



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 22, 2020 – 08:27 AM BST

PDB ID : 4H6P
Title : Crystal structure of a putative chromate reductase from *Gluconacetobacter hansenii*, Gh-ChrR, containing a R101A substitution.
Authors : Zhang, Y.; Robinson, H.; Buchko, G.W.
Deposited on : 2012-09-19
Resolution : 2.56 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 ⓘ 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 : 2.13.1
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) : 2.13.1

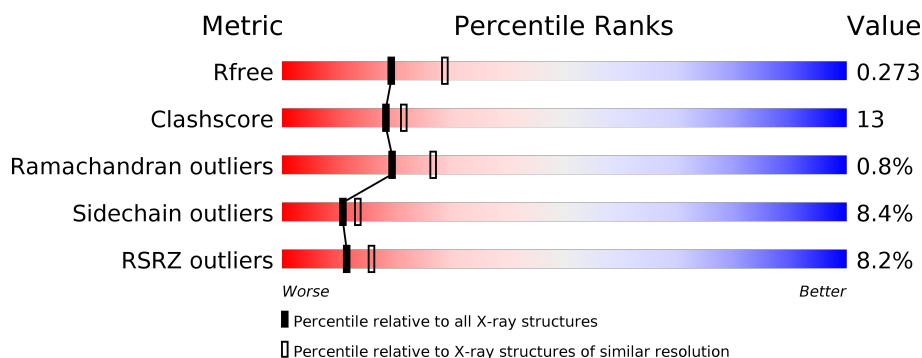
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.56 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	1279 (2.58-2.54)
Clashscore	141614	1327 (2.58-2.54)
Ramachandran outliers	138981	1312 (2.58-2.54)
Sidechain outliers	138945	1312 (2.58-2.54)
RSRZ outliers	127900	1269 (2.58-2.54)

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 $\leq 5\%$. 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	199	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>13%</div> <div>•</div> <div>8%</div> </div> </div>
1	B	199	<div> <div></div> <div>79%</div> <div>12%</div> <div>•</div> <div>8%</div> </div>
1	C	199	<div> <div>%</div> <div> <div></div> <div>75%</div> <div>16%</div> <div>•</div> <div>8%</div> </div> </div>
1	D	199	<div> <div>2%</div> <div> <div></div> <div>79%</div> <div>10%</div> <div>• •</div> <div>8%</div> </div> </div>
1	E	199	<div> <div>%</div> <div> <div></div> <div>71%</div> <div>19%</div> <div>• •</div> <div>8%</div> </div> </div>
1	F	199	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>16%</div> <div>•</div> <div>8%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	199	<div><div></div><div>8%58%33%8%</div></div>
1	H	199	<div><div></div><div>18%53%36%8%</div></div>
1	I	199	<div><div></div><div>18%58%30%5%8%</div></div>
1	J	199	<div><div></div><div>21%40%48%8%</div></div>
1	K	199	<div><div></div><div>12%62%27%8%</div></div>
1	L	199	<div><div></div><div>9%56%32%5%8%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 17193 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 Chromate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			
1	B	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			
1	C	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			
1	D	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			
1	E	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			
1	F	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			
1	G	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			
1	H	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			
1	I	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			
1	J	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			
1	K	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			
1	L	184	Total	C	N	O	S	0	0	0
			1370	878	233	256	3			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	101	ALA	ARG	CONFLICT	UNP D5QFC5
A	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
A	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
A	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
A	197	HIS	-	EXPRESSION TAG	UNP D5QFC5

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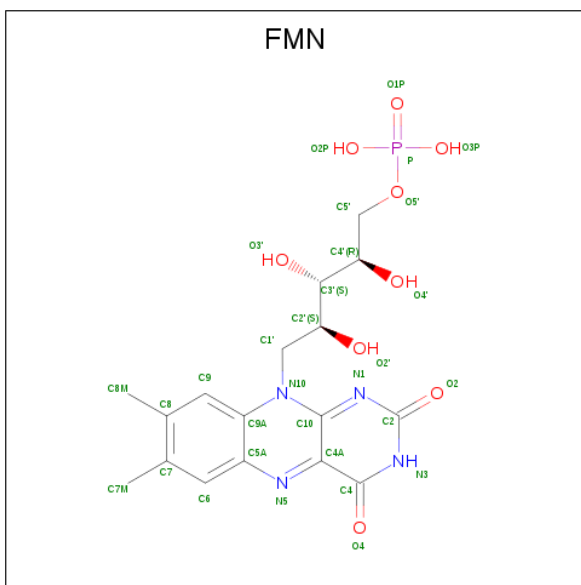
Chain	Residue	Modelled	Actual	Comment	Reference
A	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
A	199	HIS	-	EXPRESSION TAG	UNP D5QFC5
B	101	ALA	ARG	CONFLICT	UNP D5QFC5
B	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
B	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
B	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
B	197	HIS	-	EXPRESSION TAG	UNP D5QFC5
B	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
B	199	HIS	-	EXPRESSION TAG	UNP D5QFC5
C	101	ALA	ARG	CONFLICT	UNP D5QFC5
C	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
C	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
C	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
C	197	HIS	-	EXPRESSION TAG	UNP D5QFC5
C	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
C	199	HIS	-	EXPRESSION TAG	UNP D5QFC5
D	101	ALA	ARG	CONFLICT	UNP D5QFC5
D	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
D	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
D	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
D	197	HIS	-	EXPRESSION TAG	UNP D5QFC5
D	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
D	199	HIS	-	EXPRESSION TAG	UNP D5QFC5
E	101	ALA	ARG	CONFLICT	UNP D5QFC5
E	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
E	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
E	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
E	197	HIS	-	EXPRESSION TAG	UNP D5QFC5
E	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
E	199	HIS	-	EXPRESSION TAG	UNP D5QFC5
F	101	ALA	ARG	CONFLICT	UNP D5QFC5
F	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
F	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
F	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
F	197	HIS	-	EXPRESSION TAG	UNP D5QFC5
F	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
F	199	HIS	-	EXPRESSION TAG	UNP D5QFC5
G	101	ALA	ARG	CONFLICT	UNP D5QFC5
G	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
G	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
G	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
G	197	HIS	-	EXPRESSION TAG	UNP D5QFC5

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Chain	Residue	Modelled	Actual	Comment	Reference
G	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
G	199	HIS	-	EXPRESSION TAG	UNP D5QFC5
H	101	ALA	ARG	CONFLICT	UNP D5QFC5
H	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
H	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
H	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
H	197	HIS	-	EXPRESSION TAG	UNP D5QFC5
H	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
H	199	HIS	-	EXPRESSION TAG	UNP D5QFC5
I	101	ALA	ARG	CONFLICT	UNP D5QFC5
I	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
I	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
I	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
I	197	HIS	-	EXPRESSION TAG	UNP D5QFC5
I	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
I	199	HIS	-	EXPRESSION TAG	UNP D5QFC5
J	101	ALA	ARG	CONFLICT	UNP D5QFC5
J	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
J	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
J	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
J	197	HIS	-	EXPRESSION TAG	UNP D5QFC5
J	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
J	199	HIS	-	EXPRESSION TAG	UNP D5QFC5
K	101	ALA	ARG	CONFLICT	UNP D5QFC5
K	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
K	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
K	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
K	197	HIS	-	EXPRESSION TAG	UNP D5QFC5
K	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
K	199	HIS	-	EXPRESSION TAG	UNP D5QFC5
L	101	ALA	ARG	CONFLICT	UNP D5QFC5
L	194	HIS	-	EXPRESSION TAG	UNP D5QFC5
L	195	HIS	-	EXPRESSION TAG	UNP D5QFC5
L	196	HIS	-	EXPRESSION TAG	UNP D5QFC5
L	197	HIS	-	EXPRESSION TAG	UNP D5QFC5
L	198	HIS	-	EXPRESSION TAG	UNP D5QFC5
L	199	HIS	-	EXPRESSION TAG	UNP D5QFC5

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula: C₁₇H₂₁N₄O₉P).



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

- Molecule 3 is water.

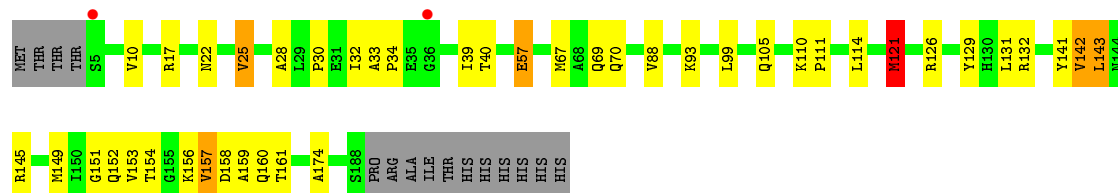
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	61	Total O 61 61	0	0
3	B	64	Total O 64 64	0	0
3	C	44	Total O 44 44	0	0
3	D	59	Total O 59 59	0	0
3	E	34	Total O 34 34	0	0
3	F	39	Total O 39 39	0	0
3	G	11	Total O 11 11	0	0
3	H	13	Total O 13 13	0	0
3	I	8	Total O 8 8	0	0
3	J	22	Total O 22 22	0	0
3	K	15	Total O 15 15	0	0
3	L	11	Total O 11 11	0	0

- Molecule 1: Chromate reductase




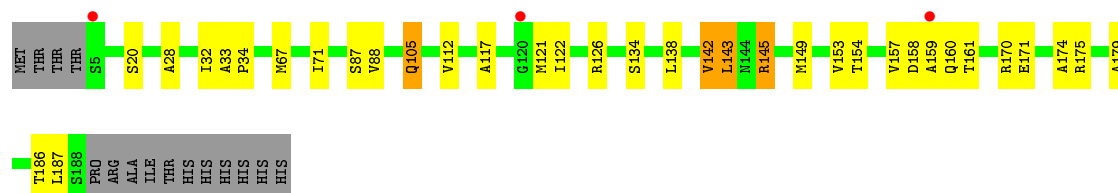
- Molecule 1: Chromate reductase

Chain E: 



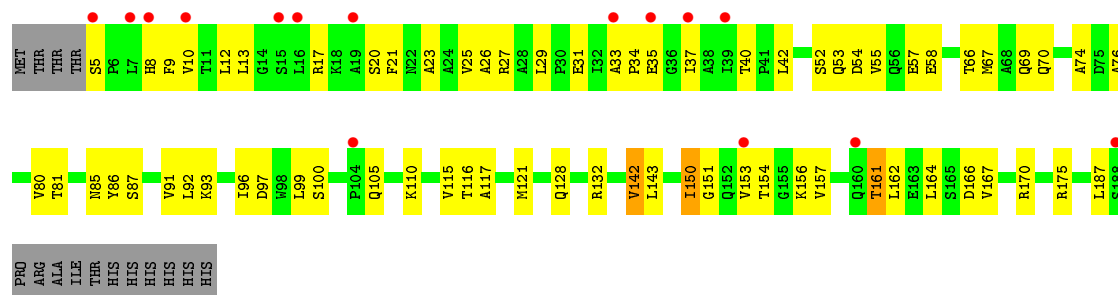
- Molecule 1: Chromate reductase

Chain F: 



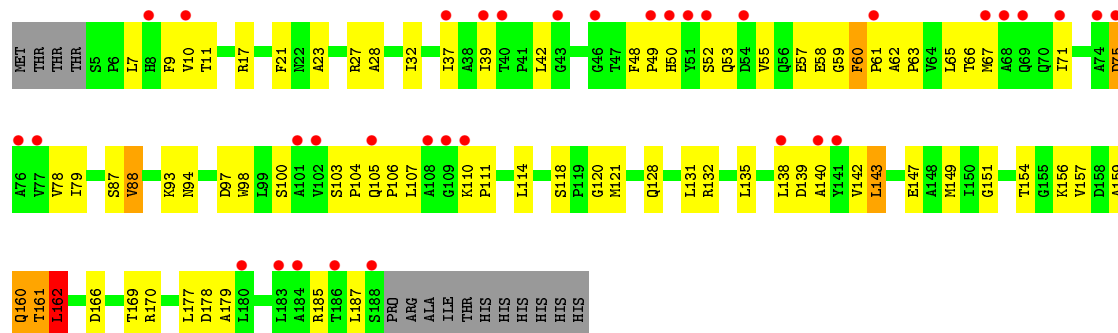
- Molecule 1: Chromate reductase

Chain G: 

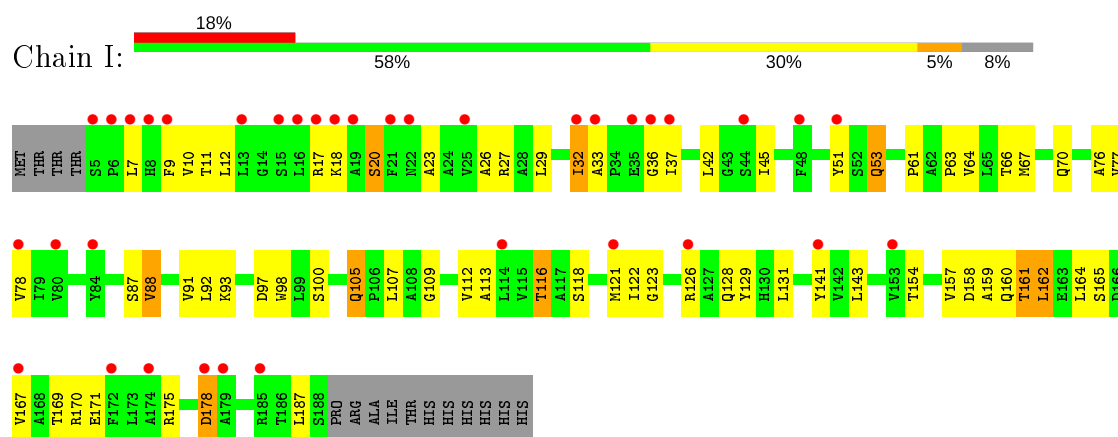


- Molecule 1: Chromate reductase

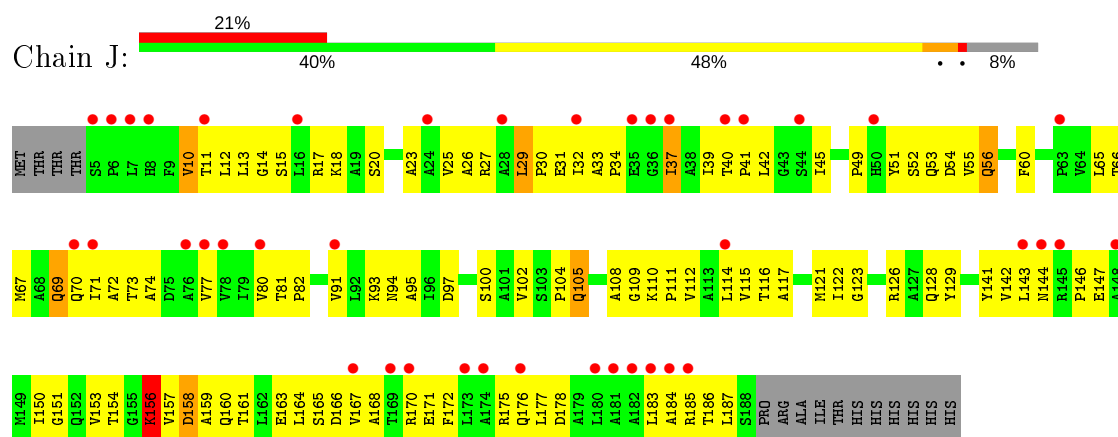
Chain H: 



