10-N1	4.84	1.39	1.33
2	D	501	FMN	C1'-N10	4.74	1.53	1.48
2	C	501	FMN	C10-N1	4.74	1.39	1.33
2	B	501	FMN	C10-N1	4.62	1.39	1.33
2	C	501	FMN	C1'-N10	4.19	1.52	1.48
2	D	501	FMN	C4A-N5	4.11	1.39	1.33
2	A	501	FMN	C10-N1	4.10	1.38	1.33
2	B	501	FMN	C4A-N5	3.93	1.38	1.33
2	A	501	FMN	C4A-N5	3.37	1.38	1.33
2	A	501	FMN	C4-N3	3.11	1.38	1.33
2	C	501	FMN	C5A-N5	3.01	1.40	1.35
2	D	501	FMN	C5'-C4'	2.96	1.56	1.51
2	B	501	FMN	C9A-N10	2.95	1.42	1.38
2	B	501	FMN	C4-N3	2.84	1.38	1.33
2	C	501	FMN	C4-N3	2.65	1.37	1.33
2	D	501	FMN	C4-N3	2.54	1.37	1.33
2	B	501	FMN	C5A-N5	2.40	1.39	1.35
2	A	501	FMN	C5A-N5	2.29	1.39	1.35
2	D	501	FMN	C5A-N5	2.20	1.39	1.35
2	A	501	FMN	C5'-C4'	2.15	1.54	1.51

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	FMN	C1'-N10-C9A	12.54	128.16	118.29
2	B	501	FMN	C4-N3-C2	6.16	120.35	115.14
2	C	501	FMN	C4-N3-C2	6.16	120.34	115.14
2	D	501	FMN	C4-N3-C2	6.15	120.33	115.14
2	D	501	FMN	P-O5'-C5'	5.90	134.53	118.30
2	B	501	FMN	C1'-N10-C10	-5.42	113.55	118.41
2	D	501	FMN	C5'-C4'-C3'	5.22	122.30	112.20
2	A	501	FMN	C4-N3-C2	5.21	119.54	115.14
2	A	501	FMN	C1'-N10-C9A	5.11	122.32	118.29
5	D	506	IND	C2-C3-C9	5.09	120.70	105.05
5	C	505	IND	C2-C3-C9	5.07	120.64	105.05
5	D	506	IND	C3-C9-C8	-4.76	102.13	106.20
5	C	505	IND	C3-C9-C8	-4.66	102.22	106.20
2	C	501	FMN	C5A-C9A-N10	4.01	120.62	117.72
2	A	501	FMN	C5A-C9A-N10	3.97	120.59	117.72

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	501	FMN	C1'-N10-C9A	3.72	121.22	118.29
2	D	501	FMN	C5A-C9A-N10	3.71	120.41	117.72
2	B	501	FMN	C4A-N5-C5A	3.66	120.43	116.77
2	D	501	FMN	O5'-P-O1P	3.59	116.55	106.47
2	A	501	FMN	C4A-N5-C5A	3.42	120.19	116.77
2	B	501	FMN	O2'-C2'-C1'	-3.16	101.98	109.59
2	D	501	FMN	C1'-N10-C9A	3.10	120.73	118.29
2	A	501	FMN	C9A-N10-C10	-3.05	117.91	121.91
2	B	501	FMN	C9A-N10-C10	-2.88	118.14	121.91
2	D	501	FMN	C4A-N5-C5A	2.76	119.53	116.77
2	D	501	FMN	C9A-N10-C10	-2.72	118.34	121.91
2	A	501	FMN	C4A-C4-N3	-2.72	119.71	123.43
2	B	501	FMN	C5A-C9A-N10	2.70	119.67	117.72
2	D	501	FMN	C4A-C4-N3	-2.67	119.78	123.43
5	C	505	IND	C7-C8-N1	2.66	138.16	130.80
5	D	506	IND	C7-C8-N1	2.53	137.81	130.80
2	C	501	FMN	C4A-N5-C5A	2.51	119.28	116.77
2	D	501	FMN	C1'-N10-C10	2.45	120.60	118.41
2	B	501	FMN	C10-C4A-N5	-2.40	119.60	121.26
2	B	501	FMN	C6-C5A-C9A	2.34	122.11	119.05
2	B	501	FMN	C4A-C4-N3	-2.26	120.33	123.43
2	D	501	FMN	O3'-C3'-C4'	2.24	114.21	108.81
2	C	501	FMN	C9A-N10-C10	-2.23	118.98	121.91
2	A	501	FMN	C10-C4A-N5	-2.16	119.76	121.26
2	D	501	FMN	O2P-P-O1P	-2.14	102.31	110.68
2	C	501	FMN	P-O5'-C5'	2.08	124.01	118.30
2	C	501	FMN	C4A-C4-N3	-2.04	120.64	123.43
2	A	501	FMN	P-O5'-C5'	2.02	123.87	118.30

