



# Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 08:07 pm BST

PDB ID : 1NEC  
Title : NITROREDUCTASE FROM ENTEROBACTER CLOACAE  
Authors : Hecht, H.J.; Bryant, C.; Erdmann, H.; Pelletier, H.; Sawaya, R.  
Deposited on : 1999-03-30  
Resolution : 1.95 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

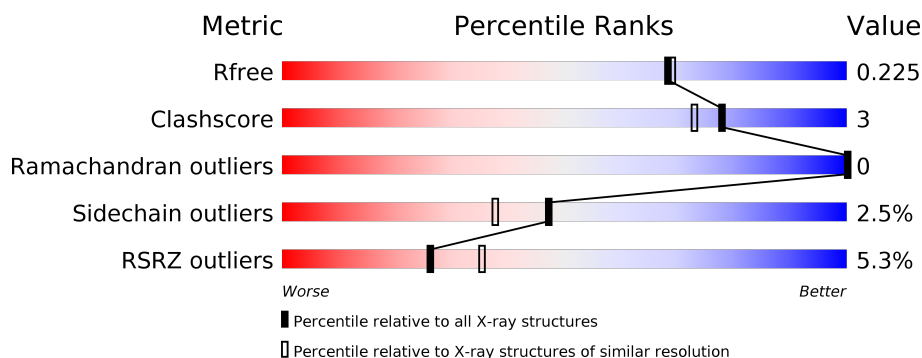
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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

Mol	Chain	Length	Quality of chain
1	A	216	<div> <div>12%</div> <div> <div></div> <div>82%</div> <div>16%</div> <div>.</div> </div> </div>
1	B	216	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>19%</div> <div>.</div> </div> </div>
1	C	216	<div> <div>8%</div> <div> <div></div> <div>81%</div> <div>16%</div> <div>.</div> </div> </div>
1	D	216	<div> <div></div> <div> <div></div> <div>81%</div> <div>17%</div> <div>.</div> </div> </div>

## 2 Entry composition [i](#)

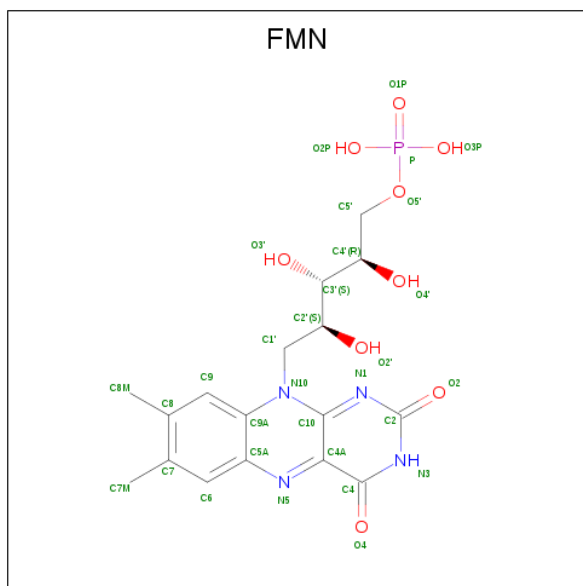
There are 3 unique types of molecules in this entry. The entry contains 7152 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROTEIN (NITROREDUCTASE).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	216	Total	C	N	O	S	0	0	0
			1677	1060	286	324	7			
1	B	216	Total	C	N	O	S	0	0	0
			1677	1060	286	324	7			
1	C	216	Total	C	N	O	S	0	0	0
			1677	1060	286	324	7			
1	D	216	Total	C	N	O	S	0	0	0
			1677	1060	286	324	7			

- 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
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		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
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		

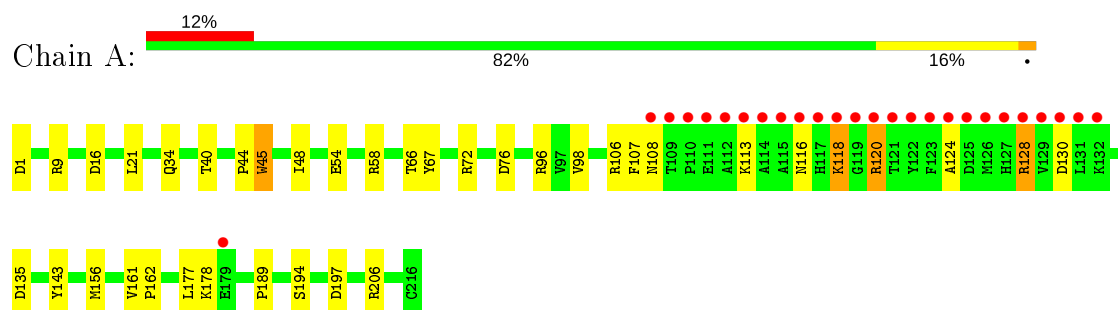
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	72	Total	O	0	0
			72	72		
3	B	79	Total	O	0	0
			79	79		
3	C	85	Total	O	0	0
			85	85		
3	D	84	Total	O	0	0
			84	84		

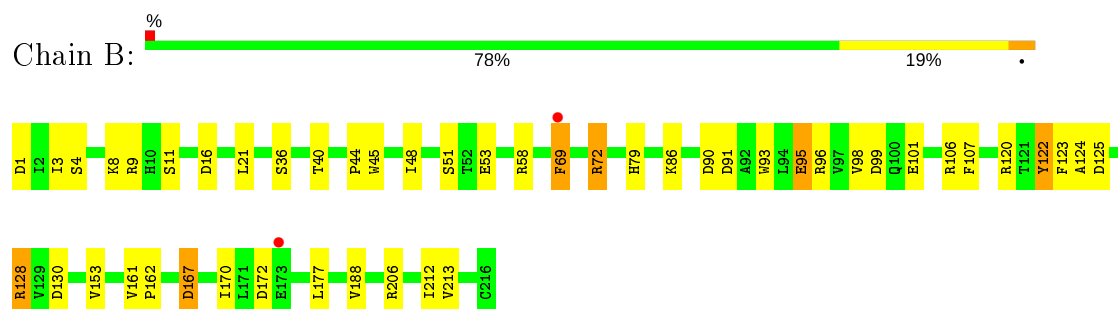
### 3 Residue-property plots [i](#)

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.

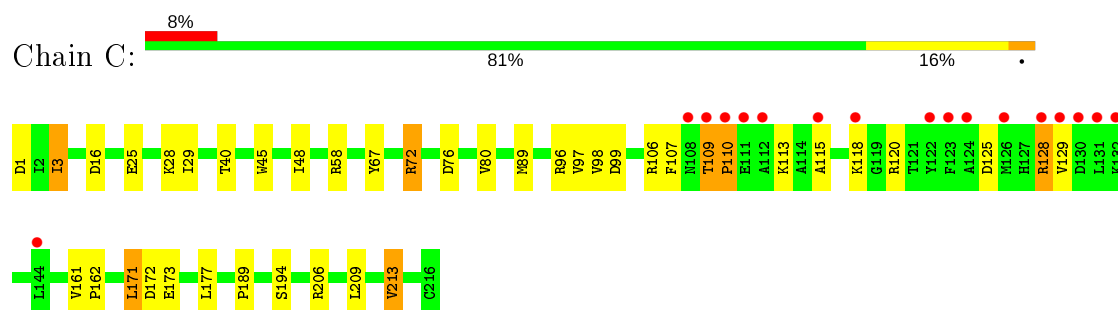
- Molecule 1: PROTEIN (NITROREDUCTASE)