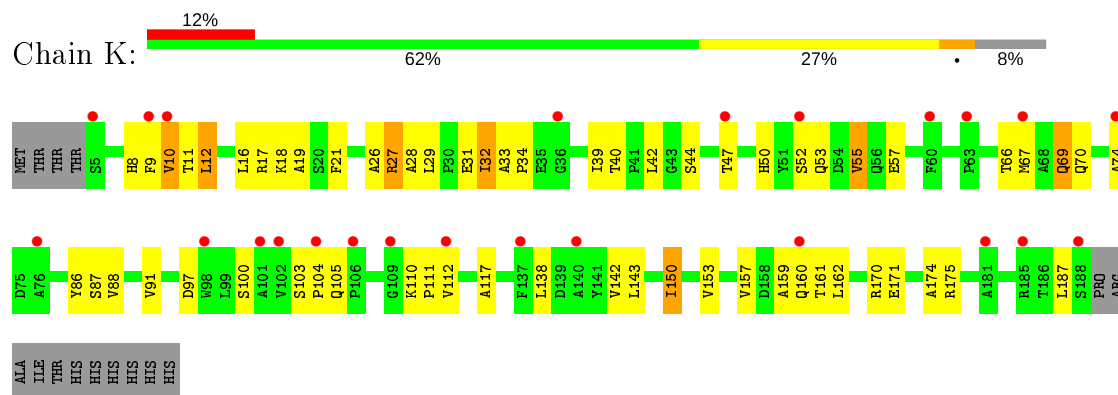
- Molecule 1: Chromate reductase



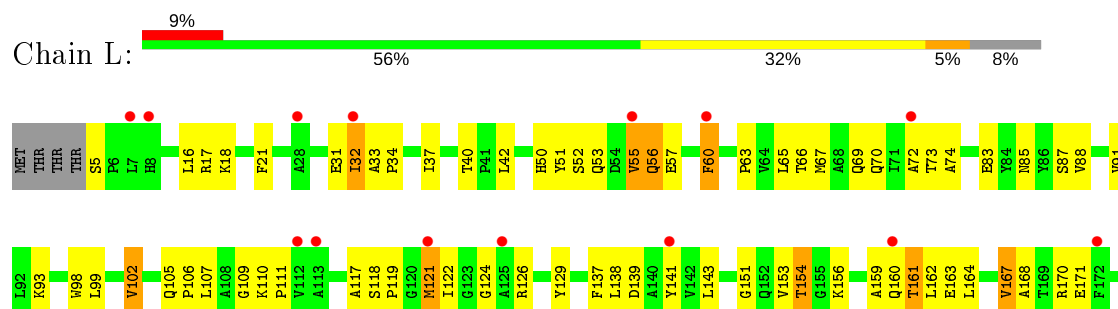
• Molecule 1: Chromate reductase



• Molecule 1: Chromate reductase



• Molecule 1: Chromate reductase





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	94.92Å 287.59Å 91.36Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.19 – 2.56 49.19 – 2.56	Depositor EDS
% Data completeness (in resolution range)	99.6 (49.19-2.56) 99.6 (49.19-2.56)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 2.54Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.233 , 0.271 0.235 , 0.273	Depositor DCC
R_{free} test set	4089 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	44.7	Xtriage
Anisotropy	0.041	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 44.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.034 for l,-k,h	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	17193	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% of the height of the origin peak. No significant pseudotranslation is detected.*

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: 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	0.42	0/1399	0.65	0/1915
1	B	0.41	0/1399	0.67	0/1915
1	C	0.39	0/1399	0.64	0/1915
1	D	0.37	0/1399	0.64	0/1915
1	E	0.40	0/1399	0.63	1/1915 (0.1%)
1	F	0.37	0/1399	0.63	1/1915 (0.1%)
1	G	0.31	0/1399	0.59	0/1915
1	H	0.33	0/1399	0.63	0/1915
1	I	0.32	0/1399	0.60	0/1915
1	J	0.40	0/1399	0.73	2/1915 (0.1%)
1	K	0.32	0/1399	0.62	0/1915
1	L	0.35	0/1399	0.65	1/1915 (0.1%)
All	All	0.37	0/16788	0.64	5/22980 (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	D	0	1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	121	MET	CB-CG-SD	7.24	134.11	112.40
1	L	121	MET	CG-SD-CE	6.96	111.33	100.20
1	J	143	LEU	CA-CB-CG	5.96	129.02	115.30
1	J	156	LYS	CA-CB-CG	5.69	125.91	113.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	145	ARG	NE-CZ-NH1	-5.25	117.68	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	162	LEU	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	1370	0	1397	14	0
1	B	1370	0	1397	15	0
1	C	1370	0	1397	22	0
1	D	1370	0	1397	19	0
1	E	1370	0	1397	30	0
1	F	1370	0	1397	20	0
1	G	1370	0	1397	47	0
1	H	1370	0	1397	62	0
1	I	1370	0	1397	54	0
1	J	1370	0	1397	100	0
1	K	1370	0	1397	43	0
1	L	1370	0	1397	60	0
2	A	31	0	19	0	0
2	B	31	0	19	0	0
2	C	31	0	19	0	0
2	D	31	0	19	0	0
2	E	31	0	19	1	0
2	F	31	0	19	0	0
2	G	31	0	19	4	0
2	H	31	0	19	5	0
2	I	31	0	19	6	0
2	J	31	0	19	6	0
2	K	31	0	19	1	0
2	L	31	0	19	3	0
3	A	61	0	0	3	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	64	0	0	3	1
3	C	44	0	0	3	0
3	D	59	0	0	1	0
3	E	34	0	0	3	0
3	F	39	0	0	3	0
3	G	11	0	0	3	0
3	H	13	0	0	6	0
3	I	8	0	0	5	0
3	J	22	0	0	15	0
3	K	15	0	0	4	0
3	L	11	0	0	7	0
All	All	17193	0	16992	449	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 13.