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	501	FMN	C2'-C1'-N10-C9A
2	B	501	FMN	C2'-C1'-N10-C10
2	B	501	FMN	N10-C1'-C2'-O2'
2	B	501	FMN	N10-C1'-C2'-C3'
2	D	501	FMN	C2'-C3'-C4'-O4'
2	D	501	FMN	O3'-C3'-C4'-O4'
2	D	501	FMN	O3'-C3'-C4'-C5'
2	D	501	FMN	C3'-C4'-C5'-O5'
2	D	501	FMN	O4'-C4'-C5'-O5'

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Mol	Chain	Res	Type	Atoms
2	A	501	FMN	C1'-C2'-C3'-C4'
2	D	501	FMN	C2'-C3'-C4'-C5'
2	A	501	FMN	O2'-C2'-C3'-O3'
2	C	501	FMN	C2'-C3'-C4'-O4'
2	A	501	FMN	O2'-C2'-C3'-C4'
2	C	501	FMN	C3'-C4'-C5'-O5'
2	A	501	FMN	C1'-C2'-C3'-O3'
2	C	501	FMN	O3'-C3'-C4'-O4'
2	C	501	FMN	C2'-C3'-C4'-C5'
2	D	501	FMN	C4'-C5'-O5'-P
2	D	501	FMN	C5'-O5'-P-O2P
2	C	501	FMN	O4'-C4'-C5'-O5'
2	C	501	FMN	O3'-C3'-C4'-C5'

There are no ring outliers.
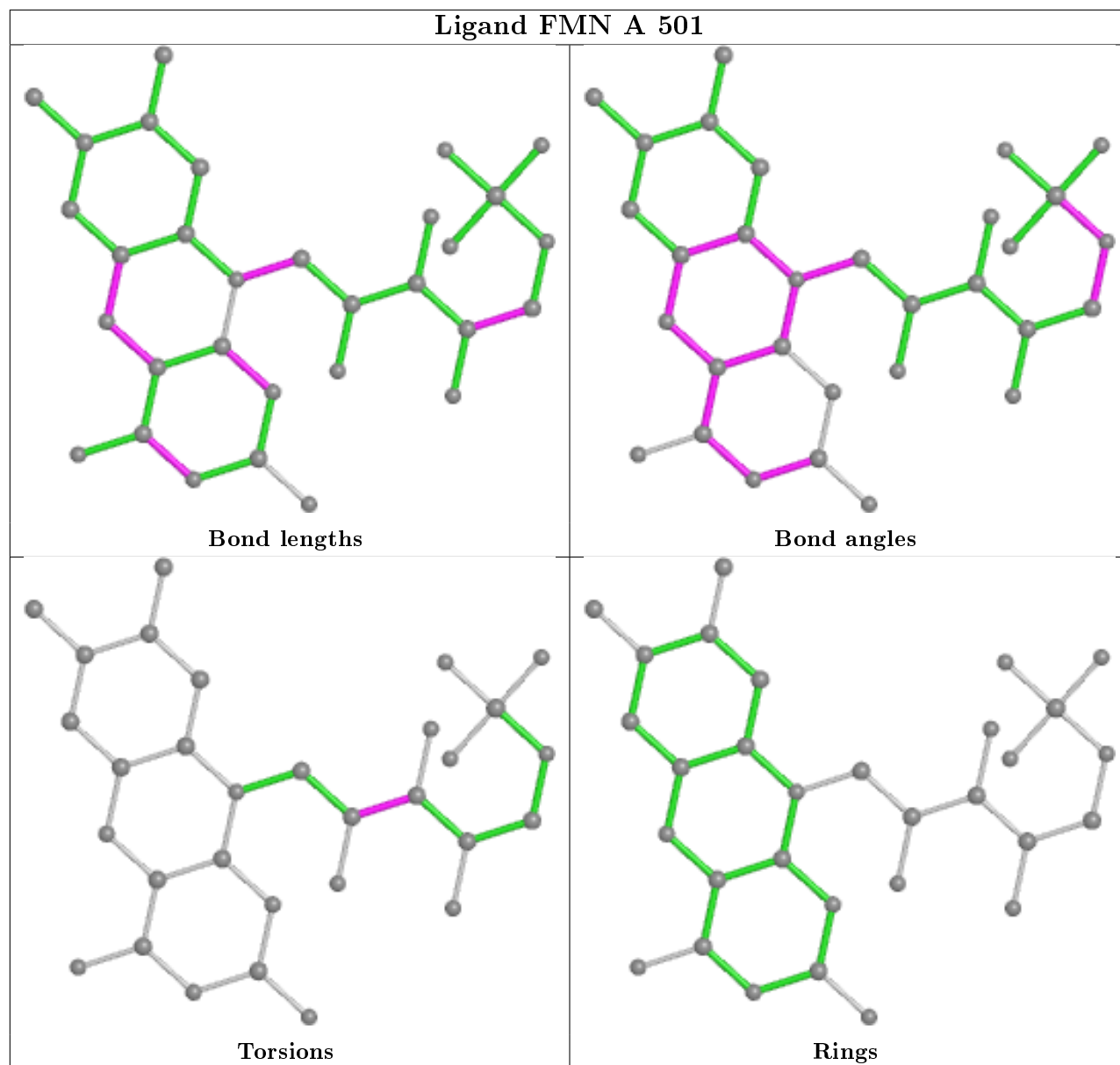
5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	FMN	1	0
2	D	501	FMN	1	0
3	A	503	SO4	1	0
3	D	503	SO4	1	0
5	D	506	IND	1	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 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 501