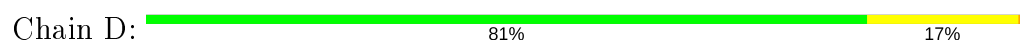
- Molecule 1: PROTEIN (NITROREDUCTASE)

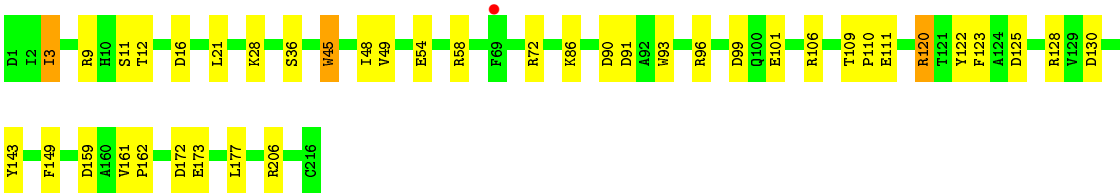


- Molecule 1: PROTEIN (NITROREDUCTASE)



- Molecule 1: PROTEIN (NITROREDUCTASE)





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	44.00 Å 92.50 Å 102.50 Å 90.00° 93.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.95 29.52 – 1.95	Depositor EDS
% Data completeness (in resolution range)	95.0 (20.00-1.95) 79.3 (29.52-1.95)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	0.00 (at 1.95 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.180 , 0.240 0.172 , 0.225	Depositor DCC
$R_{free}$ test set	2644 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	19.0	Xtriage
Anisotropy	0.350	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 50.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.54$ , $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	7152	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	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 47.64 % 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 9.6460e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

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

## 5 Model quality

### 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	0.61	0/1711	1.72	31/2314 (1.3%)
1	B	0.64	0/1711	1.78	39/2314 (1.7%)
1	C	0.64	0/1711	1.76	39/2314 (1.7%)
1	D	0.66	0/1711	1.66	32/2314 (1.4%)
All	All	0.64	0/6844	1.73	141/9256 (1.5%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
All	All	0	3

There are no bond length outliers.