All (449) 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:L:170:ARG:NH1	3:L:310:HOH:O	1.83	1.10
1:F:145:ARG:HH12	1:L:139:ASP:HB2	1.28	0.99
1:J:156:LYS:NZ	1:J:166:ASP:H	1.60	0.98
1:I:10:VAL:HB	1:I:42:LEU:HD11	1.46	0.97
1:K:10:VAL:HG13	1:K:42:LEU:HD11	1.49	0.94
1:K:171:GLU:OE1	3:K:309:HOH:O	1.88	0.89
1:J:156:LYS:HZ3	1:J:166:ASP:H	1.20	0.88
1:J:12:LEU:HA	3:J:308:HOH:O	1.74	0.85
1:H:107:LEU:HB3	1:H:138:LEU:HD13	1.59	0.85
1:H:10:VAL:HB	1:H:42:LEU:HD11	1.57	0.84
2:H:201:FMN:O2P	3:H:312:HOH:O	1.96	0.84
1:I:126:ARG:NH1	1:J:129:TYR:OH	2.11	0.83
1:J:65:LEU:O	3:J:319:HOH:O	1.97	0.82
1:H:52:SER:HB3	1:H:55:VAL:HG12	1.63	0.81
2:L:201:FMN:N5	3:L:309:HOH:O	2.13	0.81
1:L:105:GLN:HE22	1:L:107:LEU:HB2	1.44	0.81
1:E:70:GLN:OE1	3:E:324:HOH:O	2.00	0.80
1:H:151:GLY:O	1:H:156:LYS:NZ	2.16	0.78
1:F:186:THR:O	3:F:312:HOH:O	2.00	0.78
1:J:153:VAL:O	1:J:156:LYS:HB3	1.84	0.77
1:G:85:ASN:H	2:G:201:FMN:H6	1.48	0.77
1:J:15:SER:OG	2:J:201:FMN:O3P	2.01	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:129:TYR:OH	1:L:126:ARG:NH1	2.19	0.75
1:K:52:SER:HB3	1:K:55:VAL:HG22	1.67	0.75
1:I:178:ASP:OD2	3:I:303:HOH:O	2.03	0.75
1:L:74:ALA:O	1:L:110:LYS:NZ	2.20	0.75
1:J:82:PRO:HB3	2:J:201:FMN:H5'2	1.69	0.74
1:C:161:THR:O	3:C:344:HOH:O	2.06	0.74
1:J:74:ALA:O	1:J:110:LYS:NZ	2.21	0.74
1:L:151:GLY:O	1:L:156:LYS:NZ	2.18	0.74
1:F:145:ARG:HH12	1:L:139:ASP:CB	2.00	0.73
1:G:74:ALA:O	1:G:110:LYS:NZ	2.21	0.73
1:H:88:VAL:HG21	1:H:131:LEU:HD13	1.71	0.73
1:H:59:GLY:HA3	3:H:310:HOH:O	1.89	0.73
1:E:114:LEU:O	3:E:313:HOH:O	2.08	0.72
1:J:14:GLY:N	1:J:80:VAL:O	2.22	0.72
1:J:156:LYS:HZ3	1:J:166:ASP:N	1.88	0.72
1:A:54:ASP:OD1	3:A:346:HOH:O	2.07	0.71
1:J:156:LYS:NZ	1:J:166:ASP:N	2.38	0.71
1:F:20:SER:O	3:F:329:HOH:O	2.08	0.71
1:H:132:ARG:NH2	1:H:143:LEU:O	2.23	0.71
1:I:129:TYR:OH	1:J:126:ARG:NH1	2.23	0.71
1:D:161:THR:OG1	1:D:162:LEU:N	2.21	0.71
1:J:156:LYS:O	1:J:156:LYS:NZ	2.24	0.71
1:H:58:GLU:O	3:H:310:HOH:O	2.07	0.71
1:D:145:ARG:NH2	3:D:322:HOH:O	2.15	0.70
1:H:32:ILE:HD11	1:H:170:ARG:HG2	1.73	0.69
1:H:60:PHE:HB3	1:H:65:LEU:HD21	1.73	0.69
1:D:32:ILE:HD11	1:D:170:ARG:HG2	1.74	0.69
1:H:79:ILE:HD12	1:H:131:LEU:HD21	1.74	0.69
1:B:17:ARG:NH1	3:B:351:HOH:O	2.10	0.69
1:A:154:THR:OG1	3:A:349:HOH:O	2.11	0.68
1:H:10:VAL:HB	1:H:42:LEU:CD1	2.24	0.67
1:I:17:ARG:NH1	2:I:201:FMN:O2P	2.28	0.67
1:J:56:GLN:O	3:J:315:HOH:O	2.12	0.67
1:A:32:ILE:HD11	1:A:170:ARG:HG2	1.76	0.67
1:J:69:GLN:N	3:J:319:HOH:O	2.11	0.67
1:J:52:SER:HB3	1:J:55:VAL:HB	1.77	0.66
1:J:178:ASP:O	3:J:312:HOH:O	2.12	0.66
1:K:31:GLU:OE2	3:K:313:HOH:O	2.12	0.66
1:E:159:ALA:O	1:E:161:THR:N	2.26	0.66
1:J:31:GLU:OE2	1:J:170:ARG:NH1	2.28	0.66
1:L:21:PHE:HE2	1:L:154:THR:HA	1.60	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:32:ILE:HD11	1:B:170:ARG:HG2	1.77	0.65
1:G:151:GLY:O	1:G:156:LYS:NZ	2.23	0.65
1:G:128:GLN:OE1	3:G:305:HOH:O	2.14	0.65
1:K:17:ARG:NH2	2:K:201:FMN:O1P	2.30	0.65
1:G:85:ASN:HA	1:L:137:PHE:HD2	1.62	0.65
1:E:17:ARG:NH2	2:E:201:FMN:O1P	2.30	0.64
1:B:159:ALA:O	1:B:161:THR:N	2.25	0.64
1:G:8:HIS:HE1	1:G:40:THR:HG23	1.63	0.64
1:G:5:SER:N	3:G:302:HOH:O	2.30	0.64
1:H:63:PRO:HA	1:H:66:THR:HB	1.79	0.63
1:I:122:ILE:HG21	1:I:126:ARG:HH21	1.63	0.63
1:I:157:VAL:HG22	1:I:164:LEU:HA	1.81	0.63
1:L:60:PHE:HB2	1:L:65:LEU:HD11	1.80	0.63
1:H:132:ARG:HG2	1:H:142:VAL:HG21	1.81	0.63
1:K:112:VAL:HB	1:K:142:VAL:HG22	1.80	0.63
1:F:32:ILE:HD11	1:F:170:ARG:HG2	1.80	0.62
1:G:85:ASN:N	2:G:201:FMN:H6	2.13	0.62
1:J:156:LYS:HZ3	1:J:165:SER:N	1.96	0.62
1:I:10:VAL:HB	1:I:42:LEU:CD1	2.26	0.62
1:B:18:LYS:NZ	3:B:338:HOH:O	2.29	0.62
1:E:67:MET:HG2	1:E:99:LEU:HD21	1.82	0.61
1:K:28:ALA:O	1:K:32:ILE:HD12	2.00	0.61
1:K:21:PHE:HB3	1:K:153:VAL:HG21	1.81	0.61
1:H:107:LEU:HD13	1:H:138:LEU:HD22	1.82	0.61
1:I:42:LEU:HD23	1:I:67:MET:HG3	1.82	0.61
1:H:28:ALA:O	1:H:32:ILE:HD12	2.01	0.61
1:G:153:VAL:HA	1:G:156:LYS:HD2	1.83	0.61
1:I:159:ALA:O	1:I:161:THR:N	2.33	0.61
1:E:32:ILE:HG23	1:E:174:ALA:HB2	1.83	0.60
1:C:175:ARG:NH2	3:C:312:HOH:O	2.22	0.60
1:E:126:ARG:NH1	1:L:129:TYR:OH	2.34	0.60
1:L:18:LYS:O	3:L:307:HOH:O	2.16	0.60
1:E:121:MET:SD	1:E:152:GLN:NE2	2.75	0.60
1:E:151:GLY:O	1:E:156:LYS:NZ	2.33	0.60
1:H:53:GLN:O	1:H:57:GLU:HG2	2.02	0.60
1:J:17:ARG:N	3:K:305:HOH:O	2.35	0.60
1:C:93:LYS:HE3	1:D:87:SER:HB3	1.84	0.59
1:H:161:THR:O	1:H:161:THR:OG1	2.19	0.59
1:J:156:LYS:HZ2	1:J:166:ASP:H	1.49	0.59
1:K:32:ILE:HD11	1:K:170:ARG:HG2	1.83	0.59
1:J:147:GLU:O	3:J:306:HOH:O	2.16	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:159:ALA:O	1:F:161:THR:N	2.32	0.59
1:K:159:ALA:O	1:K:161:THR:N	2.32	0.59
1:C:91:VAL:HG13	1:D:50:HIS:CG	2.38	0.59
1:J:156:LYS:HZ3	1:J:165:SER:H	1.50	0.59
1:H:87:SER:HB3	1:I:93:LYS:HE3	1.85	0.58
1:J:18:LYS:O	3:J:317:HOH:O	2.17	0.58
1:B:53:GLN:O	1:B:57:GLU:HG2	2.03	0.58
1:K:32:ILE:HG23	1:K:174:ALA:HB2	1.84	0.58
1:J:167:VAL:O	1:J:171:GLU:N	2.34	0.58
1:J:13:LEU:O	1:J:45:ILE:HD13	2.03	0.58
1:G:87:SER:HB3	1:L:93:LYS:HE3	1.86	0.58
1:I:29:LEU:O	1:I:33:ALA:N	2.36	0.58
1:L:159:ALA:O	1:L:161:THR:N	2.32	0.58
1:J:186:THR:O	3:J:311:HOH:O	2.17	0.58
1:L:161:THR:O	1:L:163:GLU:HG2	2.03	0.58
1:K:16:LEU:HD11	1:K:91:VAL:HG11	1.84	0.57
1:K:66:THR:O	1:K:70:GLN:HG3	2.04	0.57
1:L:151:GLY:N	3:L:301:HOH:O	1.91	0.57
1:C:87:SER:HB3	1:D:93:LYS:HE3	1.85	0.57
1:J:32:ILE:HD11	1:J:170:ARG:HG2	1.85	0.57
1:F:171:GLU:O	1:F:175:ARG:HG2	2.05	0.57
1:I:23:ALA:O	1:I:27:ARG:HG3	2.05	0.57
1:J:49:PRO:O	1:J:94:ASN:ND2	2.32	0.57
1:L:67:MET:HG2	1:L:99:LEU:HD21	1.87	0.57
1:G:97:ASP:O	1:G:100:SER:OG	2.20	0.57
1:H:135:LEU:HD23	1:H:138:LEU:HD21	1.85	0.57
1:J:53:GLN:HB3	1:K:17:ARG:HD3	1.87	0.56
1:G:150:ILE:HG13	1:G:150:ILE:O	2.05	0.56
1:J:128:GLN:HG3	3:J:302:HOH:O	2.06	0.56
1:K:74:ALA:O	1:K:110:LYS:NZ	2.37	0.56
1:J:156:LYS:NZ	1:J:165:SER:H	2.03	0.56
1:L:122:ILE:HG13	1:L:126:ARG:NH2	2.19	0.56
1:G:17:ARG:HG2	1:G:20:SER:HB2	1.88	0.56
1:K:11:THR:O	1:K:42:LEU:HD13	2.05	0.56
1:I:122:ILE:HG23	1:J:122:ILE:HG23	1.85	0.56
1:J:45:ILE:HG23	1:J:95:ALA:HB2	1.86	0.56
1:C:50:HIS:CG	1:D:91:VAL:HG13	2.41	0.56
1:F:28:ALA:O	1:F:32:ILE:HD12	2.06	0.55
1:I:26:ALA:HA	1:I:29:LEU:HD12	1.88	0.55
1:I:7:LEU:N	1:I:36:GLY:O	2.39	0.55
1:E:110:LYS:HD3	1:E:111:PRO:HD2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:97:ASP:O	1:I:100:SER:OG	2.19	0.55
1:L:52:SER:HB3	1:L:55:VAL:HG23	1.89	0.55
1:J:10:VAL:CG2	1:J:42:LEU:HG	2.37	0.55
1:J:105:GLN:OE1	1:J:108:ALA:N	2.39	0.55
1:G:93:LYS:HE3	1:L:87:SER:HB3	1.88	0.55
1:H:93:LYS:O	1:H:97:ASP:N	2.37	0.55
1:I:109:GLY:HA2	1:I:141:TYR:CE2	2.42	0.55
1:K:53:GLN:O	1:K:57:GLU:HG3	2.07	0.55
1:K:117:ALA:HA	1:K:150:ILE:HG13	1.88	0.55
1:J:30:PRO:HG3	1:J:39:ILE:O	2.08	0.54
1:K:110:LYS:HD3	1:K:111:PRO:HD2	1.89	0.54
1:B:151:GLY:O	1:B:156:LYS:NZ	2.38	0.54
1:H:42:LEU:HD23	1:H:67:MET:HG3	1.89	0.54
1:J:109:GLY:HA2	1:J:141:TYR:CE2	2.43	0.54
1:H:162:LEU:HD22	1:H:162:LEU:H	1.72	0.54
1:J:146:PRO:O	1:J:176:GLN:NE2	2.38	0.54
1:J:159:ALA:O	1:J:161:THR:N	2.39	0.54
1:H:140:ALA:O	1:J:144:ASN:ND2	2.41	0.53
1:H:110:LYS:HD3	1:H:111:PRO:HD2	1.89	0.53
1:L:72:ALA:O	3:L:311:HOH:O	2.19	0.53
1:G:53:GLN:O	1:G:57:GLU:HG3	2.08	0.53
1:J:69:GLN:HA	1:J:72:ALA:HB3	1.89	0.53
1:E:153:VAL:HG22	1:E:157:VAL:HG12	1.91	0.53
1:K:12:LEU:HA	1:K:42:LEU:HD22	1.89	0.53
1:H:9:PHE:HB2	1:H:39:ILE:HD13	1.90	0.53
1:L:16:LEU:HD11	1:L:91:VAL:HG21	1.90	0.53
1:G:42:LEU:HD22	1:G:67:MET:HG3	1.90	0.53
1:I:159:ALA:C	1:I:161:THR:H	2.12	0.53
1:J:23:ALA:O	1:J:27:ARG:HG3	2.09	0.53
1:B:32:ILE:HG23	1:B:174:ALA:HB2	1.91	0.52
1:J:82:PRO:CB	2:J:201:FMN:H5'2	2.37	0.52
1:I:129:TYR:CE2	1:J:126:ARG:HD3	2.45	0.52
1:I:32:ILE:HD11	1:I:170:ARG:HD3	1.91	0.52
1:H:71:ILE:O	1:H:110:LYS:NZ	2.41	0.52
1:I:78:VAL:HG22	1:I:113:ALA:HB3	1.92	0.52
1:J:26:ALA:HB2	1:J:80:VAL:HG11	1.90	0.52
1:J:29:LEU:H	1:J:30:PRO:HD2	1.75	0.52
1:B:128:GLN:OE1	3:B:305:HOH:O	2.18	0.52
1:L:168:ALA:HA	3:L:308:HOH:O	2.09	0.52
1:L:53:GLN:O	1:L:53:GLN:HG2	2.10	0.52
1:G:161:THR:O	1:G:162:LEU:HD23	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:25:VAL:HA	1:J:164:LEU:HD22	1.92	0.52
1:G:52:SER:HB3	1:G:55:VAL:HB	1.91	0.52
1:J:17:ARG:O	1:J:20:SER:OG	2.28	0.52
1:E:69:GLN:NE2	3:E:331:HOH:O	2.23	0.51
1:H:159:ALA:O	1:H:161:THR:N	2.42	0.51
1:I:20:SER:HA	2:I:201:FMN:O1P	2.10	0.51
1:E:30:PRO:HA	1:E:39:ILE:HB	1.92	0.51
1:C:48:PHE:O	1:D:50:HIS:HE1	1.93	0.51
1:I:51:TYR:HA	1:I:98:TRP:CZ2	2.46	0.51
1:J:166:ASP:OD2	1:J:168:ALA:HB3	2.10	0.51
1:J:13:LEU:HD11	1:J:23:ALA:HA	1.93	0.51
1:L:167:VAL:O	1:L:171:GLU:N	2.39	0.51
1:J:156:LYS:HD2	1:J:166:ASP:HB2	1.93	0.51
1:K:42:LEU:HG	1:K:67:MET:SD	2.51	0.51
1:I:126:ARG:HD3	1:J:129:TYR:CE2	2.46	0.51
1:A:159:ALA:O	1:A:161:THR:N	2.43	0.51
1:F:122:ILE:HG13	1:F:126:ARG:NH2	2.26	0.51
1:I:105:GLN:HE22	1:I:107:LEU:HB2	1.76	0.51
1:J:51:TYR:N	1:J:94:ASN:OD1	2.40	0.51
1:B:121:MET:HE1	1:B:152:GLN:HG3	1.93	0.50
1:H:17:ARG:HD3	1:I:53:GLN:HB3	1.93	0.50
1:I:129:TYR:CZ	1:J:126:ARG:HD3	2.46	0.50
1:B:105:GLN:HE21	1:B:106:PRO:HD2	1.75	0.50
1:H:21:PHE:CD2	2:H:201:FMN:H5'2	2.46	0.50
2:I:201:FMN:HM81	3:I:306:HOH:O	2.11	0.50
1:J:27:ARG:NH2	3:J:317:HOH:O	2.29	0.50
1:H:23:ALA:O	1:H:27:ARG:HG3	2.12	0.50
1:E:153:VAL:HG22	1:E:157:VAL:CG1	2.42	0.50
1:E:93:LYS:HE3	1:F:87:SER:HB3	1.93	0.50
1:G:23:ALA:O	1:G:27:ARG:HG3	2.12	0.50
1:J:115:VAL:HB	3:J:321:HOH:O	2.11	0.50
1:A:105:GLN:HE21	1:A:106:PRO:HD2	1.77	0.50
1:B:66:THR:O	1:B:69:GLN:HG3	2.11	0.50
1:I:77:VAL:HG13	1:I:112:VAL:HG13	1.93	0.50
1:J:17:ARG:HD3	1:K:53:GLN:HB3	1.94	0.50
1:L:21:PHE:HD2	1:L:153:VAL:HG22	1.77	0.50
1:J:183:LEU:C	1:J:185:ARG:H	2.15	0.49
1:J:93:LYS:HE3	1:K:87:SER:HB3	1.94	0.49
1:H:147:GLU:OE1	1:K:86:TYR:OH	2.19	0.49
1:H:21:PHE:HD2	2:H:201:FMN:H5'2	1.76	0.49