Bond lengths

Bond angles

Torsions

Rings

Ligand FMN D 501

Bond lengths

Bond angles

Torsions

Rings

Ligand FMN C 501

Bond lengths

Bond angles

Torsions

Rings

Ligand FMN A 501

Bond lengths

Bond angles

Torsions

Rings
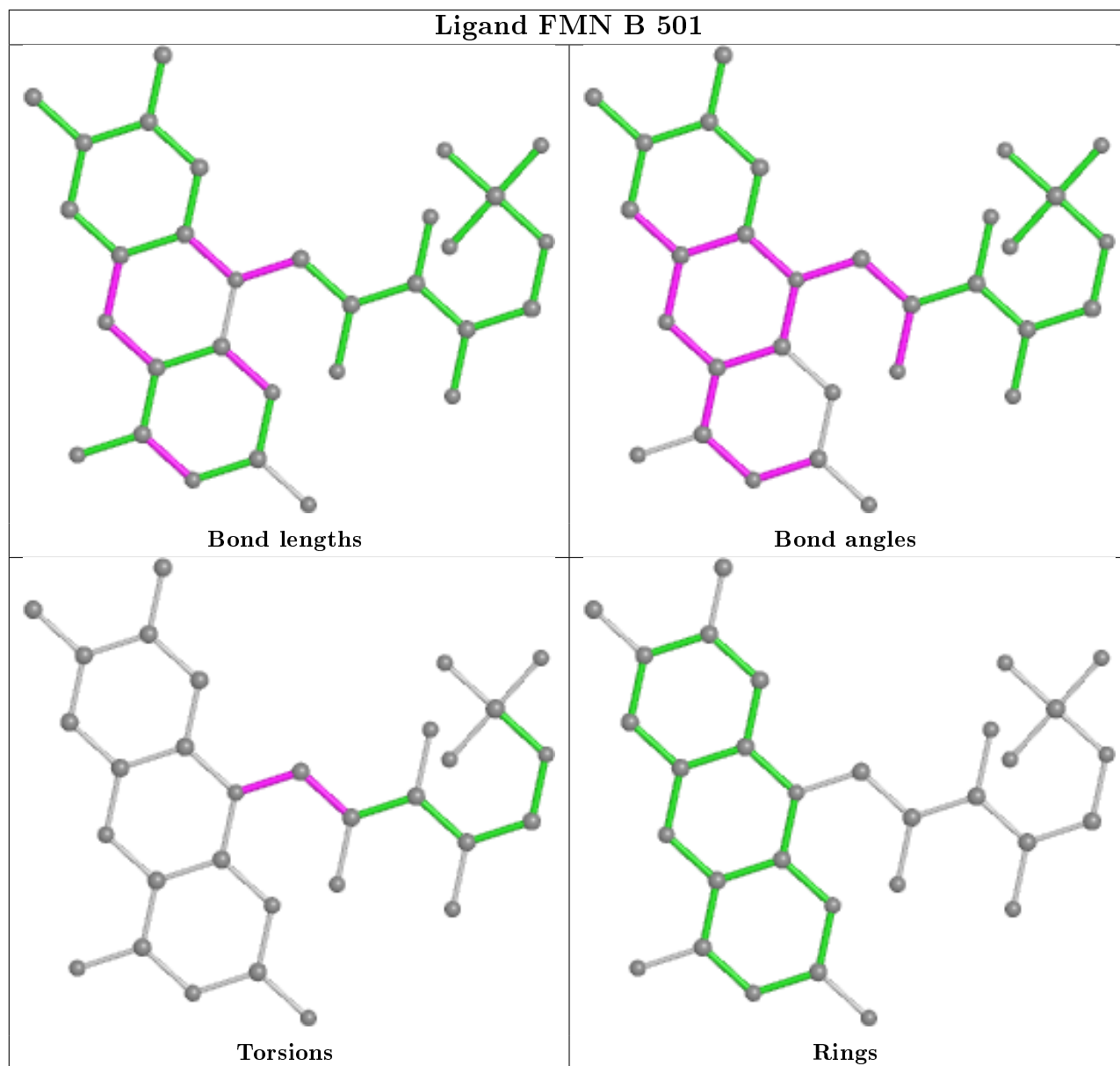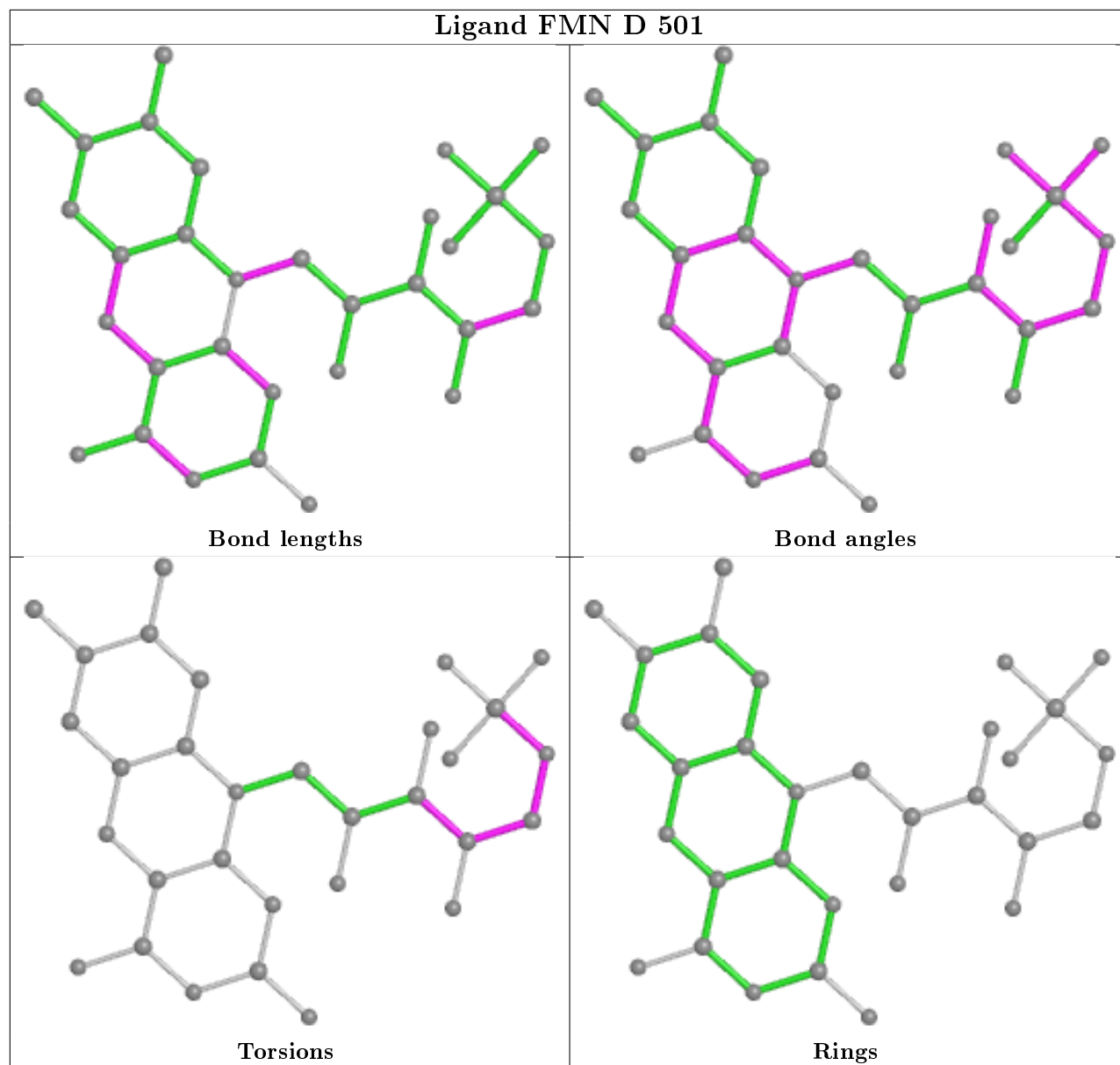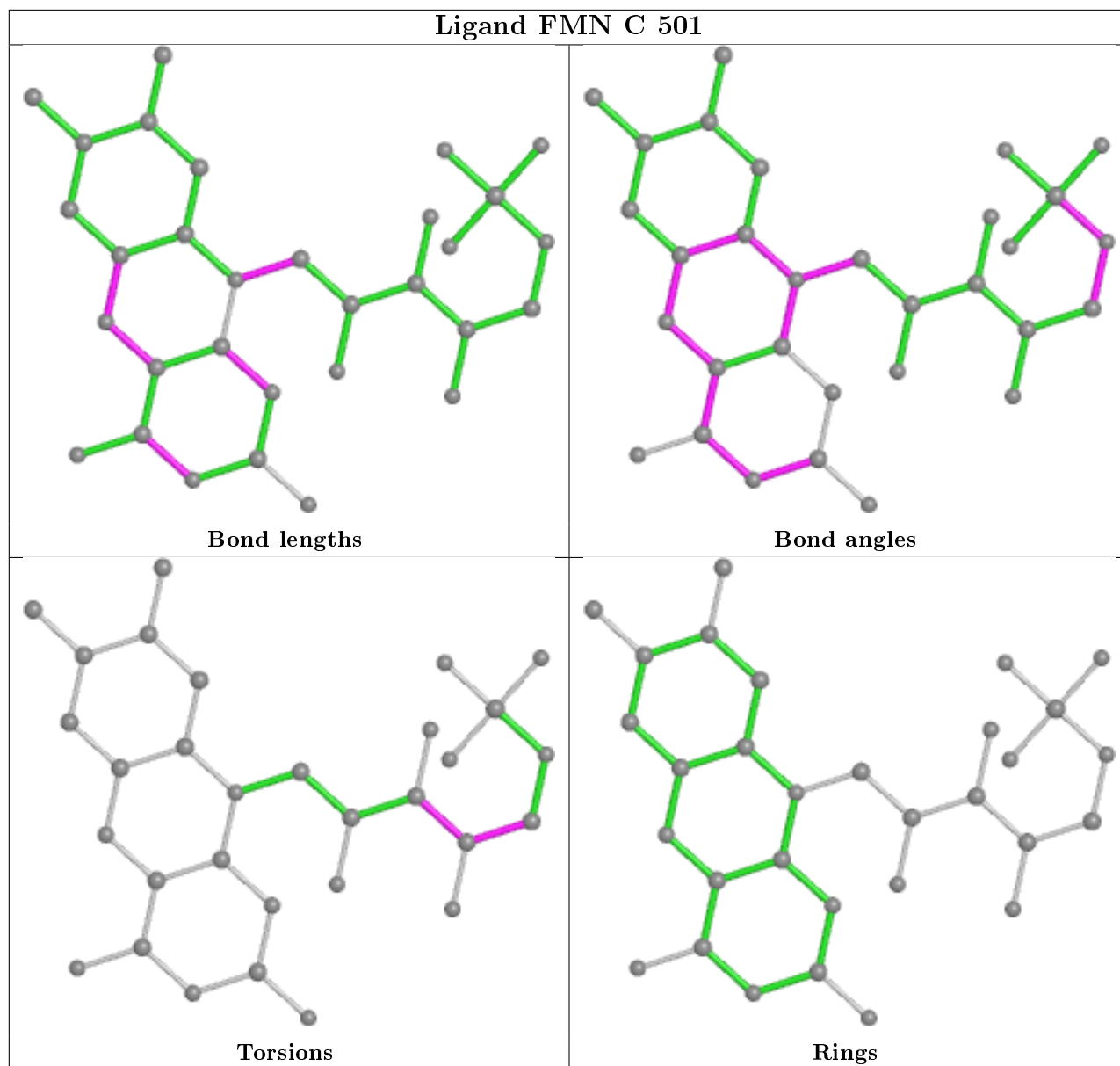
5.7 Other polymers (i)