All (141) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	96	ARG	NE-CZ-NH1	18.27	129.44	120.30
1	C	72	ARG	NE-CZ-NH1	14.75	127.67	120.30
1	A	120	ARG	NE-CZ-NH1	14.04	127.32	120.30
1	C	72	ARG	NE-CZ-NH2	-12.77	113.92	120.30
1	B	72	ARG	NE-CZ-NH1	12.45	126.52	120.30
1	B	128	ARG	NE-CZ-NH1	12.31	126.45	120.30
1	C	128	ARG	NE-CZ-NH1	11.81	126.21	120.30
1	B	96	ARG	CD-NE-CZ	10.92	138.89	123.60
1	B	96	ARG	NE-CZ-NH2	-10.82	114.89	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	128	ARG	NE-CZ-NH2	-10.75	114.92	120.30
1	A	128	ARG	NE-CZ-NH1	10.48	125.54	120.30
1	A	1	ASP	CB-CG-OD1	10.06	127.35	118.30
1	A	96	ARG	NE-CZ-NH2	-9.77	115.42	120.30
1	D	90	ASP	CB-CG-OD2	9.71	127.04	118.30
1	B	72	ARG	NE-CZ-NH2	-9.61	115.49	120.30
1	C	72	ARG	CD-NE-CZ	9.42	136.79	123.60
1	D	120	ARG	NE-CZ-NH1	9.35	124.97	120.30
1	A	58	ARG	CD-NE-CZ	9.23	136.52	123.60
1	C	120	ARG	CD-NE-CZ	9.12	136.37	123.60
1	A	206	ARG	NE-CZ-NH2	-9.07	115.76	120.30
1	A	96	ARG	NE-CZ-NH1	8.88	124.74	120.30
1	C	171	LEU	CA-CB-CG	8.74	135.40	115.30
1	C	96	ARG	NE-CZ-NH1	8.58	124.59	120.30
1	A	120	ARG	NE-CZ-NH2	-8.57	116.02	120.30
1	A	156	MET	CA-CB-CG	8.55	127.83	113.30
1	C	172	ASP	CB-CG-OD1	8.52	125.97	118.30
1	D	99	ASP	CB-CG-OD1	8.26	125.73	118.30
1	B	99	ASP	CB-CG-OD1	8.23	125.70	118.30
1	A	106	ARG	CD-NE-CZ	8.19	135.07	123.60
1	A	76	ASP	CB-CG-OD2	-8.15	110.96	118.30
1	B	120	ARG	NE-CZ-NH2	-8.14	116.23	120.30
1	C	206	ARG	CD-NE-CZ	8.13	134.99	123.60
1	B	58	ARG	CD-NE-CZ	8.09	134.93	123.60
1	C	106	ARG	CD-NE-CZ	8.03	134.84	123.60
1	C	120	ARG	NE-CZ-NH1	8.02	124.31	120.30
1	D	106	ARG	CD-NE-CZ	7.94	134.72	123.60
1	D	72	ARG	CD-NE-CZ	7.86	134.60	123.60
1	C	125	ASP	CB-CG-OD1	7.85	125.36	118.30
1	A	108	ASN	C-N-CA	7.82	141.25	121.70
1	A	72	ARG	CD-NE-CZ	7.80	134.51	123.60
1	C	1	ASP	CB-CG-OD1	7.79	125.31	118.30
1	B	72	ARG	CD-NE-CZ	7.64	134.30	123.60
1	A	128	ARG	CD-NE-CZ	7.63	134.28	123.60
1	C	96	ARG	CD-NE-CZ	7.61	134.26	123.60
1	B	125	ASP	CB-CG-OD1	7.61	125.15	118.30
1	A	120	ARG	CD-NE-CZ	7.60	134.24	123.60
1	D	128	ARG	NE-CZ-NH1	7.54	124.07	120.30
1	A	16	ASP	CB-CG-OD2	-7.45	111.59	118.30
1	B	69	PHE	O-C-N	-7.29	111.04	122.70
1	C	106	ARG	NE-CZ-NH1	7.24	123.92	120.30
1	B	1	ASP	CB-CG-OD1	7.22	124.80	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	122	TYR	CB-CG-CD1	-7.22	116.67	121.00
1	C	58	ARG	CD-NE-CZ	7.20	133.68	123.60
1	D	206	ARG	NE-CZ-NH2	-7.13	116.74	120.30
1	B	1	ASP	CB-CG-OD2	-7.05	111.95	118.30
1	B	206	ARG	CD-NE-CZ	6.87	133.22	123.60
1	A	40	THR	O-C-N	-6.81	111.80	122.70
1	C	106	ARG	NE-CZ-NH2	-6.81	116.90	120.30
1	D	58	ARG	NE-CZ-NH1	6.80	123.70	120.30
1	B	9	ARG	CD-NE-CZ	6.75	133.06	123.60
1	D	101	GLU	OE1-CD-OE2	-6.68	115.28	123.30
1	A	76	ASP	CB-CG-OD1	6.67	124.30	118.30
1	D	96	ARG	NE-CZ-NH2	-6.60	117.00	120.30
1	A	96	ARG	CD-NE-CZ	6.59	132.82	123.60
1	C	128	ARG	CG-CD-NE	6.58	125.62	111.80
1	B	130	ASP	CB-CG-OD2	-6.57	112.39	118.30
1	D	123	PHE	CB-CG-CD1	6.55	125.39	120.80
1	C	120	ARG	NE-CZ-NH2	-6.52	117.04	120.30
1	B	90	ASP	CB-CG-OD1	6.52	124.17	118.30
1	C	173	GLU	CA-CB-CG	6.44	127.56	113.40
1	B	172	ASP	CB-CG-OD1	6.40	124.06	118.30
1	B	120	ARG	CD-NE-CZ	6.39	132.55	123.60
1	D	96	ARG	CD-NE-CZ	6.38	132.53	123.60
1	C	72	ARG	CA-CB-CG	6.34	127.35	113.40
1	D	3	ILE	CB-CG1-CD1	6.29	131.50	113.90
1	A	9	ARG	CD-NE-CZ	6.25	132.35	123.60
1	A	54	GLU	CB-CG-CD	6.22	131.00	114.20
1	A	21	LEU	CA-CB-CG	6.19	129.55	115.30
1	D	54	GLU	CA-CB-CG	6.17	126.98	113.40
1	D	16	ASP	CB-CG-OD1	6.14	123.82	118.30
1	C	110	PRO	C-N-CA	6.12	137.01	121.70
1	B	40	THR	O-C-N	-6.11	112.92	122.70
1	D	122	TYR	CB-CG-CD2	6.10	124.66	121.00
1	D	91	ASP	CB-CG-OD1	6.02	123.72	118.30
1	A	143	TYR	CB-CG-CD2	-6.02	117.39	121.00
1	A	9	ARG	NE-CZ-NH2	6.00	123.30	120.30
1	D	206	ARG	CD-NE-CZ	5.99	131.98	123.60
1	C	76	ASP	CB-CG-OD2	-5.97	112.92	118.30
1	B	188	VAL	CA-CB-CG2	5.95	119.82	110.90
1	B	153	VAL	CA-CB-CG2	5.90	119.74	110.90
1	D	36	SER	N-CA-CB	-5.88	101.68	110.50
1	D	45	TRP	CA-CB-CG	5.87	124.85	113.70
1	D	58	ARG	NE-CZ-NH2	-5.86	117.37	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	67	TYR	CB-CG-CD1	-5.83	117.50	121.00
1	A	156	MET	O-C-N	-5.82	113.30	123.20
1	B	122	TYR	CB-CG-CD1	-5.82	117.51	121.00
1	D	21	LEU	CA-CB-CG	5.80	128.65	115.30
1	B	128	ARG	NE-CZ-NH2	-5.79	117.40	120.30
1	B	91	ASP	CB-CG-OD1	5.76	123.48	118.30
1	C	40	THR	O-C-N	-5.74	113.51	122.70
1	C	80	VAL	CG1-CB-CG2	-5.71	101.76	110.90
1	D	9	ARG	CD-NE-CZ	5.71	131.60	123.60
1	D	172	ASP	CB-CG-OD1	5.71	123.44	118.30
1	C	25	GLU	OE1-CD-OE2	-5.70	116.46	123.30
1	B	3	ILE	CB-CG1-CD1	5.67	129.78	113.90
1	B	167	ASP	CB-CG-OD1	5.67	123.40	118.30
1	D	143	TYR	CB-CG-CD2	-5.65	117.61	121.00
1	B	53	GLU	CB-CA-C	5.64	121.68	110.40
1	A	45	TRP	CA-CB-CG	5.62	124.37	113.70
1	C	99	ASP	CB-CG-OD2	5.60	123.34	118.30
1	D	128	ARG	CD-NE-CZ	5.60	131.44	123.60
1	C	3	ILE	CB-CG1-CD1	5.57	129.50	113.90
1	C	99	ASP	CB-CG-OD1	5.57	123.31	118.30
1	D	149	PHE	CB-CG-CD2	-5.54	116.92	120.80
1	A	197	ASP	CB-CG-OD1	5.50	123.25	118.30
1	A	143	TYR	CB-CG-CD1	5.42	124.25	121.00
1	C	209	LEU	CB-CG-CD2	5.41	120.20	111.00
1	B	106	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	A	206	ARG	CD-NE-CZ	5.38	131.14	123.60
1	C	128	ARG	CD-NE-CZ	5.34	131.08	123.60
1	C	16	ASP	CB-CG-OD1	5.28	123.05	118.30
1	A	130	ASP	CA-CB-CG	5.26	124.97	113.40
1	C	99	ASP	OD1-CG-OD2	-5.25	113.33	123.30
1	C	67	TYR	CA-CB-CG	5.24	123.35	113.40
1	D	125	ASP	CB-CG-OD1	5.23	123.01	118.30
1	C	76	ASP	CB-CG-OD1	5.21	122.99	118.30
1	D	111	GLU	OE1-CD-OE2	-5.21	117.05	123.30
1	B	107	PHE	CB-CG-CD2	-5.19	117.17	120.80
1	B	101	GLU	OE1-CD-OE2	-5.18	117.08	123.30
1	C	213	VAL	CG1-CB-CG2	-5.17	102.64	110.90
1	A	72	ARG	NE-CZ-NH1	5.15	122.87	120.30
1	B	36	SER	N-CA-CB	-5.14	102.79	110.50
1	B	213	VAL	CG1-CB-CG2	-5.14	102.68	110.90
1	D	130	ASP	CB-CG-OD1	5.14	122.92	118.30
1	C	129	VAL	CA-CB-CG1	5.12	118.58	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	16	ASP	CB-CG-OD1	5.12	122.91	118.30
1	C	97	VAL	CA-CB-CG1	5.11	118.56	110.90
1	B	120	ARG	NE-CZ-NH1	5.10	122.85	120.30
1	B	106	ARG	CD-NE-CZ	5.04	130.66	123.60
1	B	21	LEU	CA-CB-CG	5.02	126.85	115.30
1	B	95	GLU	OE1-CD-OE2	5.01	129.32	123.30

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	194	SER	Mainchain
1	B	69	PHE	Mainchain
1	C	194	SER	Mainchain

## 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	1677	0	1653	10	0
1	B	1677	0	1653	13	0
1	C	1677	0	1653	12	0
1	D	1677	0	1653	11	0
2	A	31	0	19	1	0
2	B	31	0	19	1	0
2	C	31	0	19	1	0
2	D	31	0	19	1	0
3	A	72	0	0	0	0
3	B	79	0	0	0	0
3	C	85	0	0	0	0
3	D	84	0	0	1	0
All	All	7152	0	6688	39	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.