1:C:50:HIS:HE1	1:D:48:PHE:O	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:109:GLY:HA2	1:J:141:TYR:HE2	1.76	0.49
1:F:158:ASP:HB2	3:F:327:HOH:O	2.11	0.49
1:G:67:MET:HG2	1:G:99:LEU:HD21	1.95	0.49
1:I:17:ARG:HH11	2:I:201:FMN:P	2.36	0.49
1:L:31:GLU:HG3	1:L:170:ARG:HH22	1.77	0.49
1:A:41:PRO:HD2	1:C:69:GLN:NE2	2.28	0.49
1:J:66:THR:O	3:J:319:HOH:O	2.20	0.49
1:G:85:ASN:HA	1:L:137:PHE:CD2	2.44	0.48
1:I:10:VAL:CB	1:I:42:LEU:HD11	2.30	0.48
1:J:51:TYR:OH	1:J:56:GLN:NE2	2.45	0.48
1:J:32:ILE:HG22	1:J:177:LEU:HD12	1.96	0.48
1:G:164:LEU:O	3:G:304:HOH:O	2.19	0.48
1:C:114:LEU:HD22	1:C:131:LEU:HD23	1.95	0.48
1:K:8:HIS:HE1	1:K:40:THR:HG23	1.79	0.48
1:L:122:ILE:HG13	1:L:126:ARG:HH21	1.78	0.48
1:J:60:PHE:CE1	1:J:102:VAL:HA	2.49	0.48
1:J:81:THR:O	1:J:116:THR:OG1	2.28	0.48
1:L:32:ILE:HG12	1:L:170:ARG:HG3	1.95	0.48
1:J:10:VAL:HG22	1:J:42:LEU:HG	1.95	0.48
1:J:158:ASP:OD1	1:J:161:THR:OG1	2.32	0.48
1:F:105:GLN:OE1	1:F:138:LEU:HB3	2.14	0.47
1:I:18:LYS:NZ	1:I:162:LEU:HD11	2.29	0.47
1:J:111:PRO:HA	1:J:141:TYR:O	2.14	0.47
1:E:121:MET:HG3	1:L:121:MET:SD	2.54	0.47
1:G:166:ASP:O	1:G:170:ARG:HG3	2.15	0.47
1:H:166:ASP:HB3	1:H:169:THR:HB	1.96	0.47
1:F:158:ASP:HB3	1:F:161:THR:HG22	1.96	0.47
1:G:117:ALA:HA	1:G:150:ILE:HG13	1.96	0.47
1:C:28:ALA:O	1:C:32:ILE:HD12	2.15	0.47
1:I:11:THR:O	1:I:42:LEU:HD13	2.14	0.47
1:D:32:ILE:HG23	1:D:174:ALA:HB2	1.96	0.47
1:G:27:ARG:CZ	1:G:162:LEU:HD12	2.44	0.47
1:G:157:VAL:HG22	1:G:164:LEU:HA	1.97	0.47
1:I:164:LEU:HD11	1:I:169:THR:HG21	1.97	0.47
1:J:171:GLU:O	1:J:175:ARG:HG2	2.15	0.47
1:L:21:PHE:CD2	1:L:153:VAL:HG13	2.50	0.47
1:B:105:GLN:NE2	1:B:106:PRO:HD2	2.29	0.47
1:L:119:PRO:HD2	2:L:201:FMN:O2'	2.14	0.47
1:G:26:ALA:HA	1:G:29:LEU:HD13	1.96	0.46
1:C:32:ILE:HD11	1:C:170:ARG:HG2	1.97	0.46
1:I:164:LEU:N	3:I:304:HOH:O	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:151:GLY:HA3	3:J:318:HOH:O	2.15	0.46
1:H:160:GLN:OE1	1:H:162:LEU:HD21	2.15	0.46
1:J:69:GLN:O	1:J:73:THR:HG23	2.16	0.46
1:C:53:GLN:O	1:C:57:GLU:HG3	2.15	0.46
1:I:109:GLY:HA2	1:I:141:TYR:HE2	1.80	0.46
1:J:77:VAL:HB	1:J:112:VAL:HG22	1.96	0.46
1:H:62:ALA:O	1:H:66:THR:N	2.36	0.46
1:L:110:LYS:HD3	1:L:111:PRO:HD2	1.98	0.46
1:L:83:GLU:CD	1:L:126:ARG:HB2	2.36	0.46
1:L:66:THR:O	1:L:70:GLN:HG3	2.16	0.46
1:H:93:LYS:NZ	1:I:88:VAL:O	2.49	0.46
1:J:128:GLN:CG	3:J:302:HOH:O	2.64	0.46
1:G:31:GLU:HG2	1:G:170:ARG:HH22	1.81	0.46
1:G:33:ALA:HA	1:G:34:PRO:HD3	1.84	0.46
1:H:114:LEU:O	3:H:303:HOH:O	2.20	0.46
1:L:117:ALA:O	3:L:306:HOH:O	2.21	0.46
1:D:132:ARG:NH1	1:D:142:VAL:HG22	2.30	0.45
1:E:28:ALA:O	1:E:32:ILE:HD13	2.15	0.45
1:I:171:GLU:O	1:I:175:ARG:HG3	2.16	0.45
1:A:31:GLU:OE1	1:A:31:GLU:N	2.49	0.45
1:C:27:ARG:NH2	1:C:162:LEU:O	2.50	0.45
1:H:50:HIS:HA	1:H:94:ASN:HD21	1.81	0.45
1:I:126:ARG:HD3	1:J:129:TYR:CZ	2.52	0.45
1:A:159:ALA:C	1:A:161:THR:H	2.19	0.45
1:C:185:ARG:HD3	3:C:342:HOH:O	2.15	0.45
1:D:105:GLN:NE2	1:D:138:LEU:HB3	2.32	0.45
1:D:28:ALA:O	1:D:32:ILE:HD12	2.16	0.45
1:D:161:THR:O	1:D:162:LEU:HB3	2.17	0.45
1:G:21:PHE:O	1:G:25:VAL:HG23	2.16	0.45
1:I:23:ALA:HB3	3:I:301:HOH:O	2.17	0.45
1:J:60:PHE:HB2	1:J:65:LEU:HD11	1.99	0.45
1:A:105:GLN:HA	1:A:106:PRO:HD3	1.88	0.45
1:C:105:GLN:HE21	1:C:106:PRO:HD2	1.80	0.45
1:J:117:ALA:HB2	1:J:150:ILE:HD12	1.98	0.45
1:J:29:LEU:H	1:J:30:PRO:CD	2.29	0.45
1:G:92:LEU:O	1:G:96:ILE:HD12	2.16	0.45
1:J:166:ASP:O	1:J:170:ARG:HB2	2.16	0.45
1:J:54:ASP:OD2	1:K:18:LYS:N	2.36	0.45
1:E:132:ARG:NH2	1:E:143:LEU:O	2.41	0.45
1:I:12:LEU:HD21	1:I:92:LEU:HD11	1.99	0.45
1:H:120:GLY:N	3:H:302:HOH:O	2.34	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:156:LYS:HZ3	1:J:165:SER:CA	2.30	0.45
1:K:150:ILE:HG13	1:K:150:ILE:O	2.16	0.45
1:A:67:MET:O	1:A:71:ILE:HG13	2.16	0.45
1:F:67:MET:O	1:F:71:ILE:HG13	2.15	0.45
1:G:35:GLU:N	1:G:35:GLU:OE1	2.44	0.45
1:H:93:LYS:HZ2	1:I:87:SER:HB3	1.82	0.45
1:I:64:VAL:HG21	1:I:98:TRP:HB3	1.98	0.45
1:L:159:ALA:C	1:L:161:THR:H	2.18	0.45
1:F:117:ALA:HB1	1:F:153:VAL:HG21	1.98	0.45
1:K:10:VAL:CG1	1:K:42:LEU:HD11	2.33	0.45
1:L:69:GLN:N	1:L:69:GLN:OE1	2.46	0.45
1:L:85:ASN:H	2:L:201:FMN:H6	1.82	0.45
1:H:114:LEU:HD22	1:H:131:LEU:HD23	1.99	0.44
1:J:154:THR:C	1:J:156:LYS:H	2.20	0.44
1:G:85:ASN:OD1	1:L:138:LEU:HD21	2.16	0.44
1:A:32:ILE:HG23	1:A:174:ALA:HB2	2.00	0.44
1:J:12:LEU:CD1	1:J:45:ILE:HD11	2.47	0.44
1:H:61:PRO:O	1:H:65:LEU:HG	2.18	0.44
1:L:56:GLN:NE2	1:L:57:GLU:HG2	2.33	0.44
1:B:105:GLN:OE1	1:B:138:LEU:HB3	2.17	0.44
3:A:349:HOH:O	1:E:57:GLU:HG3	2.17	0.44
2:G:201:FMN:H1'1	2:G:201:FMN:H9	1.60	0.44
1:H:177:LEU:C	1:H:179:ALA:H	2.21	0.44
1:H:9:PHE:HB2	1:H:39:ILE:CD1	2.47	0.44
1:J:33:ALA:HB1	1:J:37:ILE:HB	1.99	0.44
1:L:60:PHE:CD1	1:L:65:LEU:HD11	2.53	0.44
1:F:112:VAL:O	1:F:142:VAL:HA	2.17	0.44
1:H:132:ARG:HA	1:H:135:LEU:HB2	1.99	0.44
1:H:107:LEU:CB	1:H:138:LEU:HD13	2.40	0.44
1:A:105:GLN:OE1	1:A:138:LEU:HB3	2.18	0.44
1:G:86:TYR:H	2:G:201:FMN:HM72	1.83	0.44
1:A:32:ILE:HD11	1:A:170:ARG:CG	2.46	0.44
1:C:78:VAL:HG22	1:C:113:ALA:HB3	2.00	0.44
1:J:17:ARG:NH2	2:J:201:FMN:O1P	2.48	0.44
1:K:33:ALA:HA	1:K:34:PRO:HD3	1.83	0.44
1:I:118:SER:HB3	1:I:123:GLY:HA2	1.99	0.43
1:J:102:VAL:HG12	1:J:104:PRO:O	2.18	0.43
1:H:17:ARG:NH2	2:H:201:FMN:O5'	2.45	0.43
1:F:143:LEU:HD11	1:F:179:ALA:HB3	2.00	0.43
1:L:55:VAL:O	1:L:55:VAL:HG12	2.18	0.43
1:C:158:ASP:OD1	1:C:161:THR:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:111:PRO:HA	1:E:141:TYR:O	2.18	0.43
1:G:81:THR:HB	1:G:92:LEU:HD22	2.00	0.43
1:I:66:THR:O	1:I:70:GLN:HG3	2.19	0.43
1:J:116:THR:HG23	1:J:123:GLY:O	2.19	0.43
1:K:26:ALA:HA	1:K:29:LEU:HD12	2.01	0.43
1:H:11:THR:O	1:H:42:LEU:HD13	2.19	0.43
1:K:87:SER:OG	1:K:88:VAL:N	2.48	0.43
1:C:15:SER:HB2	1:C:82:PRO:HG3	2.00	0.43
1:F:105:GLN:O	1:F:105:GLN:HG3	2.19	0.43
1:K:105:GLN:NE2	1:K:138:LEU:O	2.52	0.43
1:K:19:ALA:HA	1:K:162:LEU:HD21	2.01	0.43
1:H:75:ASP:OD1	1:H:75:ASP:N	2.52	0.43
1:I:97:ASP:OD2	3:I:305:HOH:O	2.21	0.43
1:J:11:THR:OG1	1:J:41:PRO:HA	2.18	0.43
1:K:9:PHE:HB2	1:K:39:ILE:HG12	2.01	0.43
1:I:9:PHE:CD1	1:I:76:ALA:HB3	2.54	0.43
1:J:114:LEU:HD12	1:J:128:GLN:HG2	2.01	0.43
1:E:114:LEU:HD22	1:E:131:LEU:HD23	2.00	0.42
1:E:132:ARG:NH1	1:E:142:VAL:HG22	2.34	0.42
1:H:160:GLN:HA	1:H:162:LEU:HD13	2.01	0.42
1:G:66:THR:O	1:G:69:GLN:HG3	2.19	0.42
1:H:49:PRO:O	1:H:98:TRP:NE1	2.43	0.42
1:J:91:VAL:HG23	1:K:50:HIS:CD2	2.54	0.42
1:K:103:SER:HA	1:K:104:PRO:HA	1.84	0.42
1:B:105:GLN:HA	1:B:106:PRO:HD3	1.93	0.42
1:G:12:LEU:HD12	1:G:13:LEU:N	2.33	0.42
1:K:44:SER:HB3	1:K:47:THR:HG23	2.01	0.42
1:L:102:VAL:HG12	1:L:106:PRO:HD3	2.02	0.42
1:E:126:ARG:HD3	1:L:129:TYR:CE2	2.54	0.42
1:E:159:ALA:C	1:E:161:THR:H	2.19	0.42
2:I:201:FMN:H9	2:I:201:FMN:H1'2	1.87	0.42
1:K:67:MET:O	1:K:70:GLN:HB2	2.19	0.42
1:G:17:ARG:CB	1:L:53:GLN:HB3	2.50	0.42
1:C:50:HIS:CD2	1:D:91:VAL:HG13	2.55	0.42
1:E:22:ASN:O	1:E:25:VAL:HG13	2.20	0.42
1:H:97:ASP:O	1:H:100:SER:OG	2.37	0.42
1:L:87:SER:OG	1:L:88:VAL:N	2.53	0.42
1:G:80:VAL:HG22	1:G:115:VAL:HB	2.01	0.42
1:J:17:ARG:NE	2:J:201:FMN:O1P	2.52	0.42
1:G:91:VAL:HG11	1:L:50:HIS:HB3	2.02	0.42
1:I:158:ASP:OD1	1:I:159:ALA:N	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:17:ARG:HD2	2:I:201:FMN:O1P	2.19	0.42
1:I:61:PRO:HB2	1:I:63:PRO:HD2	2.01	0.42
1:E:158:ASP:OD1	1:E:159:ALA:N	2.53	0.41
1:G:8:HIS:HE1	1:G:40:THR:CG2	2.32	0.41
1:J:159:ALA:C	1:J:161:THR:H	2.22	0.41
1:J:97:ASP:O	1:J:100:SER:OG	2.32	0.41
1:G:10:VAL:HB	1:G:42:LEU:HG	2.02	0.41
1:H:7:LEU:HD11	1:H:185:ARG:HD2	2.01	0.41
1:J:66:THR:O	1:J:70:GLN:HG3	2.20	0.41
1:G:54:ASP:O	1:G:58:GLU:HG3	2.20	0.41
1:L:118:SER:C	1:L:153:VAL:HG12	2.40	0.41
1:G:66:THR:O	1:G:70:GLN:HG3	2.20	0.41
1:H:139:ASP:OD1	1:J:144:ASN:HB2	2.20	0.41
1:I:116:THR:OG1	1:I:128:GLN:NE2	2.53	0.41
1:L:33:ALA:HA	1:L:34:PRO:HD3	1.91	0.41
1:E:121:MET:CG	1:L:121:MET:SD	3.09	0.41
1:H:103:SER:HA	1:H:104:PRO:HA	1.89	0.41
1:H:7:LEU:HD23	1:H:75:ASP:OD2	2.19	0.41
1:L:69:GLN:O	1:L:73:THR:N	2.34	0.41
1:D:143:LEU:HD11	1:D:179:ALA:HB3	2.02	0.41
1:E:10:VAL:HA	1:E:40:THR:O	2.21	0.41
1:F:33:ALA:HA	1:F:34:PRO:HD3	1.88	0.41
1:H:93:LYS:NZ	1:I:87:SER:HB3	2.36	0.41
1:L:109:GLY:HA2	1:L:141:TYR:CE2	2.56	0.41
1:D:33:ALA:HA	1:D:34:PRO:HD3	1.93	0.41
1:C:132:ARG:NH1	1:C:142:VAL:HG22	2.36	0.41
1:G:9:PHE:CE1	1:G:76:ALA:HB3	2.56	0.41
2:J:201:FMN:H1'2	2:J:201:FMN:H9	1.76	0.41
1:J:67:MET:O	1:J:71:ILE:HG13	2.20	0.41
1:L:51:TYR:HA	1:L:98:TRP:CZ2	2.55	0.41
1:D:10:VAL:HA	1:D:40:THR:O	2.21	0.41
1:E:33:ALA:HA	1:E:34:PRO:HD3	1.96	0.41
1:G:132:ARG:NH1	1:G:142:VAL:HG22	2.36	0.41
1:H:128:GLN:NE2	3:H:303:HOH:O	2.53	0.41
1:K:69:GLN:HG3	1:K:70:GLN:N	2.34	0.41
1:A:10:VAL:HB	1:A:42:LEU:HG	2.03	0.41
1:K:97:ASP:O	1:K:100:SER:OG	2.33	0.41
1:K:159:ALA:C	1:K:161:THR:H	2.20	0.40
1:L:31:GLU:HG3	1:L:170:ARG:NH2	2.36	0.40
1:J:60:PHE:HE1	1:J:102:VAL:HA	1.86	0.40
1:B:135:LEU:HA	1:B:135:LEU:HD23	1.88	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:48:PHE:N	1:H:48:PHE:CD2	2.89	0.40
1:H:48:PHE:HZ	1:H:67:MET:HB3	1.86	0.40
2:H:201:FMN:HM83	1:I:51:TYR:CE2	2.57	0.40
1:K:27:ARG:O	3:K:310:HOH:O	2.22	0.40
1:L:56:GLN:OE1	1:L:56:GLN:N	2.55	0.40
1:C:35:GLU:OE1	1:C:35:GLU:HA	2.21	0.40
1:D:9:PHE:CD1	1:D:76:ALA:HB3	2.57	0.40
1:F:32:ILE:HG23	1:F:174:ALA:HB2	2.02	0.40
1:G:116:THR:HG23	1:G:128:GLN:HE21	1.86	0.40
3:J:320:HOH:O	1:K:17:ARG:HB3	2.20	0.40
1:L:83:GLU:HB2	1:L:124:GLY:HA2	2.04	0.40
1:H:105:GLN:HA	1:H:106:PRO:HD3	1.93	0.40
1:I:45:ILE:HD12	1:I:91:VAL:HG22	2.04	0.40
1:J:80:VAL:HG22	1:J:115:VAL:CG1	2.51	0.40
1:J:15:SER:HB2	1:J:82:PRO:HG3	2.03	0.40
1:L:161:THR:O	1:L:162:LEU:HD23	2.22	0.40
1:L:17:ARG:NH2	1:L:21:PHE:CE1	2.89	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)
3:A:304:HOH:O	3:B:312:HOH:O[2_455]	2.12	0.08