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(Å^2)	Q<0.9
1	A	400/414 (96%)	0.29	25 (6%)	20	22	13, 22, 49, 78	0
1	B	394/414 (95%)	0.16	20 (5%)	28	31	13, 20, 44, 72	0
1	C	401/414 (96%)	0.21	13 (3%)	47	50	12, 22, 41, 79	0
1	D	400/414 (96%)	0.03	9 (2%)	60	63	14, 20, 38, 60	0
All	All	1595/1656 (96%)	0.17	67 (4%)	36	39	12, 21, 43, 79	0

All (67) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	133	ALA	7.1
1	A	134	HIS	6.9
1	A	15	HIS	6.8
1	C	15	HIS	5.7
1	A	135	VAL	5.5
1	A	131	ASN	5.4
1	A	132	ASN	5.4
1	C	14	ASP	5.4
1	B	282	THR	5.4
1	A	282	THR	5.2
1	A	283	LEU	5.1
1	A	146	PRO	4.9
1	D	136	LEU	4.8
1	A	14	ASP	4.7
1	B	15	HIS	4.3
1	A	178	PRO	4.2
1	B	279	ARG	4.2
1	B	285	GLY	4.1
1	B	283	LEU	4.1
1	B	284	ALA	3.8
1	D	137	ASP	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	286	VAL	3.6
1	A	176	ARG	3.6
1	D	146	PRO	3.5
1	B	137	ASP	3.5
1	A	138	TRP	3.4
1	B	287	GLU	3.2
1	A	145	LEU	3.1
1	A	16	ARG	3.1
1	A	182	GLY	3.0
1	C	16	ARG	3.0
1	A	286	VAL	2.9
1	D	16	ARG	2.9
1	D	287	GLU	2.8
1	C	176	ARG	2.8
1	C	62	GLU	2.7
1	A	285	GLY	2.7
1	B	288	LYS	2.7
1	B	138	TRP	2.7
1	D	134	HIS	2.5
1	B	278	SER	2.5
1	B	139	ARG	2.5
1	B	129	SER	2.4
1	C	137	ASP	2.4
1	A	136	LEU	2.3
1	B	281	PHE	2.2
1	C	146	PRO	2.2
1	B	347	VAL	2.2
1	A	277	HIS	2.2
1	D	133	ALA	2.2
1	C	37	ARG	2.2
1	D	138	TRP	2.2
1	C	302	PHE	2.1
1	B	280	PRO	2.1
1	A	279	ARG	2.1
1	B	176	ARG	2.1
1	A	284	ALA	2.1
1	B	16	ARG	2.1
1	D	286	VAL	2.1
1	C	133	ALA	2.1
1	B	353	ILE	2.1
1	C	63	ARG	2.0
1	A	218	SER	2.0

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Mol	Chain	Res	Type	RSRZ
1	C	139	ARG	2.0
1	A	409	THR	2.0
1	C	132	ASN	2.0
1	A	406	ARG	2.0

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

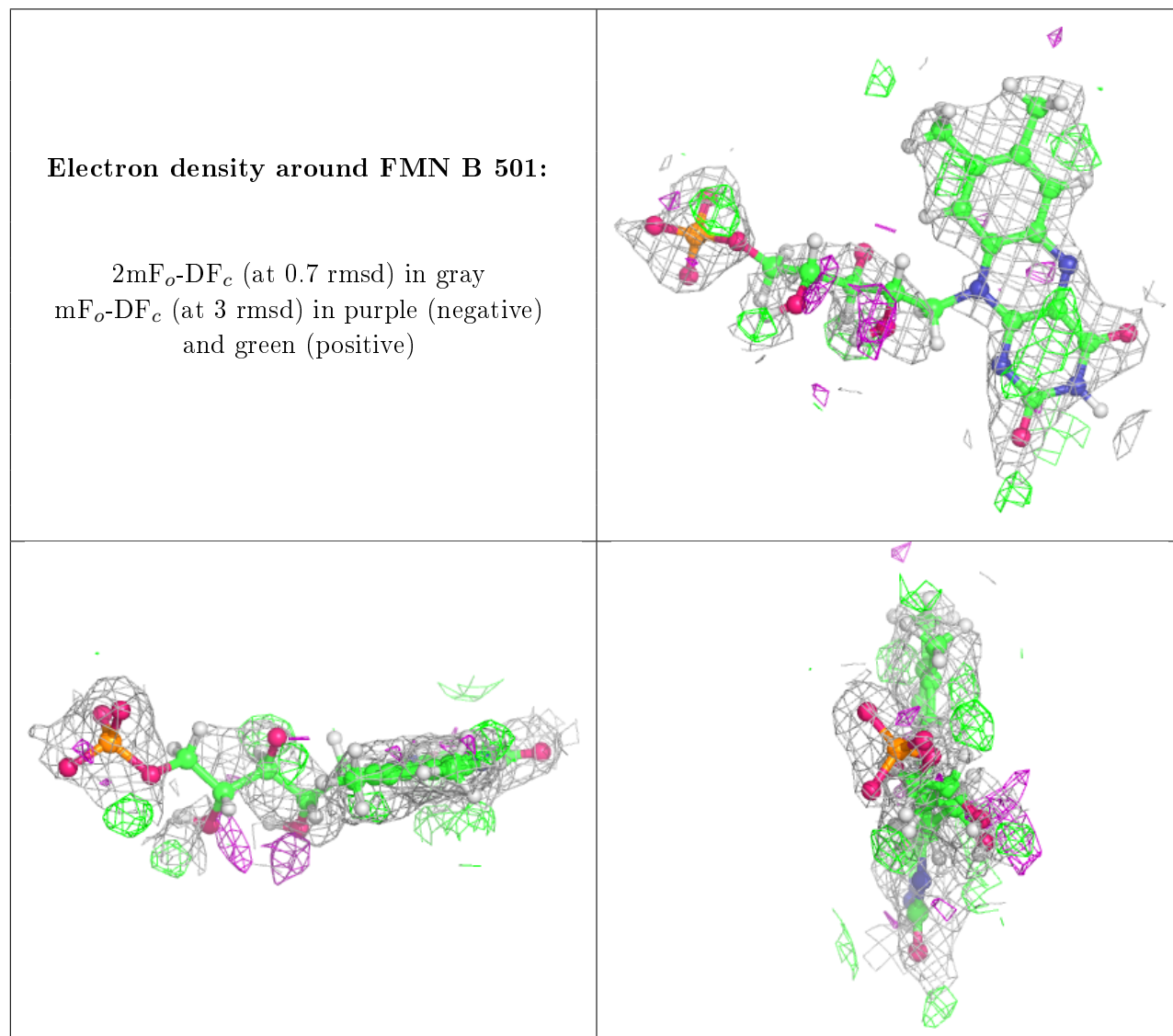There are no non-standard protein/DNA/RNA residues in this entry.