All (39) 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:11:SER:HB3	1:B:161:VAL:HG23	1.71	0.72
1:A:48:ILE:HD11	1:A:177:LEU:HD21	1.80	0.62
1:C:115:ALA:HA	1:C:118:LYS:HE2	1.84	0.59
1:B:162:PRO:HG2	2:B:218:FMN:C9	2.35	0.55
1:C:162:PRO:HG2	2:C:219:FMN:C9	2.38	0.54
1:D:162:PRO:HG2	2:D:220:FMN:C9	2.39	0.53
1:A:162:PRO:HG2	2:A:217:FMN:C9	2.39	0.53
1:A:98:VAL:HG22	1:A:120:ARG:NH2	2.25	0.52
1:C:48:ILE:HD11	1:C:177:LEU:HD21	1.92	0.51
1:C:28:LYS:HB3	1:D:3:ILE:HD11	1.92	0.50
1:B:167:ASP:HB3	1:B:170:ILE:HD12	1.94	0.50
1:D:86:LYS:HE2	1:D:93:TRP:CG	2.47	0.50
1:D:48:ILE:HD11	1:D:177:LEU:HD21	1.96	0.47
1:C:89:MET:O	1:C:128:ARG:HD2	2.15	0.47
1:C:109:THR:HG22	1:C:110:PRO:HD2	1.97	0.47
1:A:34:GLN:HB2	1:B:212:ILE:HD11	1.96	0.47
1:B:124:ALA:O	1:B:128:ARG:HG3	2.16	0.46
1:C:98:VAL:HG13	1:C:113:LYS:HE3	1.97	0.45
1:C:3:ILE:HD11	1:D:28:LYS:HB3	1.98	0.45
1:A:124:ALA:O	1:A:128:ARG:HG2	2.15	0.45
1:C:115:ALA:HA	1:C:118:LYS:HG2	1.99	0.45
1:D:12:THR:OG1	1:D:159:ASP:HB3	2.17	0.45
1:B:48:ILE:HD11	1:B:177:LEU:HD21	1.98	0.45
1:D:11:SER:HB3	1:D:161:VAL:HG23	1.99	0.45
1:A:161:VAL:O	1:A:189:PRO:HD2	2.16	0.45
1:B:4:SER:O	1:B:8:LYS:HG2	2.17	0.45
1:B:51:SER:HB3	1:B:79:HIS:CD2	2.52	0.44
1:A:118:LYS:O	1:A:118:LYS:HE2	2.17	0.44
1:B:95:GLU:HA	1:B:98:VAL:HG12	1.99	0.43
1:A:66:THR:HG21	1:B:122:TYR:OH	2.18	0.43
1:D:109:THR:HB	1:D:110:PRO:HD2	2.01	0.43
1:A:98:VAL:HG12	1:A:113:LYS:HG2	2.01	0.43
1:A:67:TYR:HE1	1:B:123:PHE:HE1	1.66	0.42
1:C:161:VAL:O	1:C:189:PRO:HD2	2.18	0.42
1:C:28:LYS:HG2	1:D:3:ILE:HG13	2.00	0.42
1:C:213:VAL:HG13	1:D:49:VAL:HG23	2.01	0.41
1:D:120:ARG:HD2	3:D:231:HOH:O	2.19	0.41
1:B:167:ASP:CB	1:B:170:ILE:HD12	2.50	0.41
1:B:86:LYS:HE2	1:B:93:TRP:CG	2.55	0.40

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	214/216 (99%)	212 (99%)	2 (1%)	0	100	100
1	B	214/216 (99%)	212 (99%)	2 (1%)	0	100	100
1	C	214/216 (99%)	211 (99%)	3 (1%)	0	100	100
1	D	214/216 (99%)	212 (99%)	2 (1%)	0	100	100
All	All	856/864 (99%)	847 (99%)	9 (1%)	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	179/179 (100%)	172 (96%)	7 (4%)	32	19
1	B	179/179 (100%)	176 (98%)	3 (2%)	60	55
1	C	179/179 (100%)	173 (97%)	6 (3%)	37	25
1	D	179/179 (100%)	177 (99%)	2 (1%)	73	71
All	All	716/716 (100%)	698 (98%)	18 (2%)	47	38

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44	PRO
1	A	45	TRP

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Mol	Chain	Res	Type
1	A	107	PHE
1	A	116	ASN
1	A	118	LYS
1	A	135	ASP
1	A	178	LYS
1	B	44	PRO
1	B	45	TRP
1	B	72	ARG
1	C	29	ILE
1	C	45	TRP
1	C	72	ARG
1	C	107	PHE
1	C	109	THR
1	C	171	LEU
1	D	45	TRP
1	D	173	GLU

Some 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 carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

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	218	-	31,33,33	1.85	6 (19%)	40,50,50	2.65	9 (22%)
2	FMN	C	219	-	31,33,33	2.08	7 (22%)	40,50,50	2.49	9 (22%)
2	FMN	A	217	-	31,33,33	1.91	7 (22%)	40,50,50	2.70	13 (32%)
2	FMN	D	220	-	31,33,33	2.06	5 (16%)	40,50,50	2.19	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	B	218	-	-	4/18/18/18	0/3/3/3
2	FMN	C	219	-	-	2/18/18/18	0/3/3/3
2	FMN	A	217	-	-	2/18/18/18	0/3/3/3
2	FMN	D	220	-	-	1/18/18/18	0/3/3/3

All (25) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	220	FMN	C4A-C10	9.41	1.48	1.38
2	C	219	FMN	C4A-C10	8.17	1.47	1.38
2	A	217	FMN	C4A-C10	7.68	1.46	1.38
2	B	218	FMN	C4A-C10	7.49	1.46	1.38
2	C	219	FMN	C4A-N5	-4.03	1.27	1.33
2	A	217	FMN	C4'-C3'	-3.15	1.47	1.53
2	D	220	FMN	C4-C4A	3.15	1.46	1.41
2	D	220	FMN	C2-N3	2.86	1.43	1.38
2	A	217	FMN	C4-C4A	2.86	1.46	1.41
2	C	219	FMN	O3'-C3'	2.80	1.49	1.43
2	B	218	FMN	C4-C4A	2.77	1.46	1.41
2	B	218	FMN	C1'-N10	-2.75	1.45	1.48
2	C	219	FMN	C4'-C3'	-2.63	1.48	1.53
2	B	218	FMN	C4A-N5	-2.47	1.29	1.33
2	C	219	FMN	C4-C4A	2.47	1.45	1.41
2	B	218	FMN	C2-N3	2.41	1.42	1.38
2	C	219	FMN	C2-N1	-2.39	1.33	1.38
2	A	217	FMN	C4A-N5	-2.28	1.30	1.33
2	A	217	FMN	C6-C5A	2.26	1.45	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	220	FMN	C4A-N5	-2.23	1.30	1.33
2	A	217	FMN	O3'-C3'	2.21	1.48	1.43
2	A	217	FMN	C2-N3	2.14	1.42	1.38
2	D	220	FMN	C7M-C7	2.14	1.55	1.51
2	C	219	FMN	C6-C7	-2.11	1.32	1.37
2	B	218	FMN	O3'-C3'	2.05	1.47	1.43

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	217	FMN	C4-N3-C2	11.47	124.83	115.14
2	C	219	FMN	C4-N3-C2	10.93	124.37	115.14
2	B	218	FMN	C4-N3-C2	10.79	124.25	115.14
2	D	220	FMN	C4-N3-C2	9.48	123.15	115.14
2	B	218	FMN	C4A-C4-N3	-6.87	114.04	123.43
2	D	220	FMN	C4A-C4-N3	-6.53	114.50	123.43
2	C	219	FMN	C4A-C4-N3	-5.73	115.60	123.43
2	A	217	FMN	C4A-C4-N3	-5.49	115.92	123.43
2	B	218	FMN	C1'-N10-C9A	5.43	122.56	118.29
2	A	217	FMN	C1'-N10-C9A	5.38	122.52	118.29
2	D	220	FMN	C4'-C3'-C2'	4.66	123.06	113.36
2	B	218	FMN	C4'-C3'-C2'	4.34	122.39	113.36
2	C	219	FMN	C1'-N10-C9A	3.99	121.44	118.29
2	A	217	FMN	C4-C4A-C10	-3.82	117.42	119.95
2	C	219	FMN	C4'-C3'-C2'	3.75	121.16	113.36
2	A	217	FMN	C7M-C7-C8	-3.51	113.54	120.74
2	B	218	FMN	C10-C4A-N5	3.33	123.56	121.26
2	C	219	FMN	O3P-P-O5'	-2.71	99.53	106.73
2	B	218	FMN	C9A-N10-C10	-2.58	118.53	121.91
2	C	219	FMN	C4A-N5-C5A	2.46	119.23	116.77
2	D	220	FMN	C5A-C9A-N10	-2.46	115.94	117.72
2	B	218	FMN	C7M-C7-C8	-2.44	115.73	120.74
2	A	217	FMN	C6-C5A-C9A	-2.42	115.87	119.05
2	D	220	FMN	C1'-N10-C10	2.39	120.55	118.41
2	A	217	FMN	O4'-C4'-C5'	-2.37	104.60	109.92
2	A	217	FMN	C10-C4A-N5	2.36	122.89	121.26
2	A	217	FMN	O2'-C2'-C1'	2.36	115.28	109.59
2	B	218	FMN	O2P-P-O5'	2.34	112.95	106.73
2	C	219	FMN	C4-C4A-C10	-2.32	118.42	119.95
2	B	218	FMN	C5'-C4'-C3'	2.30	116.64	112.20
2	A	217	FMN	O2P-P-O5'	2.21	112.62	106.73
2	A	217	FMN	C4'-C3'-C2'	2.19	117.91	113.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	217	FMN	C9A-C5A-N5	2.10	125.65	122.36
2	C	219	FMN	C5A-C9A-N10	-2.07	116.22	117.72
2	A	217	FMN	O2'-C2'-C3'	-2.06	104.08	109.10
2	C	219	FMN	C7M-C7-C8	-2.04	116.56	120.74