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	182/199 (92%)	178 (98%)	3 (2%)	1 (0%)	29	39
1	B	182/199 (92%)	178 (98%)	3 (2%)	1 (0%)	29	39
1	C	182/199 (92%)	181 (100%)	1 (0%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	D	182/199 (92%)	176 (97%)	4 (2%)	2 (1%)	14	19
1	E	182/199 (92%)	178 (98%)	3 (2%)	1 (0%)	29	39
1	F	182/199 (92%)	178 (98%)	3 (2%)	1 (0%)	29	39
1	G	182/199 (92%)	176 (97%)	6 (3%)	0	100	100
1	H	182/199 (92%)	172 (94%)	7 (4%)	3 (2%)	9	12
1	I	182/199 (92%)	175 (96%)	6 (3%)	1 (0%)	29	39
1	J	182/199 (92%)	174 (96%)	4 (2%)	4 (2%)	6	7
1	K	182/199 (92%)	176 (97%)	5 (3%)	1 (0%)	29	39
1	L	182/199 (92%)	171 (94%)	9 (5%)	2 (1%)	14	19
All	All	2184/2388 (92%)	2113 (97%)	54 (2%)	17 (1%)	19	27

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	160	GLN
1	B	160	GLN
1	D	162	LEU
1	E	160	GLN
1	F	160	GLN
1	H	160	GLN
1	I	160	GLN
1	J	160	GLN
1	K	160	GLN
1	L	160	GLN
1	D	160	GLN
1	J	29	LEU
1	J	184	ALA
1	H	162	LEU
1	H	178	ASP
1	J	34	PRO
1	L	55	VAL

5.3.2 Protein sidechains ⓘ

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	144/158 (91%)	130 (90%)	14 (10%)	8	9
1	B	144/158 (91%)	136 (94%)	8 (6%)	21	27
1	C	144/158 (91%)	133 (92%)	11 (8%)	13	17
1	D	144/158 (91%)	136 (94%)	8 (6%)	21	27
1	E	144/158 (91%)	133 (92%)	11 (8%)	13	17
1	F	144/158 (91%)	134 (93%)	10 (7%)	15	19
1	G	144/158 (91%)	133 (92%)	11 (8%)	13	17
1	H	144/158 (91%)	130 (90%)	14 (10%)	8	9
1	I	144/158 (91%)	127 (88%)	17 (12%)	5	5
1	J	144/158 (91%)	130 (90%)	14 (10%)	8	9
1	K	144/158 (91%)	133 (92%)	11 (8%)	13	17
1	L	144/158 (91%)	128 (89%)	16 (11%)	6	6
All	All	1728/1896 (91%)	1583 (92%)	145 (8%)	11	13

All (145) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ILE
1	A	37	ILE
1	A	88	VAL
1	A	103	SER
1	A	105	GLN
1	A	121	MET
1	A	122	ILE
1	A	143	LEU
1	A	149	MET
1	A	154	THR
1	A	157	VAL
1	A	158	ASP
1	A	175	ARG
1	A	187	LEU
1	B	37	ILE
1	B	69	GLN
1	B	88	VAL
1	B	105	GLN
1	B	143	LEU
1	B	154	THR

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Mol	Chain	Res	Type
1	B	161	THR
1	B	187	LEU
1	C	37	ILE
1	C	88	VAL
1	C	91	VAL
1	C	105	GLN
1	C	142	VAL
1	C	143	LEU
1	C	149	MET
1	C	154	THR
1	C	157	VAL
1	C	161	THR
1	C	187	LEU
1	D	32	ILE
1	D	91	VAL
1	D	105	GLN
1	D	142	VAL
1	D	143	LEU
1	D	154	THR
1	D	161	THR
1	D	187	LEU
1	E	25	VAL
1	E	57	GLU
1	E	88	VAL
1	E	105	GLN
1	E	121	MET
1	E	142	VAL
1	E	143	LEU
1	E	145	ARG
1	E	149	MET
1	E	154	THR
1	E	157	VAL
1	F	88	VAL
1	F	105	GLN
1	F	121	MET
1	F	134	SER
1	F	142	VAL
1	F	143	LEU
1	F	149	MET
1	F	154	THR
1	F	157	VAL
1	F	187	LEU

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Mol	Chain	Res	Type
1	G	37	ILE
1	G	105	GLN
1	G	121	MET
1	G	142	VAL
1	G	143	LEU
1	G	150	ILE
1	G	154	THR
1	G	161	THR
1	G	167	VAL
1	G	175	ARG
1	G	187	LEU
1	H	37	ILE
1	H	60	PHE
1	H	75	ASP
1	H	78	VAL
1	H	88	VAL
1	H	118	SER
1	H	121	MET
1	H	143	LEU
1	H	149	MET
1	H	154	THR
1	H	157	VAL
1	H	161	THR
1	H	162	LEU
1	H	187	LEU
1	I	20	SER
1	I	32	ILE
1	I	37	ILE
1	I	53	GLN
1	I	88	VAL
1	I	105	GLN
1	I	116	THR
1	I	121	MET
1	I	131	LEU
1	I	143	LEU
1	I	154	THR
1	I	161	THR
1	I	162	LEU
1	I	165	SER
1	I	167	VAL
1	I	178	ASP
1	I	187	LEU

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Mol	Chain	Res	Type
1	J	10	VAL
1	J	37	ILE
1	J	40	THR
1	J	56	GLN
1	J	69	GLN
1	J	105	GLN
1	J	121	MET
1	J	142	VAL
1	J	156	LYS
1	J	157	VAL
1	J	158	ASP
1	J	163	GLU
1	J	172	PHE
1	J	187	LEU
1	K	10	VAL
1	K	12	LEU
1	K	27	ARG
1	K	32	ILE
1	K	55	VAL
1	K	69	GLN
1	K	143	LEU
1	K	150	ILE
1	K	157	VAL
1	K	175	ARG
1	K	187	LEU
1	L	5	SER
1	L	32	ILE
1	L	37	ILE
1	L	40	THR
1	L	42	LEU
1	L	56	GLN
1	L	60	PHE
1	L	63	PRO
1	L	102	VAL
1	L	143	LEU
1	L	154	THR
1	L	161	THR
1	L	164	LEU
1	L	167	VAL
1	L	175	ARG
1	L	187	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	105	GLN
1	B	69	GLN
1	B	152	GLN
1	C	50	HIS
1	D	50	HIS
1	E	152	GLN
1	G	8	HIS
1	G	50	HIS
1	G	128	GLN
1	H	8	HIS
1	I	105	GLN
1	I	128	GLN
1	J	50	HIS
1	J	56	GLN
1	J	69	GLN
1	J	152	GLN
1	K	8	HIS
1	K	50	HIS
1	L	50	HIS
1	L	105	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