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
5	IND	C	505	9/9	0.61	0.25	41,49,59,60	0
3	SO4	D	502	5/5	0.73	0.23	40,41,60,65	5
2	FMN	B	501	31/31	0.75	0.24	29,52,78,86	0
5	IND	D	506	9/9	0.76	0.22	31,46,63,63	0
2	FMN	A	501	31/31	0.78	0.23	27,48,71,73	0
2	FMN	D	501	31/31	0.86	0.19	24,36,60,80	0
3	SO4	B	503	5/5	0.90	0.25	41,44,56,62	5
3	SO4	C	503	5/5	0.92	0.27	43,44,62,64	0
2	FMN	C	501	31/31	0.92	0.12	21,32,40,42	0
3	SO4	A	503	5/5	0.93	0.12	39,50,62,66	0
3	SO4	D	505	5/5	0.94	0.23	47,54,68,76	0
3	SO4	D	503	5/5	0.94	0.14	36,38,42,45	5
4	GOL	B	504	6/6	0.94	0.09	20,28,33,34	0
3	SO4	C	502	5/5	0.95	0.14	29,33,54,67	0
3	SO4	B	502	5/5	0.95	0.22	40,42,46,47	0
3	SO4	A	504	5/5	0.97	0.09	37,38,43,44	0
3	SO4	A	502	5/5	0.97	0.11	39,43,45,49	0
3	SO4	D	504	5/5	0.97	0.11	36,36,41,44	0
3	SO4	C	504	5/5	0.98	0.12	38,39,41,65	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.
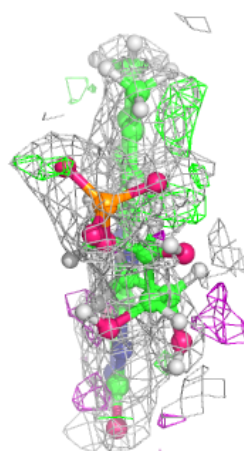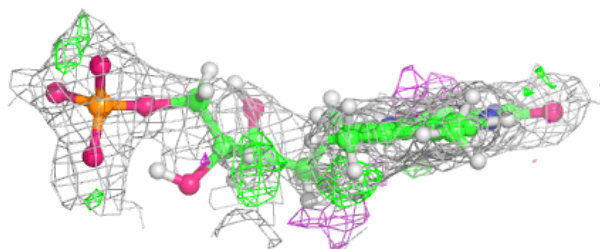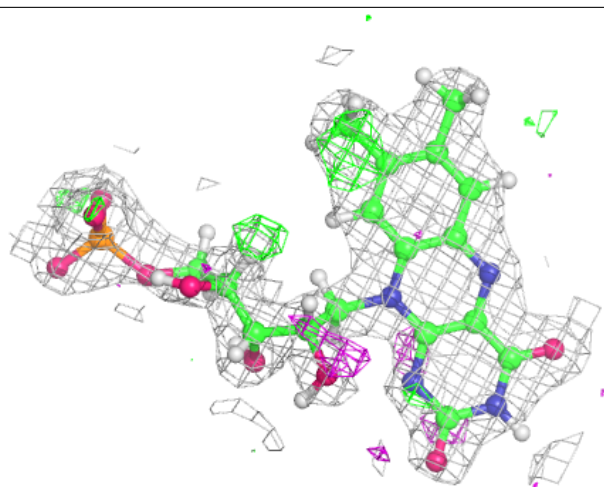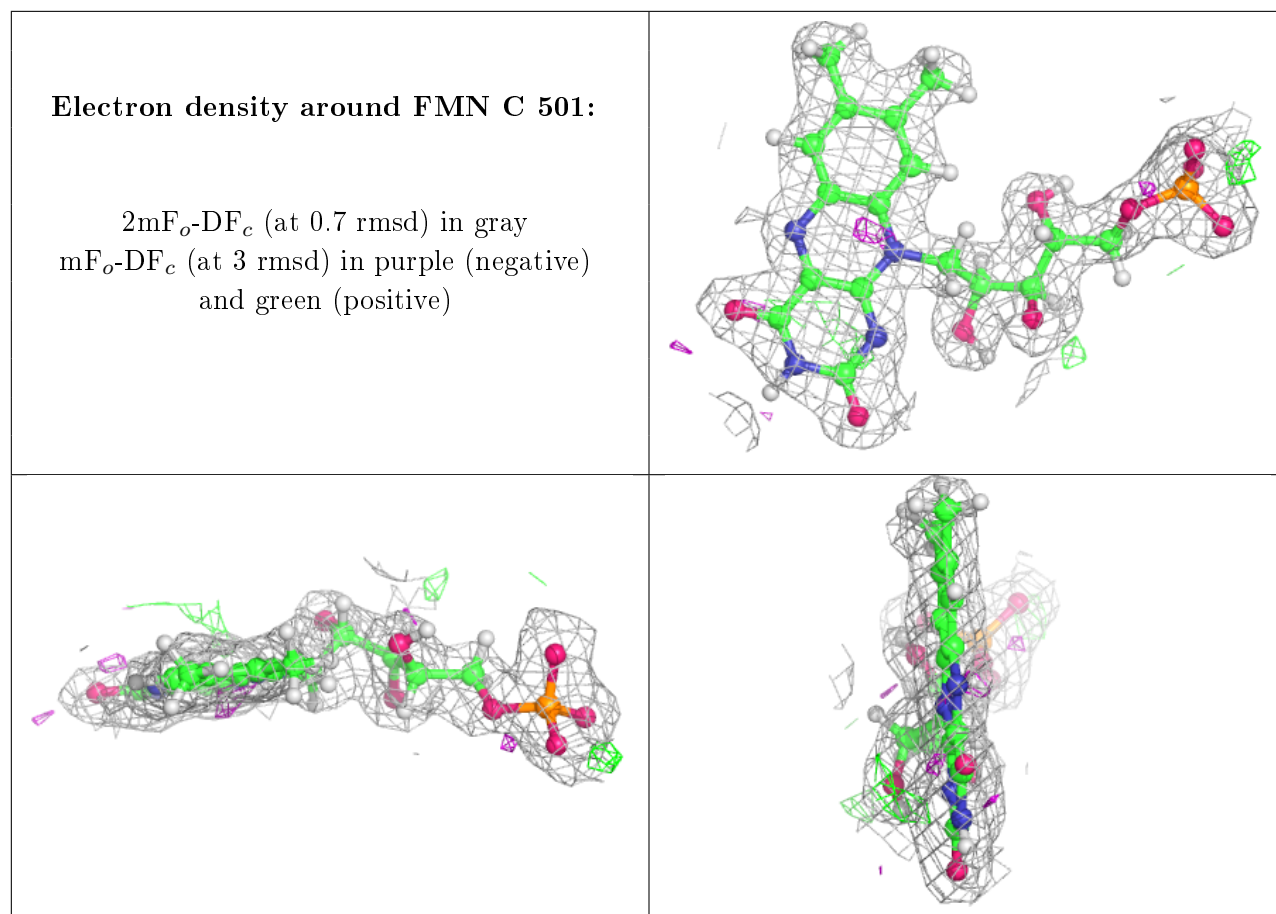
Electron density around FMN B 501:

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

Electron density around FMN A 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)
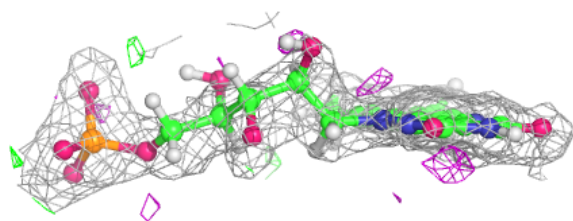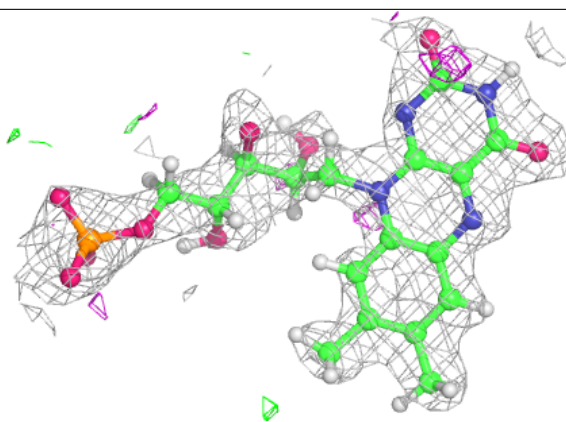
Electron density around FMN D 501:

$2mF_o$-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 501:

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

6.5 Other polymers (i)

There are no such residues in this entry.