There are no chirality outliers.

All (9) torsion outliers are listed below:

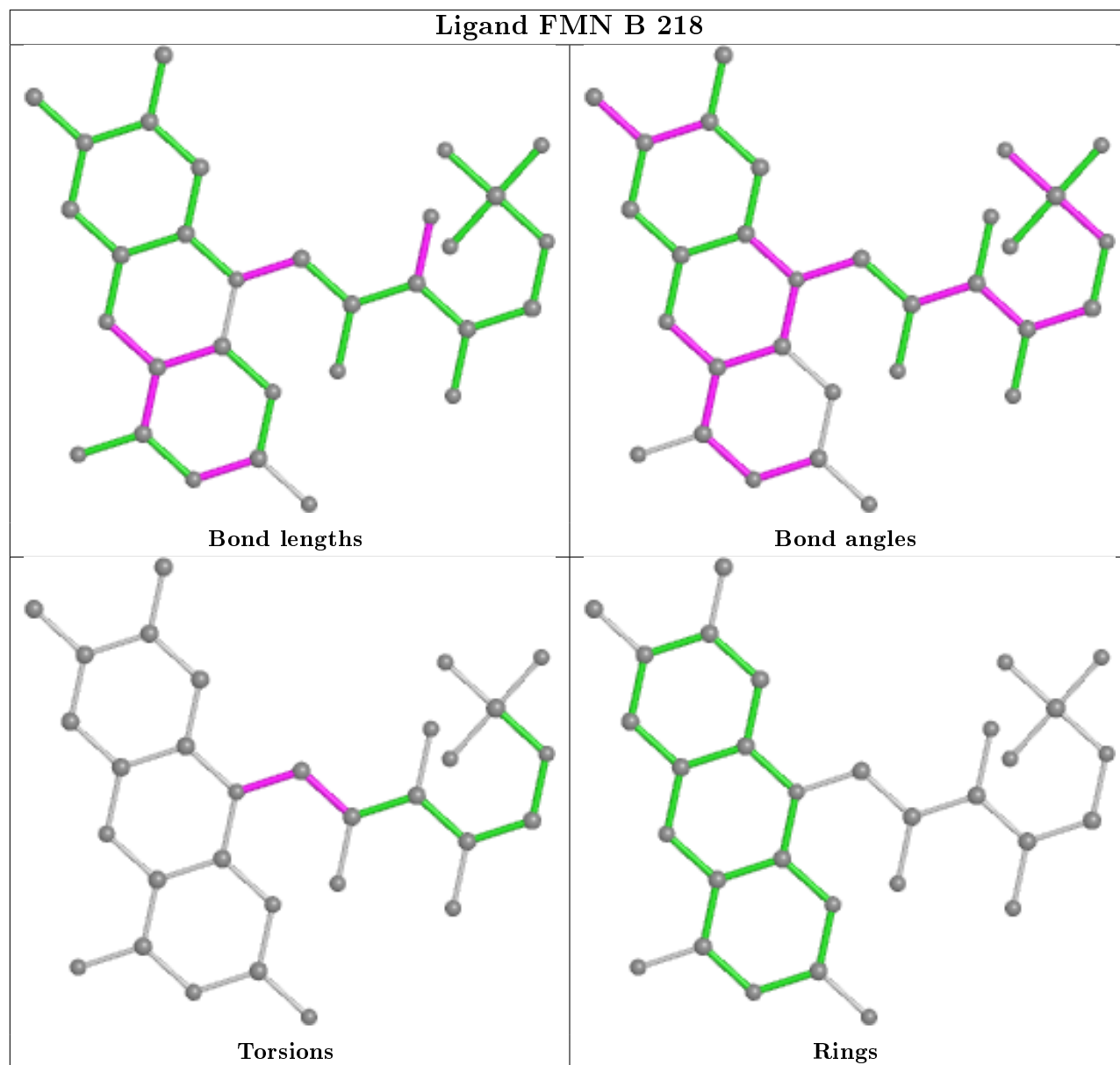
Mol	Chain	Res	Type	Atoms
2	B	218	FMN	C2'-C1'-N10-C9A
2	B	218	FMN	C2'-C1'-N10-C10
2	B	218	FMN	N10-C1'-C2'-O2'
2	B	218	FMN	N10-C1'-C2'-C3'
2	C	219	FMN	N10-C1'-C2'-O2'
2	A	217	FMN	C2'-C1'-N10-C9A
2	A	217	FMN	N10-C1'-C2'-O2'
2	D	220	FMN	N10-C1'-C2'-O2'
2	C	219	FMN	C4'-C5'-O5'-P