12 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	K	201	-	31,33,33	1.38	3 (9%)	40,50,50	1.63	7 (17%)
2	FMN	I	201	-	31,33,33	1.42	4 (12%)	40,50,50	1.43	5 (12%)
2	FMN	C	201	-	31,33,33	1.33	3 (9%)	40,50,50	1.70	6 (15%)
2	FMN	A	201	-	31,33,33	1.31	5 (16%)	40,50,50	1.61	7 (17%)
2	FMN	G	201	-	31,33,33	1.47	3 (9%)	40,50,50	1.66	5 (12%)
2	FMN	E	201	-	31,33,33	1.35	4 (12%)	40,50,50	1.50	5 (12%)
2	FMN	J	201	-	31,33,33	1.39	4 (12%)	40,50,50	1.55	6 (15%)
2	FMN	H	201	-	31,33,33	1.55	4 (12%)	40,50,50	1.84	6 (15%)
2	FMN	L	201	-	31,33,33	1.44	4 (12%)	40,50,50	1.54	5 (12%)
2	FMN	B	201	-	31,33,33	1.31	4 (12%)	40,50,50	1.61	7 (17%)
2	FMN	F	201	-	31,33,33	1.39	3 (9%)	40,50,50	1.73	9 (22%)
2	FMN	D	201	-	31,33,33	1.37	4 (12%)	40,50,50	1.66	6 (15%)

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	K	201	-	-	7/18/18/18	0/3/3/3
2	FMN	I	201	-	-	3/18/18/18	0/3/3/3
2	FMN	C	201	-	-	1/18/18/18	0/3/3/3
2	FMN	A	201	-	-	1/18/18/18	0/3/3/3
2	FMN	G	201	-	-	14/18/18/18	0/3/3/3
2	FMN	E	201	-	-	2/18/18/18	0/3/3/3
2	FMN	J	201	-	-	6/18/18/18	0/3/3/3
2	FMN	H	201	-	-	15/18/18/18	0/3/3/3
2	FMN	L	201	-	-	4/18/18/18	0/3/3/3
2	FMN	B	201	-	-	1/18/18/18	0/3/3/3
2	FMN	F	201	-	-	1/18/18/18	0/3/3/3
2	FMN	D	201	-	-	1/18/18/18	0/3/3/3

All (45) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	201	FMN	C10-N1	4.50	1.39	1.33
2	G	201	FMN	C10-N1	4.43	1.38	1.33
2	H	201	FMN	C10-N1	4.42	1.38	1.33
2	I	201	FMN	C10-N1	4.30	1.38	1.33
2	L	201	FMN	C10-N1	4.30	1.38	1.33
2	J	201	FMN	C10-N1	4.25	1.38	1.33
2	K	201	FMN	C10-N1	4.02	1.38	1.33
2	H	201	FMN	C1'-N10	4.00	1.52	1.48
2	E	201	FMN	C10-N1	3.88	1.38	1.33
2	B	201	FMN	C10-N1	3.88	1.38	1.33
2	L	201	FMN	C4A-N5	3.73	1.38	1.33
2	C	201	FMN	C10-N1	3.68	1.38	1.33
2	G	201	FMN	C4A-N5	3.67	1.38	1.33
2	K	201	FMN	C4A-N5	3.63	1.38	1.33
2	D	201	FMN	C4-N3	3.53	1.39	1.33
2	D	201	FMN	C10-N1	3.50	1.37	1.33
2	A	201	FMN	C10-N1	3.50	1.37	1.33
2	D	201	FMN	C4A-N5	3.48	1.38	1.33
2	I	201	FMN	C4A-N5	3.42	1.38	1.33
2	C	201	FMN	C4A-N5	3.40	1.38	1.33
2	H	201	FMN	C4A-N5	3.28	1.38	1.33
2	F	201	FMN	C4A-N5	3.26	1.38	1.33
2	E	201	FMN	C4-N3	3.22	1.38	1.33
2	G	201	FMN	C4-N3	3.21	1.38	1.33
2	I	201	FMN	C4-N3	3.14	1.38	1.33
2	A	201	FMN	C4A-N5	3.13	1.37	1.33
2	L	201	FMN	C4-N3	3.12	1.38	1.33
2	K	201	FMN	C4-N3	3.11	1.38	1.33
2	J	201	FMN	C4A-N5	3.09	1.37	1.33
2	B	201	FMN	C4A-N5	2.95	1.37	1.33
2	E	201	FMN	C4A-N5	2.93	1.37	1.33
2	C	201	FMN	C4-N3	2.92	1.38	1.33
2	A	201	FMN	C4-N3	2.89	1.38	1.33
2	H	201	FMN	C4-N3	2.88	1.38	1.33
2	J	201	FMN	C4-N3	2.76	1.37	1.33
2	J	201	FMN	C1'-N10	2.71	1.51	1.48
2	F	201	FMN	C4-N3	2.70	1.37	1.33
2	B	201	FMN	C1'-N10	2.46	1.50	1.48
2	D	201	FMN	C1'-N10	2.38	1.50	1.48
2	I	201	FMN	C1'-N10	2.33	1.50	1.48
2	A	201	FMN	C1'-N10	2.31	1.50	1.48
2	B	201	FMN	C4-N3	2.27	1.37	1.33
2	A	201	FMN	C9A-C5A	-2.08	1.38	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	201	FMN	C5A-N5	2.00	1.38	1.35
2	E	201	FMN	C1'-N10	2.00	1.50	1.48

All (74) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	201	FMN	C4-N3-C2	6.33	120.49	115.14
2	C	201	FMN	C4-N3-C2	6.16	120.35	115.14
2	K	201	FMN	C4-N3-C2	6.08	120.28	115.14
2	H	201	FMN	C1'-N10-C9A	5.96	122.98	118.29
2	D	201	FMN	C4-N3-C2	5.92	120.14	115.14
2	L	201	FMN	C4-N3-C2	5.81	120.04	115.14
2	G	201	FMN	C4-N3-C2	5.75	120.00	115.14
2	E	201	FMN	C4-N3-C2	5.74	119.98	115.14
2	H	201	FMN	C4-N3-C2	5.60	119.87	115.14
2	I	201	FMN	C4-N3-C2	5.57	119.85	115.14
2	J	201	FMN	C4-N3-C2	5.04	119.40	115.14
2	B	201	FMN	C4-N3-C2	4.78	119.18	115.14
2	A	201	FMN	C4-N3-C2	4.77	119.17	115.14
2	F	201	FMN	C4A-N5-C5A	4.71	121.48	116.77
2	G	201	FMN	C1'-N10-C10	4.65	122.57	118.41
2	B	201	FMN	C4A-N5-C5A	4.55	121.32	116.77
2	C	201	FMN	C4A-N5-C5A	4.33	121.09	116.77
2	L	201	FMN	C4A-N5-C5A	4.16	120.92	116.77
2	K	201	FMN	C4A-N5-C5A	4.10	120.87	116.77
2	H	201	FMN	C5A-C9A-N10	4.03	120.63	117.72
2	A	201	FMN	C4A-N5-C5A	4.01	120.78	116.77
2	D	201	FMN	C4A-N5-C5A	3.87	120.64	116.77
2	A	201	FMN	C1'-N10-C10	3.78	121.79	118.41
2	J	201	FMN	C4A-N5-C5A	3.61	120.38	116.77
2	J	201	FMN	C1'-N10-C10	3.43	121.48	118.41
2	D	201	FMN	C4A-C4-N3	-3.30	118.92	123.43
2	C	201	FMN	C4A-C4-N3	-3.28	118.94	123.43
2	H	201	FMN	C4A-N5-C5A	3.28	120.05	116.77
2	J	201	FMN	C5A-C9A-N10	3.22	120.05	117.72
2	E	201	FMN	C4A-C4-N3	-3.14	119.13	123.43
2	C	201	FMN	C1'-N10-C10	3.13	121.21	118.41
2	G	201	FMN	C4A-N5-C5A	3.11	119.88	116.77
2	E	201	FMN	C5A-C9A-N10	3.10	119.96	117.72
2	L	201	FMN	C1'-N10-C9A	3.09	120.73	118.29
2	I	201	FMN	C5A-C9A-N10	3.06	119.93	117.72
2	K	201	FMN	C1'-N10-C9A	3.03	120.68	118.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	201	FMN	C5A-C9A-N10	3.03	119.91	117.72
2	F	201	FMN	C5A-C9A-N10	3.01	119.90	117.72
2	E	201	FMN	C4A-N5-C5A	3.00	119.77	116.77
2	G	201	FMN	C5A-C9A-N10	2.99	119.88	117.72
2	I	201	FMN	C4A-N5-C5A	2.99	119.76	116.77
2	B	201	FMN	C5A-C9A-N10	2.99	119.88	117.72
2	C	201	FMN	C5A-C9A-N10	2.95	119.85	117.72
2	A	201	FMN	C10-C4A-N5	-2.93	119.23	121.26
2	I	201	FMN	C1'-N10-C9A	2.92	120.59	118.29
2	D	201	FMN	C10-C4A-N5	-2.89	119.26	121.26
2	K	201	FMN	C5A-C9A-N10	2.88	119.81	117.72
2	D	201	FMN	C1'-N10-C9A	2.84	120.53	118.29
2	A	201	FMN	C5A-C9A-N10	2.83	119.77	117.72
2	E	201	FMN	C1'-N10-C10	2.83	120.94	118.41
2	A	201	FMN	C4A-C4-N3	-2.79	119.61	123.43
2	F	201	FMN	C4A-C4-N3	-2.79	119.62	123.43
2	H	201	FMN	C9A-N10-C10	-2.68	118.40	121.91
2	B	201	FMN	C10-C4A-N5	-2.65	119.42	121.26
2	K	201	FMN	C4A-C4-N3	-2.57	119.92	123.43
2	J	201	FMN	O3P-P-O5'	2.53	113.47	106.73
2	B	201	FMN	C1'-N10-C10	2.50	120.65	118.41
2	C	201	FMN	C10-C4A-N5	-2.49	119.53	121.26
2	G	201	FMN	C4A-C4-N3	-2.48	120.04	123.43
2	B	201	FMN	C4A-C4-N3	-2.45	120.08	123.43
2	K	201	FMN	C10-C4A-N5	-2.45	119.56	121.26
2	H	201	FMN	C4A-C4-N3	-2.42	120.12	123.43
2	F	201	FMN	C1'-N10-C10	2.41	120.57	118.41
2	F	201	FMN	C4'-C3'-C2'	-2.41	108.35	113.36
2	L	201	FMN	C4A-C4-N3	-2.40	120.15	123.43
2	L	201	FMN	C10-C4A-N5	-2.28	119.68	121.26
2	I	201	FMN	C4A-C4-N3	-2.27	120.33	123.43
2	J	201	FMN	C4A-C4-N3	-2.26	120.34	123.43
2	F	201	FMN	C10-C4A-N5	-2.11	119.80	121.26
2	A	201	FMN	C5'-C4'-C3'	-2.09	108.17	112.20
2	F	201	FMN	C9A-C5A-N5	-2.09	119.10	122.36
2	K	201	FMN	C4-C4A-N5	2.07	120.96	118.60
2	B	201	FMN	C5'-C4'-C3'	-2.01	108.32	112.20
2	F	201	FMN	C5'-C4'-C3'	-2.00	108.34	112.20

There are no chirality outliers.

All (56) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	K	201	FMN	N10-C1'-C2'-O2'
2	I	201	FMN	N10-C1'-C2'-O2'
2	I	201	FMN	N10-C1'-C2'-C3'
2	G	201	FMN	N10-C1'-C2'-O2'
2	G	201	FMN	N10-C1'-C2'-C3'
2	G	201	FMN	C1'-C2'-C3'-O3'
2	G	201	FMN	C1'-C2'-C3'-C4'
2	G	201	FMN	O2'-C2'-C3'-C4'
2	G	201	FMN	C2'-C3'-C4'-O4'
2	G	201	FMN	C5'-O5'-P-O1P
2	G	201	FMN	C5'-O5'-P-O2P
2	G	201	FMN	C5'-O5'-P-O3P
2	J	201	FMN	O4'-C4'-C5'-O5'
2	J	201	FMN	C5'-O5'-P-O2P
2	J	201	FMN	C5'-O5'-P-O3P
2	H	201	FMN	C2'-C1'-N10-C9A
2	H	201	FMN	C2'-C1'-N10-C10
2	H	201	FMN	N10-C1'-C2'-O2'
2	H	201	FMN	N10-C1'-C2'-C3'
2	H	201	FMN	C1'-C2'-C3'-O3'
2	H	201	FMN	C1'-C2'-C3'-C4'
2	H	201	FMN	O2'-C2'-C3'-O3'
2	H	201	FMN	O2'-C2'-C3'-C4'
2	H	201	FMN	C2'-C3'-C4'-C5'
2	H	201	FMN	O3'-C3'-C4'-O4'
2	H	201	FMN	O3'-C3'-C4'-C5'
2	H	201	FMN	C3'-C4'-C5'-O5'
2	H	201	FMN	O4'-C4'-C5'-O5'
2	L	201	FMN	N10-C1'-C2'-O2'
2	L	201	FMN	C5'-O5'-P-O1P
2	G	201	FMN	O2'-C2'-C3'-O3'
2	G	201	FMN	O3'-C3'-C4'-O4'
2	H	201	FMN	C2'-C3'-C4'-O4'
2	G	201	FMN	O3'-C3'-C4'-C5'
2	G	201	FMN	C2'-C3'-C4'-C5'
2	K	201	FMN	O3'-C3'-C4'-O4'
2	K	201	FMN	O3'-C3'-C4'-C5'
2	K	201	FMN	C2'-C3'-C4'-C5'
2	K	201	FMN	C2'-C3'-C4'-O4'
2	J	201	FMN	C3'-C4'-C5'-O5'
2	J	201	FMN	C5'-O5'-P-O1P
2	A	201	FMN	C4'-C5'-O5'-P
2	G	201	FMN	C4'-C5'-O5'-P

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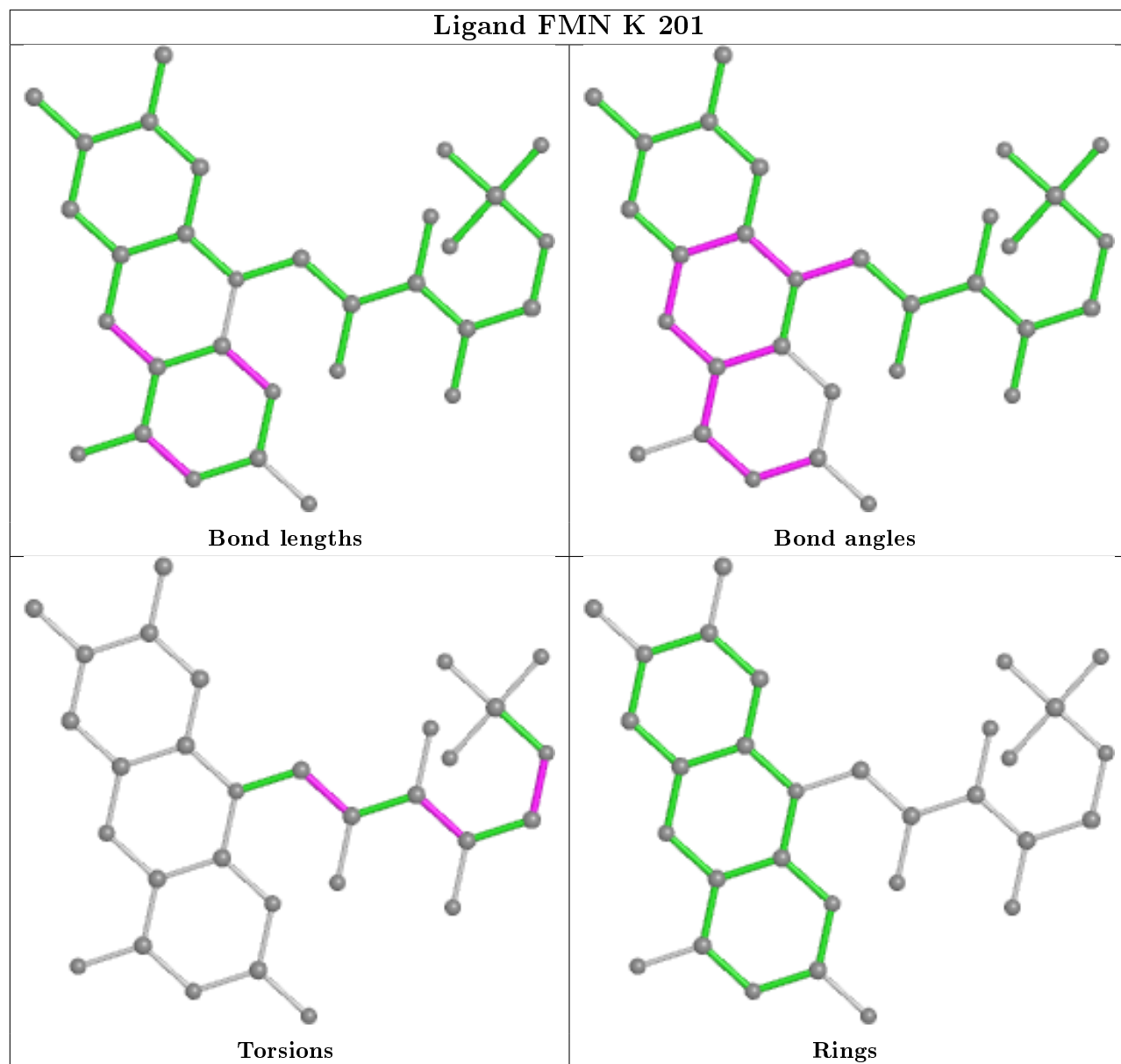
Mol	Chain	Res	Type	Atoms
2	D	201	FMN	C4'-C5'-O5'-P
2	C	201	FMN	C4'-C5'-O5'-P
2	F	201	FMN	C4'-C5'-O5'-P
2	K	201	FMN	C4'-C5'-O5'-P
2	K	201	FMN	N10-C1'-C2'-C3'
2	L	201	FMN	N10-C1'-C2'-C3'
2	E	201	FMN	C4'-C5'-O5'-P
2	E	201	FMN	C5'-O5'-P-O1P
2	J	201	FMN	C4'-C5'-O5'-P
2	B	201	FMN	C4'-C5'-O5'-P
2	L	201	FMN	C5'-O5'-P-O3P
2	H	201	FMN	C4'-C5'-O5'-P
2	I	201	FMN	O4'-C4'-C5'-O5'

There are no ring outliers.