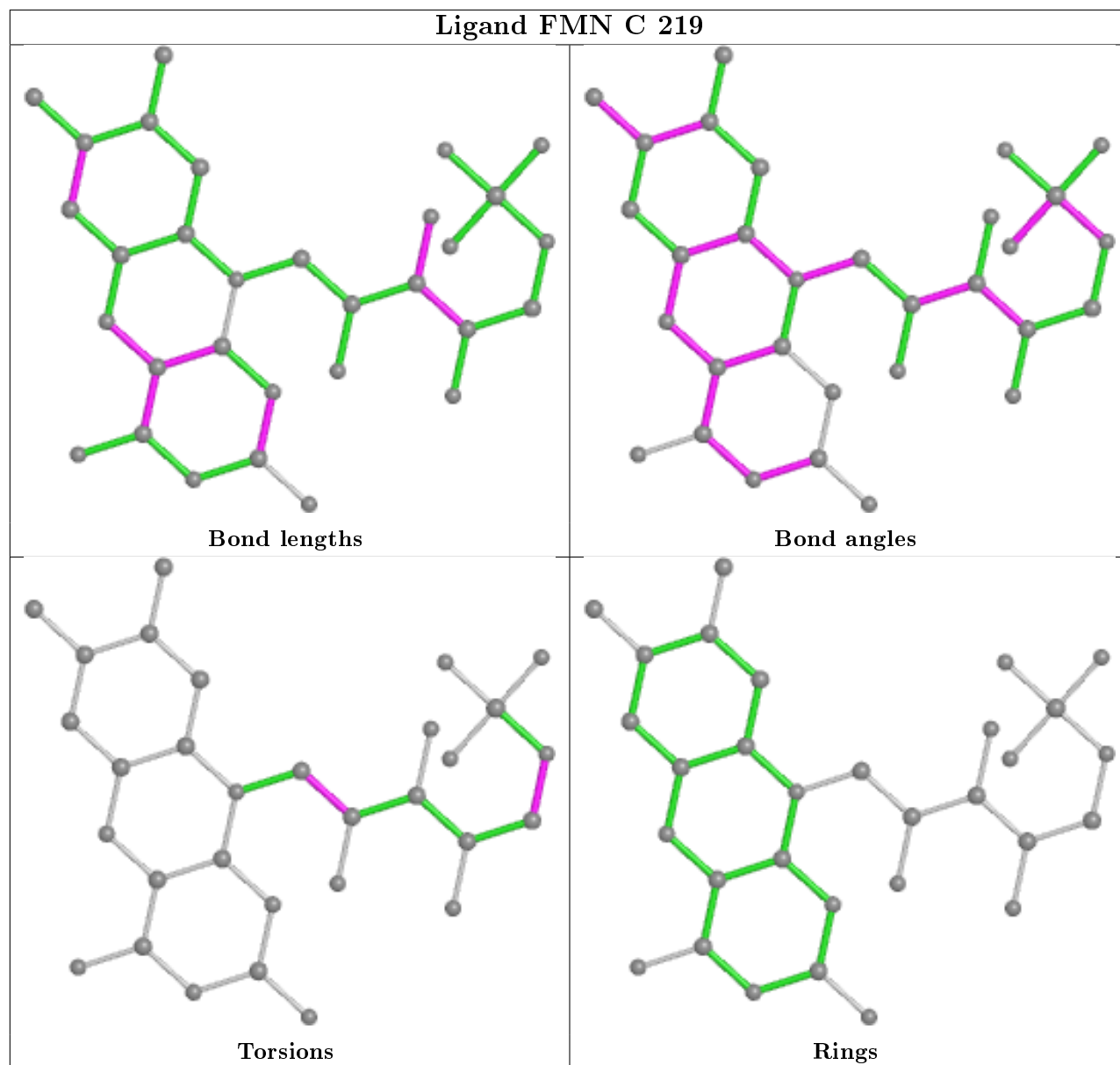
There are no ring outliers.

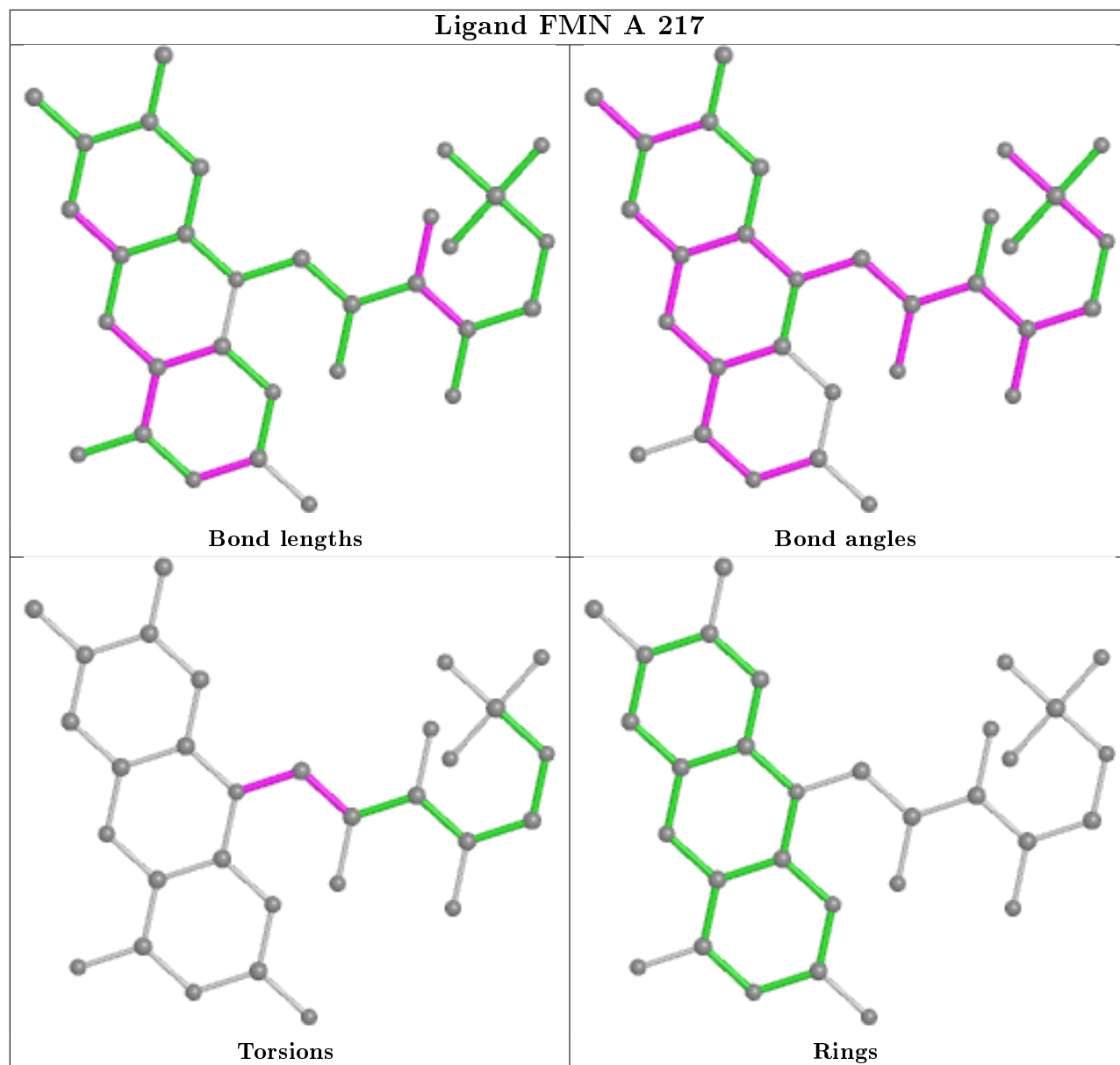
4 monomers are involved in 4 short contacts:

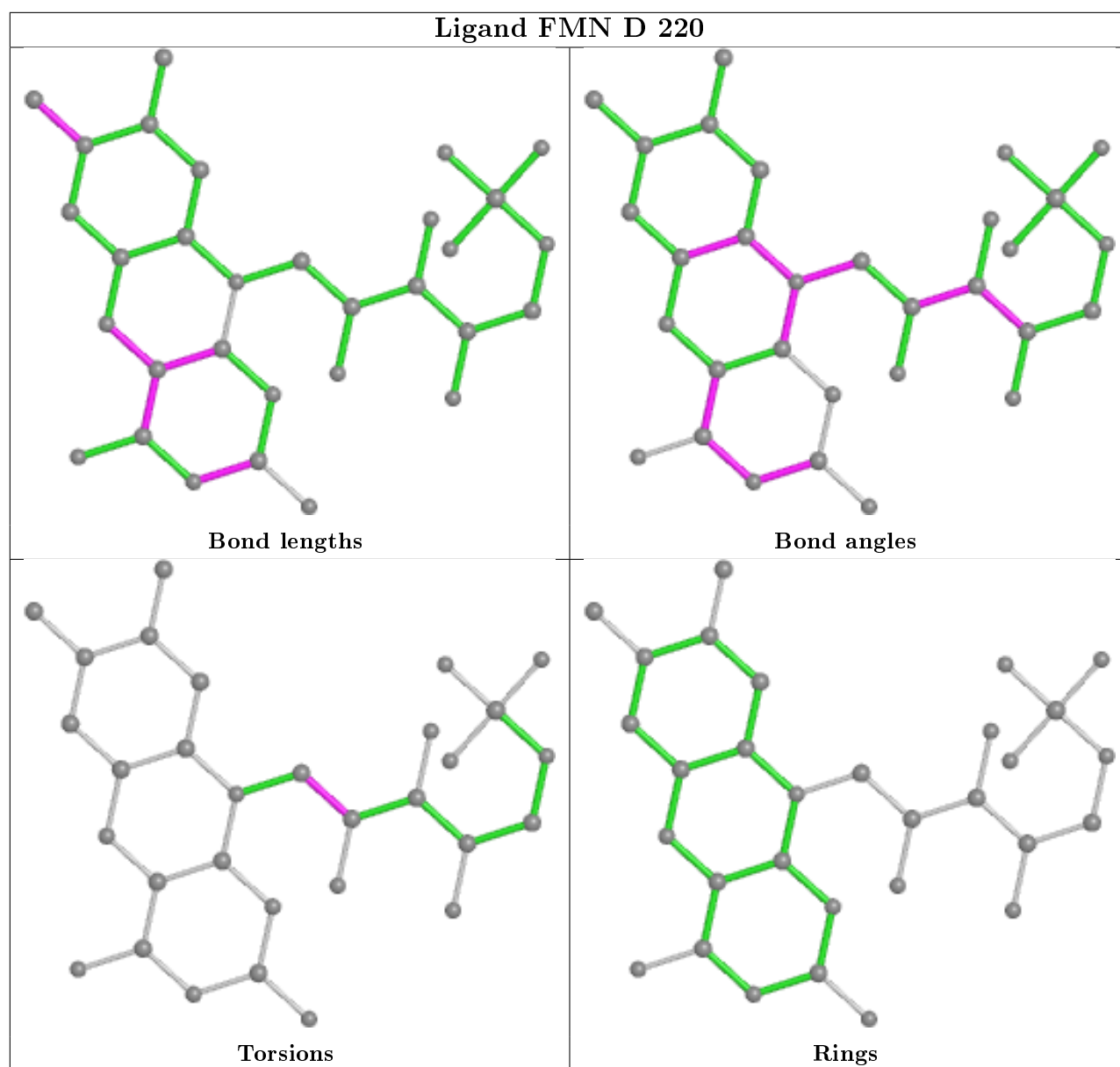
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	218	FMN	1	0
2	C	219	FMN	1	0
2	A	217	FMN	1	0
2	D	220	FMN	1	0

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 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 ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

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	216/216 (100%)	0.65	26 (12%) 4 7	11, 23, 100, 110	0
1	B	216/216 (100%)	-0.13	2 (0%) 84 89	11, 22, 35, 47	0
1	C	216/216 (100%)	0.17	17 (7%) 12 19	10, 21, 60, 81	0
1	D	216/216 (100%)	-0.13	1 (0%) 91 94	11, 22, 35, 45	0
All	All	864/864 (100%)	0.14	46 (5%) 26 35	10, 22, 56, 110	0

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	118	LYS	11.4
1	A	131	LEU	10.0
1	A	112	ALA	9.3
1	A	109	THR	9.1
1	A	129	VAL	8.6
1	A	116	ASN	8.4
1	A	117	HIS	8.3
1	A	123	PHE	7.3
1	C	129	VAL	6.8
1	A	122	TYR	6.6
1	A	114	ALA	6.4
1	C	109	THR	6.3
1	C	131	LEU	6.2
1	A	110	PRO	5.7
1	A	115	ALA	5.4
1	A	119	GLY	5.2
1	A	130	ASP	5.0
1	C	111	GLU	5.0
1	C	108	ASN	4.9
1	C	115	ALA	4.7
1	A	111	GLU	4.7

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Mol	Chain	Res	Type	RSRZ
1	A	132	LYS	4.6
1	A	121	THR	4.6
1	A	125	ASP	4.4
1	A	108	ASN	4.3
1	C	132	LYS	4.1
1	A	113	LYS	4.0
1	A	128	ARG	4.0
1	C	110	PRO	3.9
1	C	126	MET	3.7
1	A	124	ALA	3.7
1	B	69	PHE	3.6
1	C	118	LYS	3.6
1	C	122	TYR	3.4
1	A	126	MET	3.3
1	C	123	PHE	3.2
1	C	112	ALA	3.2
1	A	120	ARG	3.1
1	C	130	ASP	3.0
1	C	128	ARG	2.8
1	A	127	HIS	2.7
1	C	124	ALA	2.7
1	D	69	PHE	2.5
1	B	173	GLU	2.4
1	C	144	LEU	2.1
1	A	179	GLU	2.0

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

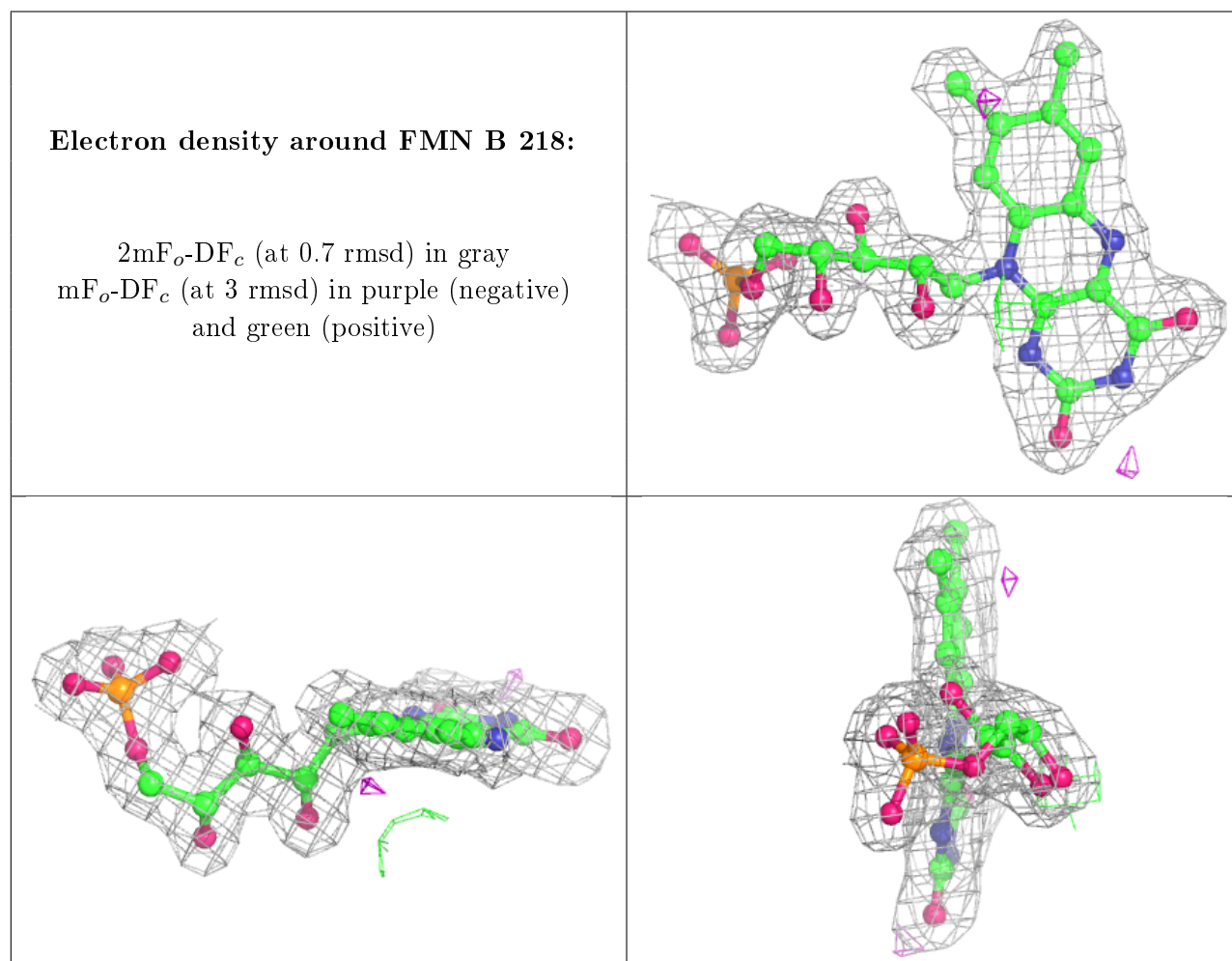
## 6.4 Ligands ⓘ

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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.



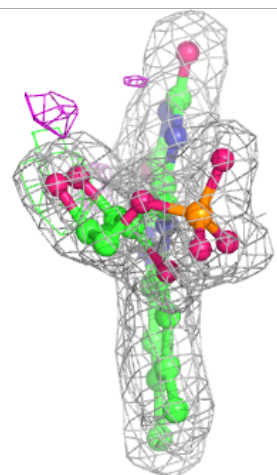
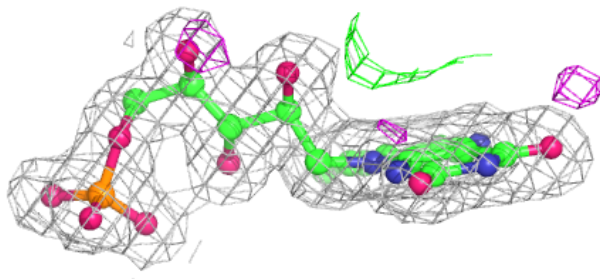
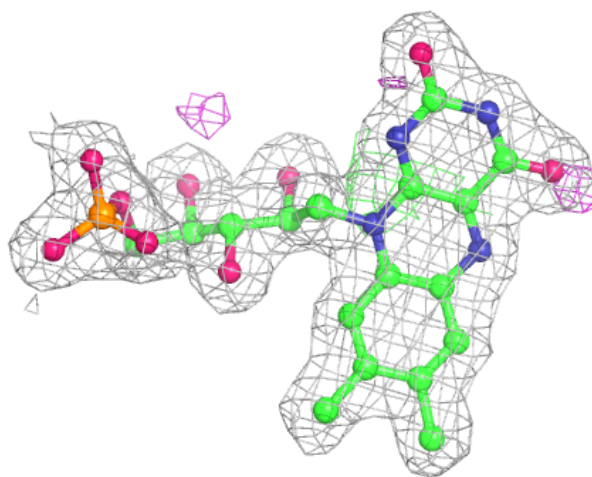
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	FMN	B	218	31/31	0.98	0.11	15,19,21,23	0
2	FMN	A	217	31/31	0.98	0.10	10,12,15,15	0
2	FMN	D	220	31/31	0.98	0.11	13,17,21,26	0
2	FMN	C	219	31/31	0.99	0.10	10,13,14,17	0

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.



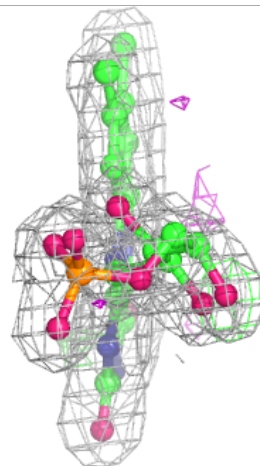
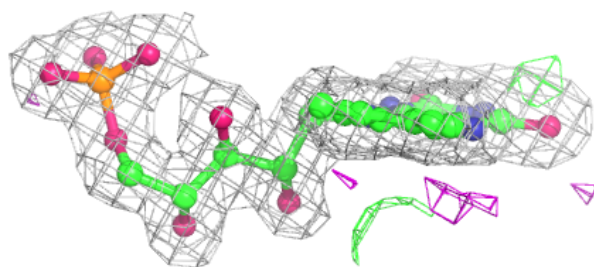
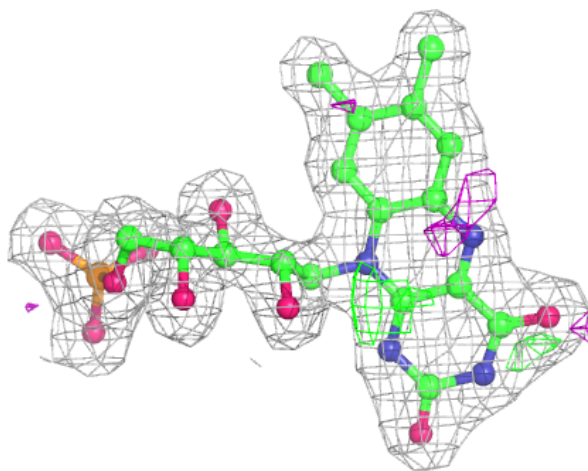
**Electron density around FMN A 217:**

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



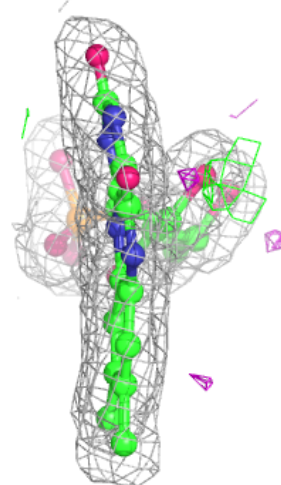
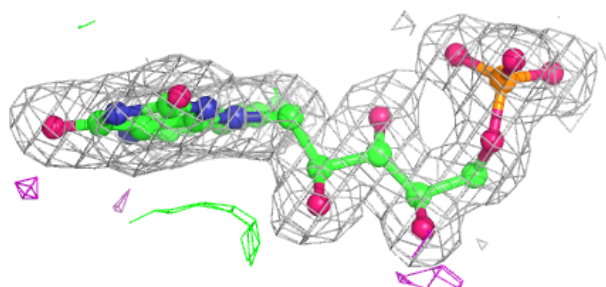
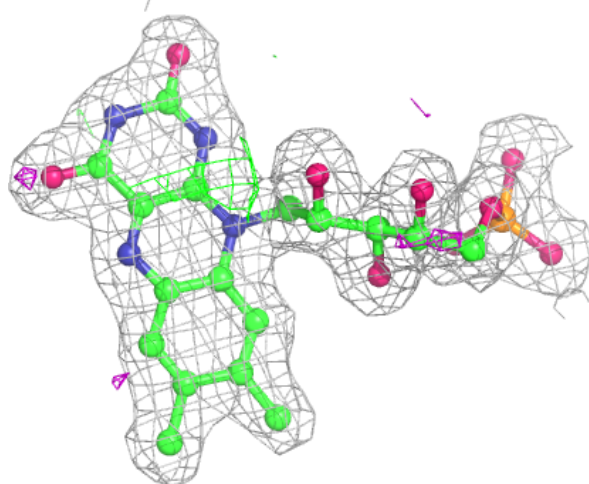
**Electron density around FMN D 220:**

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



**Electron density around FMN C 219:**

$2mF_o-DF_c$  (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.