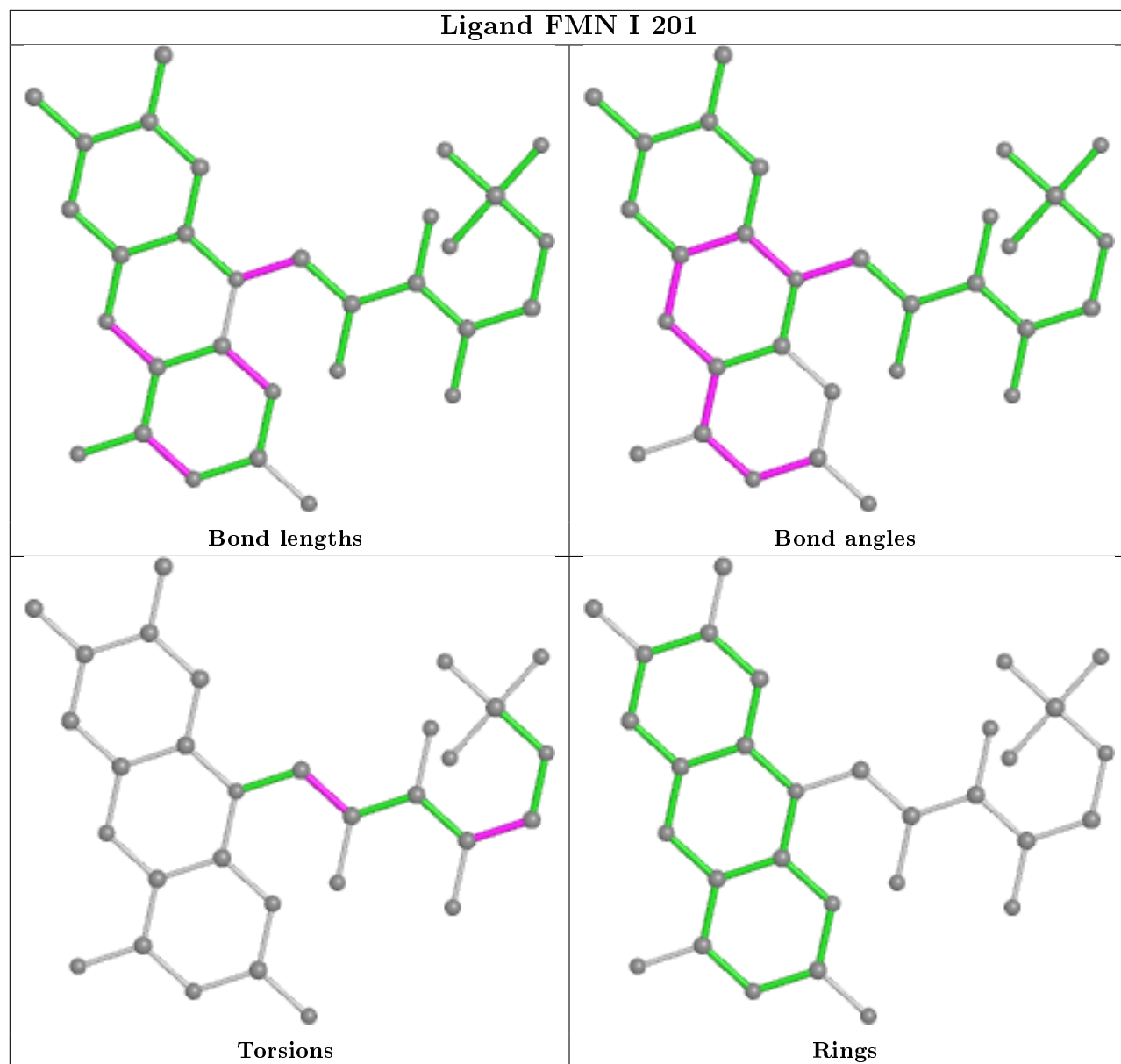
7 monomers are involved in 26 short contacts:

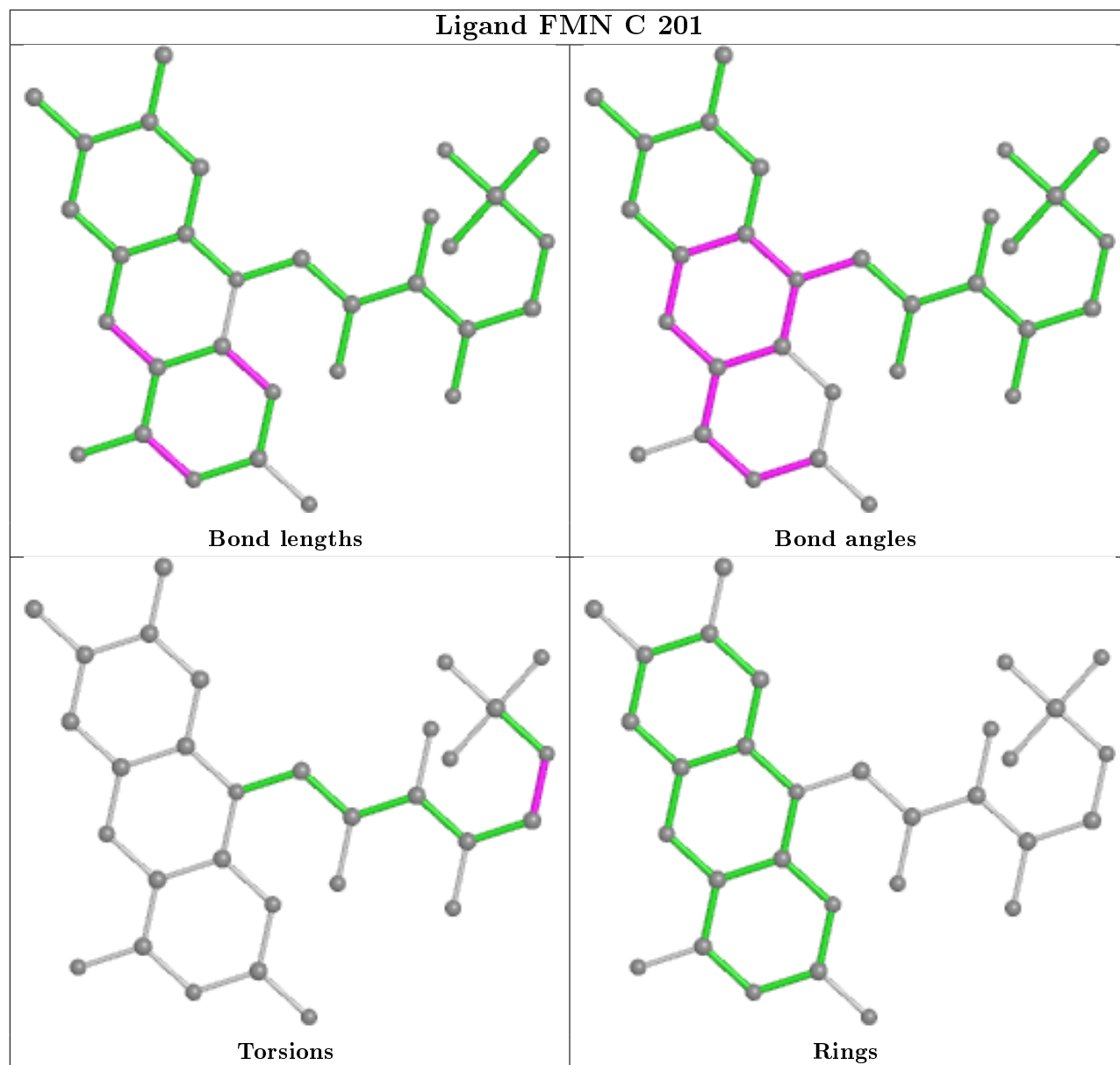
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	201	FMN	1	0
2	I	201	FMN	6	0
2	G	201	FMN	4	0
2	E	201	FMN	1	0
2	J	201	FMN	6	0
2	H	201	FMN	5	0
2	L	201	FMN	3	0

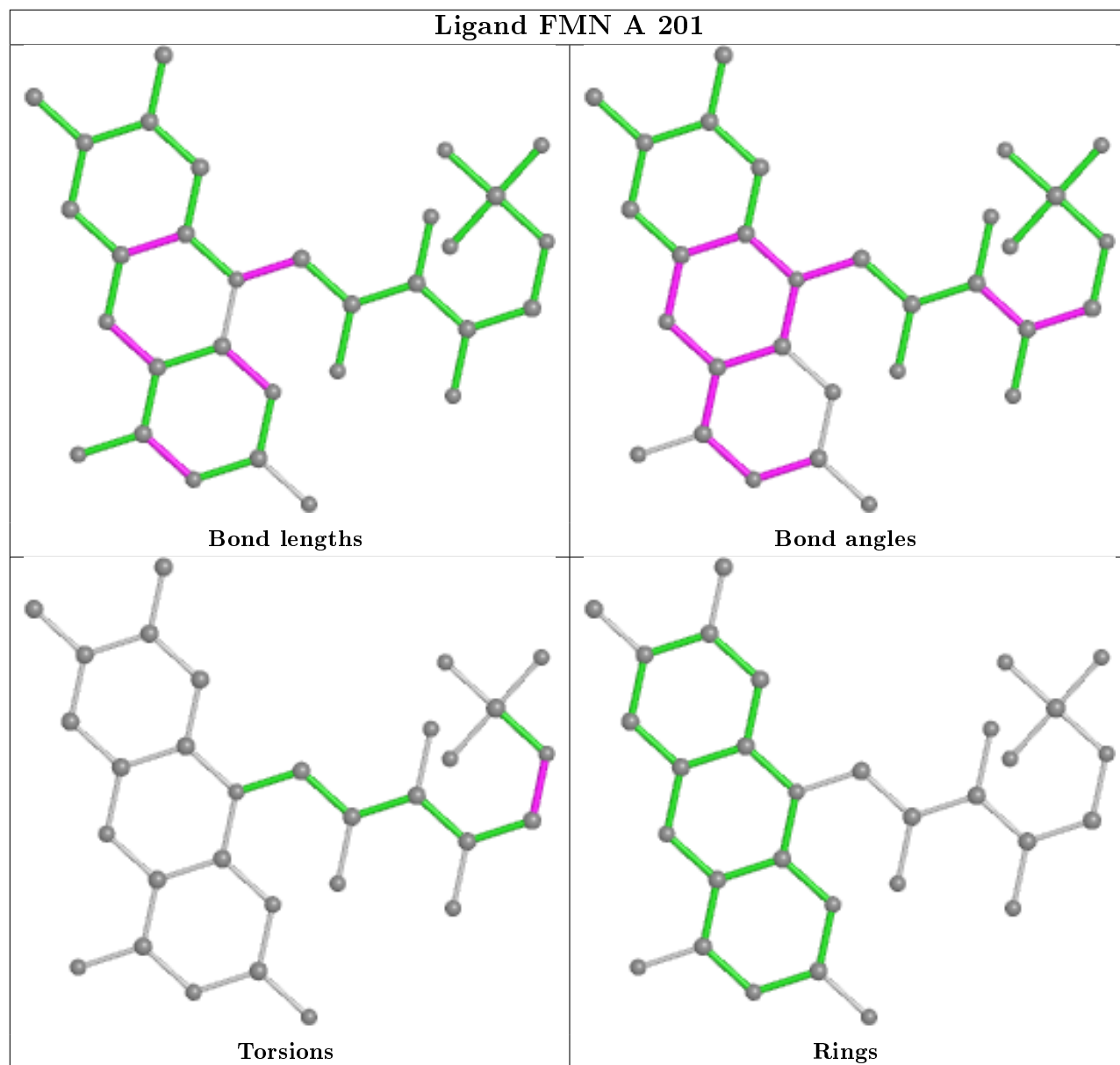
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

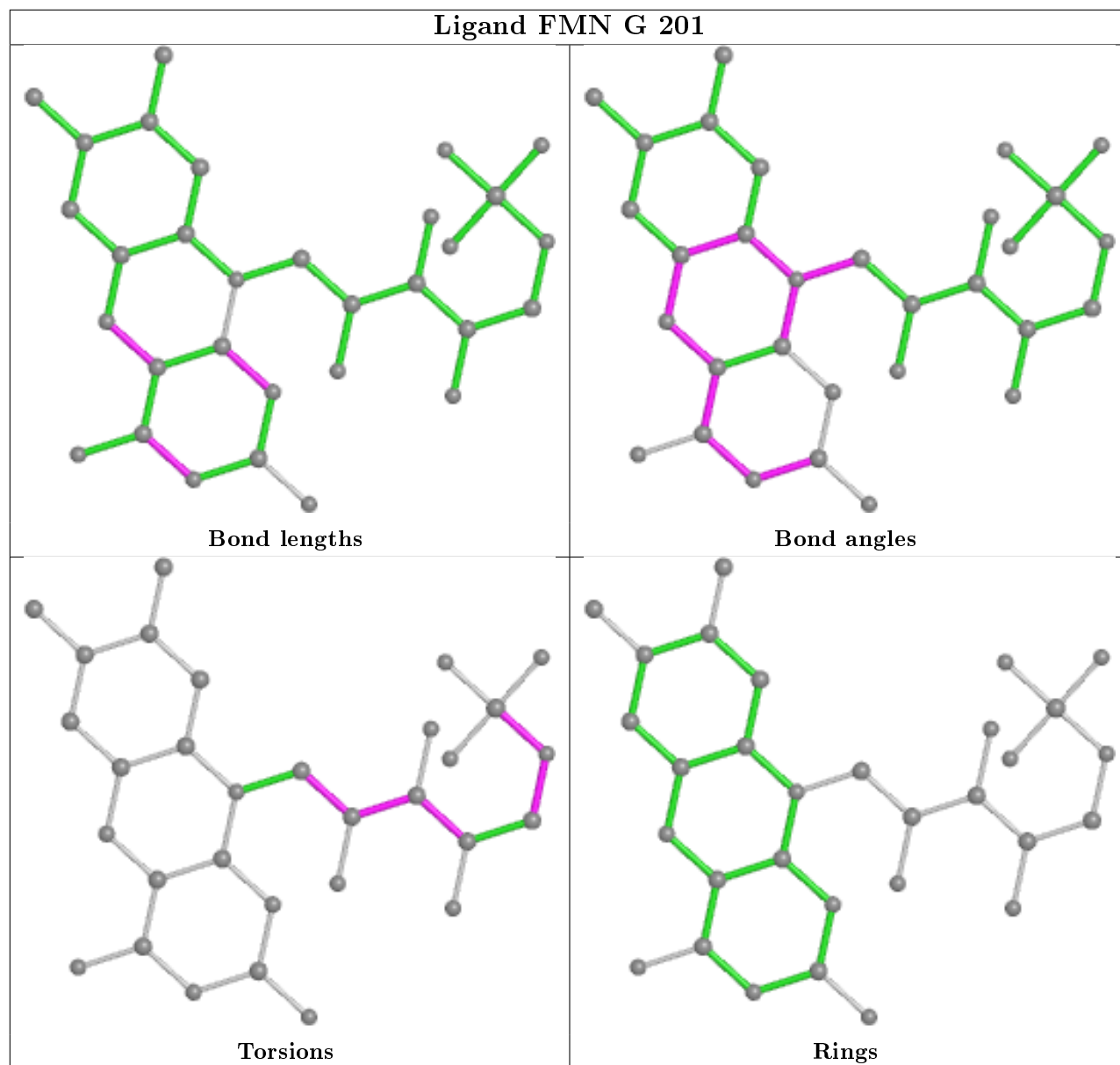


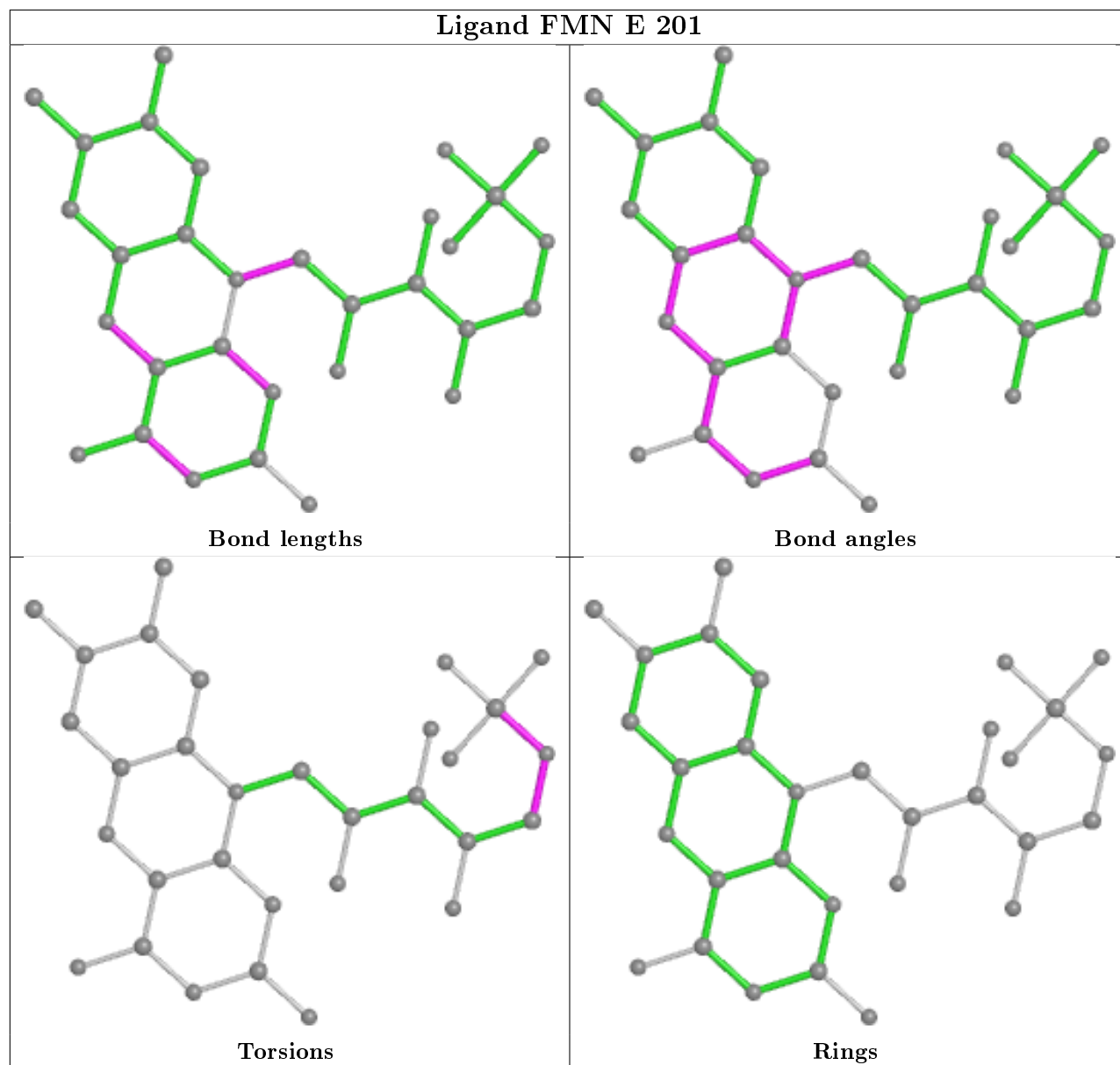
Ligand FMN I 201

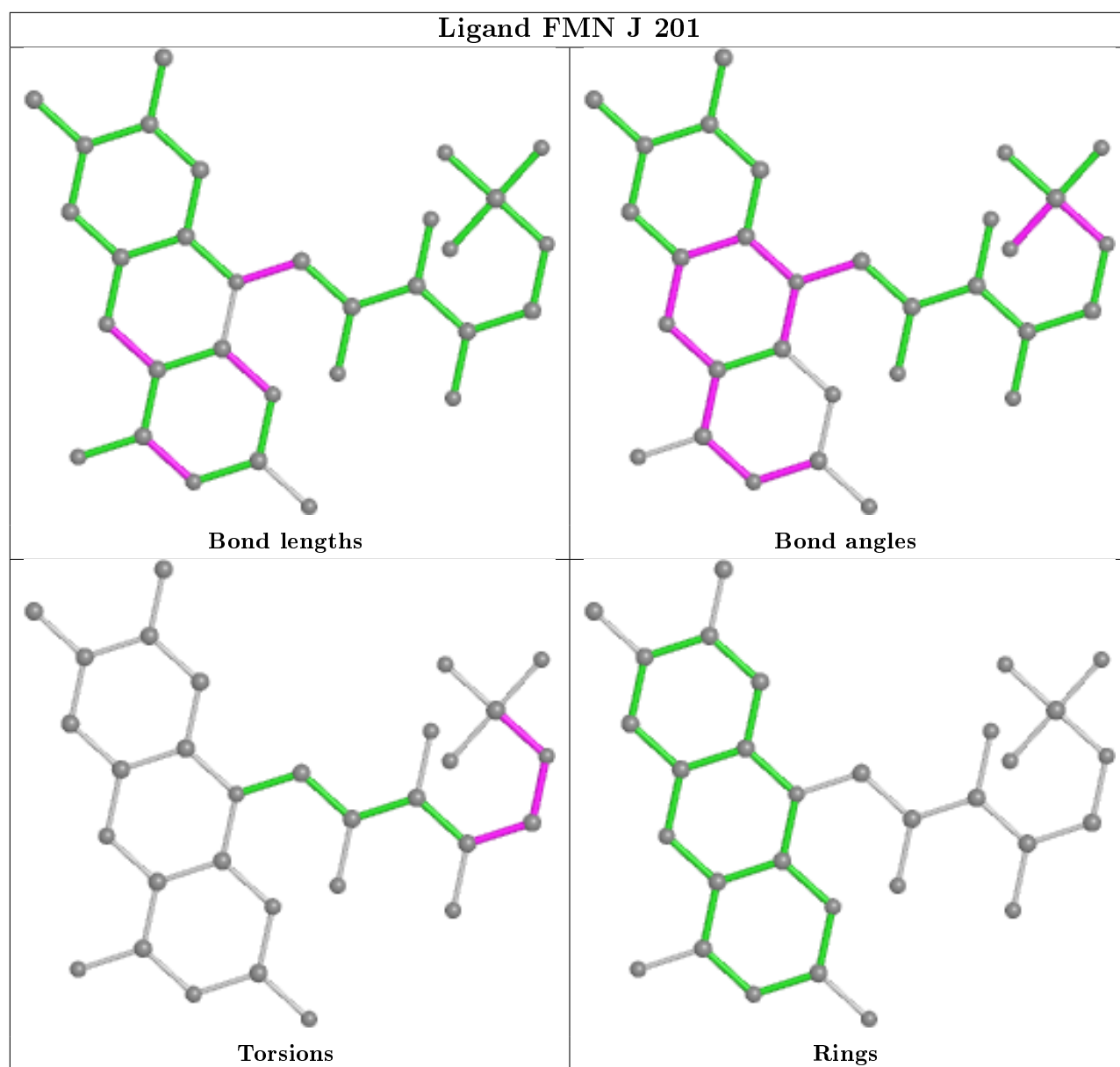


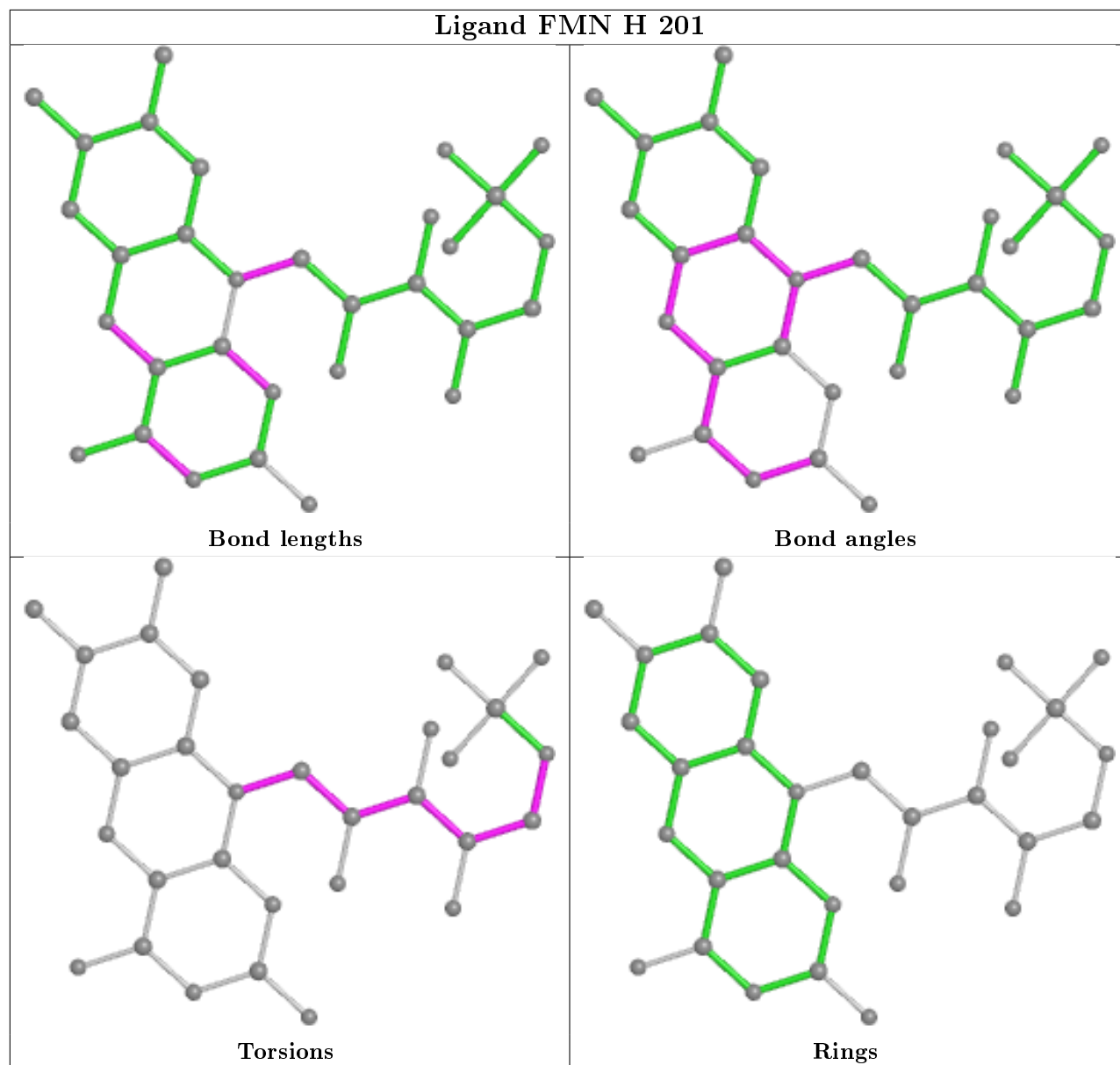


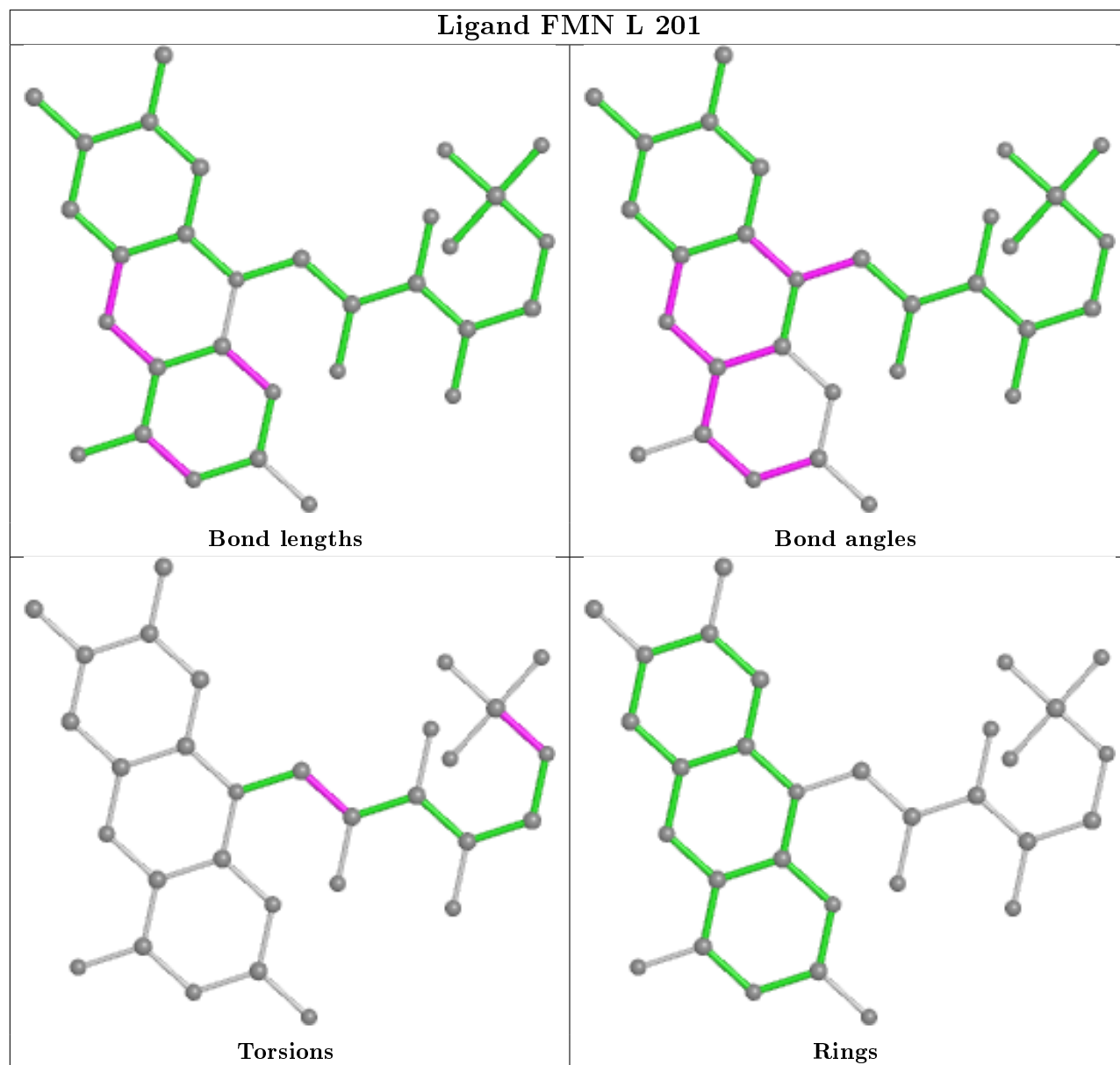


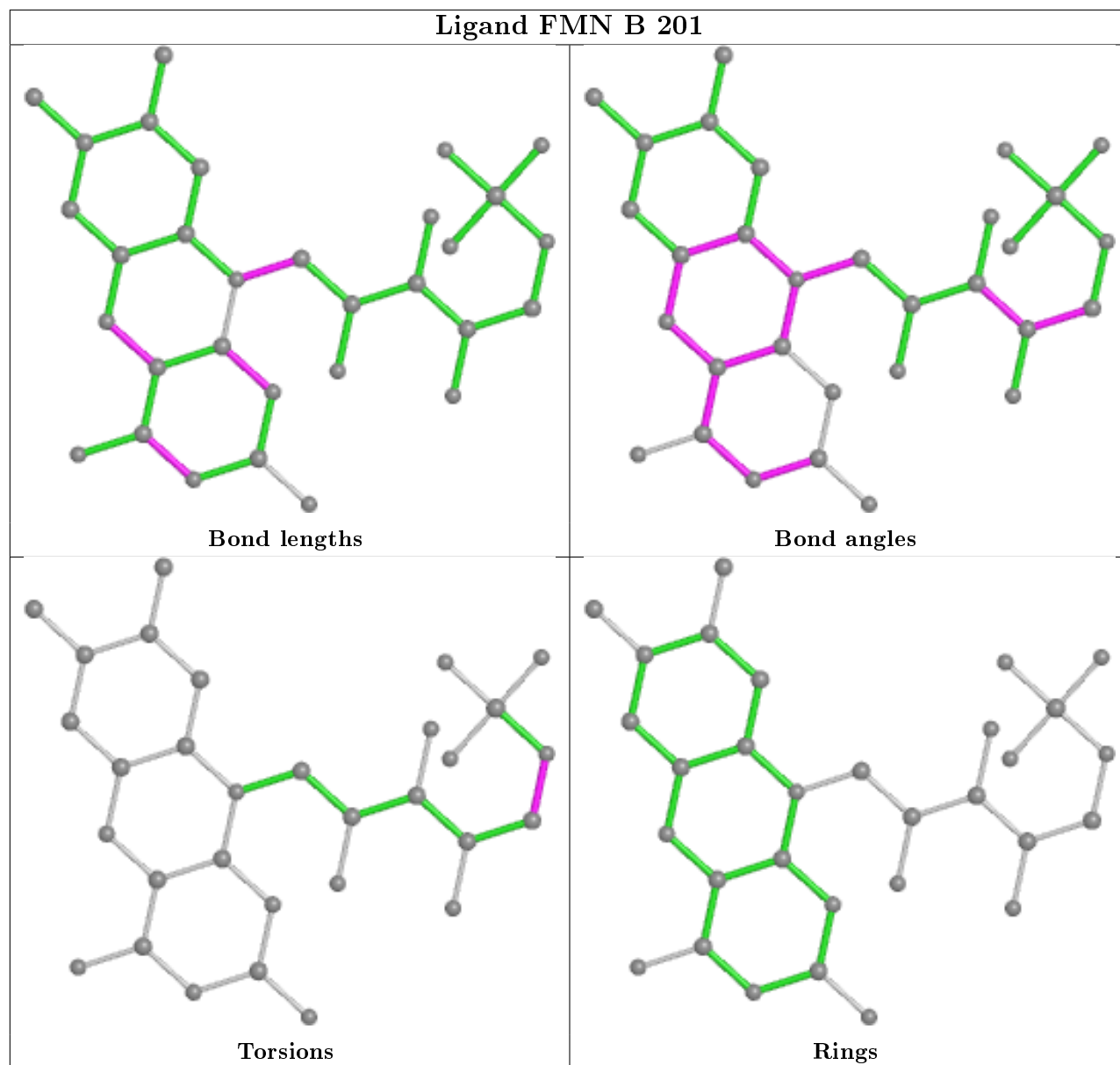


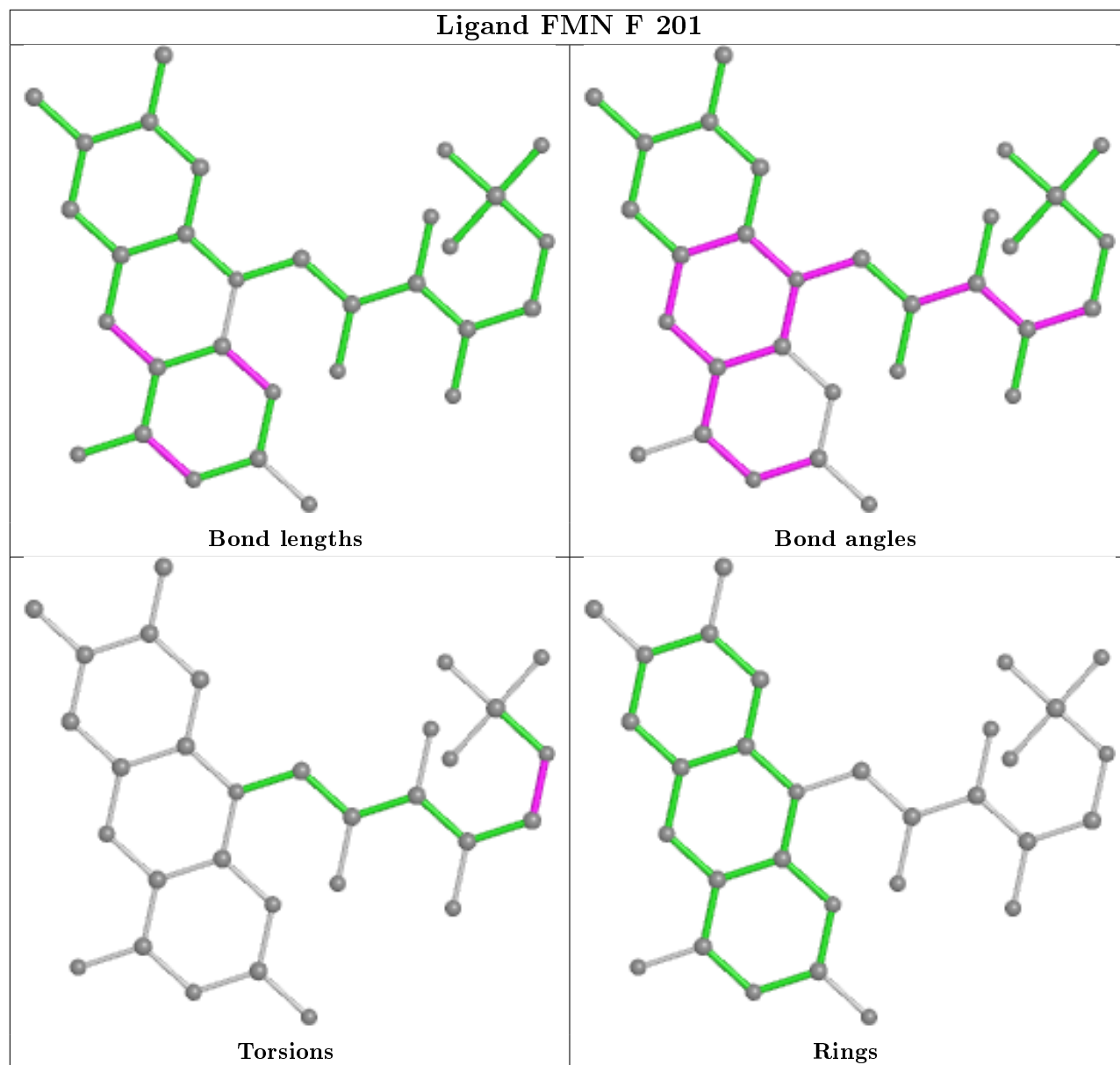


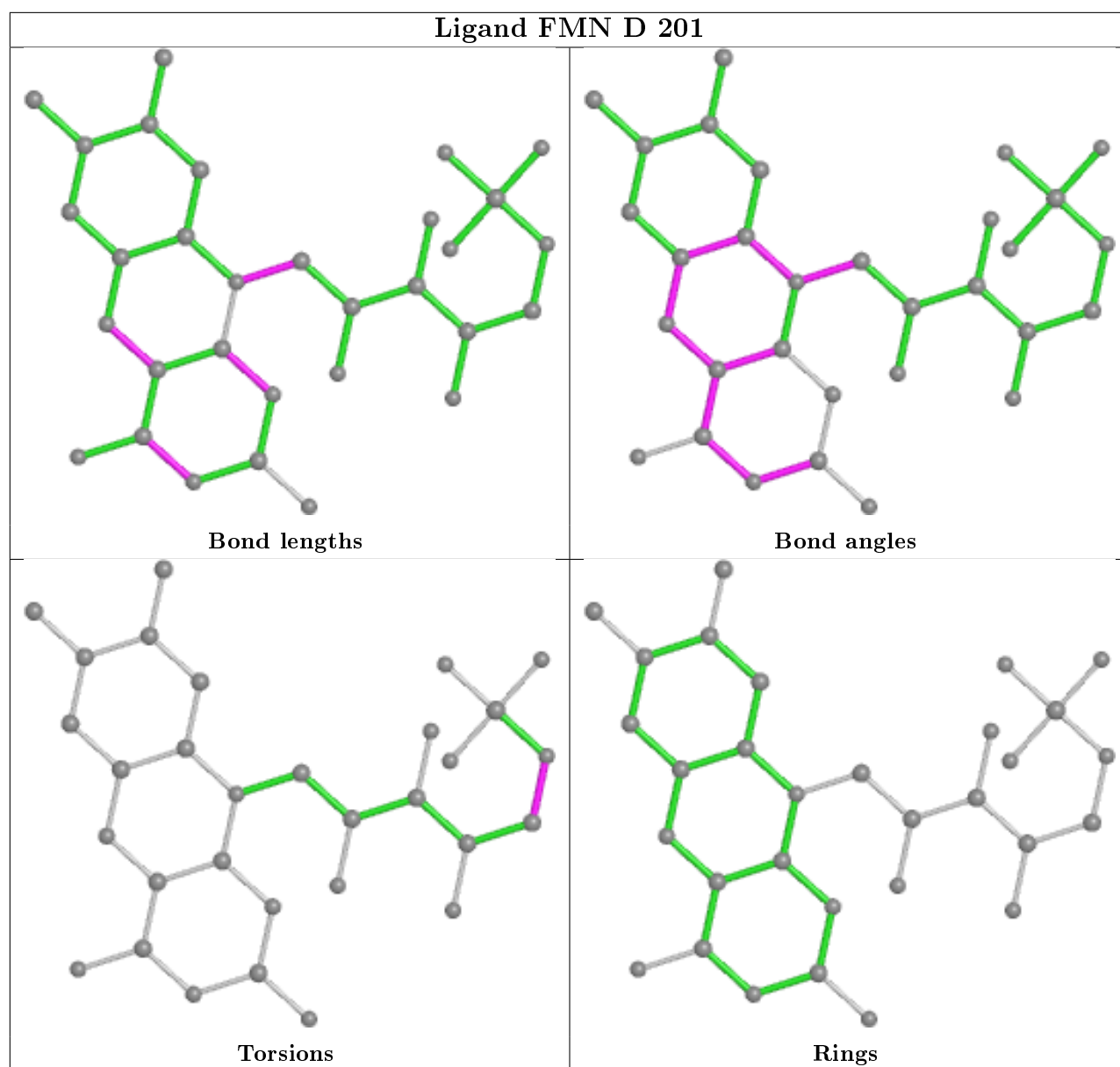












5.7 Other polymers [i](#)

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, 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	184/199 (92%)	-0.28	2 (1%) 80 85	17, 30, 52, 75	0
1	B	184/199 (92%)	-0.35	0 100 100	16, 29, 53, 75	0
1	C	184/199 (92%)	-0.29	1 (0%) 91 94	21, 33, 57, 74	0
1	D	184/199 (92%)	-0.20	4 (2%) 62 70	21, 36, 68, 82	0
1	E	184/199 (92%)	-0.18	2 (1%) 80 85	21, 38, 64, 82	0
1	F	184/199 (92%)	0.08	3 (1%) 72 78	26, 46, 67, 87	0
1	G	184/199 (92%)	0.59	15 (8%) 11 16	36, 72, 94, 102	0
1	H	184/199 (92%)	1.19	35 (19%) 1 1	42, 76, 105, 110	0
1	I	184/199 (92%)	1.17	36 (19%) 1 1	51, 81, 102, 109	0
1	J	184/199 (92%)	1.27	41 (22%) 0 1	52, 82, 104, 119	0
1	K	184/199 (92%)	0.72	24 (13%) 3 4	39, 77, 98, 108	0
1	L	184/199 (92%)	0.74	18 (9%) 7 10	40, 79, 102, 112	0
All	All	2208/2388 (92%)	0.37	181 (8%) 11 16	16, 56, 99, 119	0

All (181) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	36	GLY	8.0
1	L	32	ILE	7.2
1	I	36	GLY	6.6
1	J	35	GLU	6.6
1	J	37	ILE	6.5
1	J	183	LEU	6.1
1	I	16	LEU	6.1
1	H	8	HIS	6.0
1	A	5	SER	5.6
1	J	6	PRO	5.5
1	H	76	ALA	5.3

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Mol	Chain	Res	Type	RSRZ
1	I	15	SER	5.2
1	I	25	VAL	5.2
1	J	169	THR	5.0
1	J	174	ALA	4.9
1	H	77	VAL	4.8
1	K	60	PHE	4.8
1	J	181	ALA	4.8
1	I	6	PRO	4.7
1	G	35	GLU	4.7
1	J	7	LEU	4.6
1	L	121	MET	4.4
1	G	5	SER	4.4
1	J	76	ALA	4.3
1	K	76	ALA	4.3
1	H	49	PRO	4.2
1	J	16	LEU	4.1
1	H	108	ALA	4.1
1	K	52	SER	4.0
1	I	80	VAL	4.0
1	H	140	ALA	4.0
1	J	41	PRO	3.9
1	D	5	SER	3.8
1	H	68	ALA	3.8
1	K	74	ALA	3.7
1	I	174	ALA	3.7
1	I	33	ALA	3.7
1	G	37	ILE	3.6
1	K	98	TRP	3.6
1	G	16	LEU	3.6
1	F	5	SER	3.5
1	J	80	VAL	3.5
1	I	5	SER	3.5
1	I	126	ARG	3.5
1	I	22	ASN	3.5
1	J	77	VAL	3.4
1	H	10	VAL	3.4
1	H	75	ASP	3.4
1	H	109	GLY	3.3
1	G	7	LEU	3.3
1	D	160	GLN	3.3
1	I	18	LYS	3.3
1	J	176	GLN	3.3

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Mol	Chain	Res	Type	RSRZ
1	H	43	GLY	3.3
1	D	35	GLU	3.2
1	H	74	ALA	3.2
1	J	78	VAL	3.2
1	J	167	VAL	3.2
1	I	48	PHE	3.2
1	E	5	SER	3.2
1	H	46	GLY	3.2
1	J	185	ARG	3.2
1	A	160	GLN	3.2
1	J	32	ILE	3.1
1	H	102	VAL	3.1
1	K	160	GLN	3.0
1	J	44	SER	3.0
1	J	173	LEU	3.0
1	H	67	MET	3.0
1	K	109	GLY	2.9
1	I	167	VAL	2.9
1	H	184	ALA	2.9
1	K	112	VAL	2.9
1	I	35	GLU	2.9
1	K	47	THR	2.9
1	J	91	VAL	2.8
1	L	141	TYR	2.8
1	I	178	ASP	2.8
1	I	17	ARG	2.8
1	I	37	ILE	2.8
1	J	11	THR	2.8
1	H	39	ILE	2.8
1	C	5	SER	2.8
1	K	185	ARG	2.7
1	F	159	ALA	2.7
1	L	72	ALA	2.7
1	I	78	VAL	2.7
1	H	50	HIS	2.7
1	J	8	HIS	2.7
1	I	19	ALA	2.6
1	K	5	SER	2.6
1	K	140	ALA	2.6
1	L	181	ALA	2.6
1	H	71	ILE	2.6
1	I	114	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	L	55	VAL	2.6
1	H	37	ILE	2.6
1	H	69	GLN	2.6
1	I	179	ALA	2.6
1	J	5	SER	2.6
1	I	153	VAL	2.6
1	I	51	TYR	2.6
1	K	36	GLY	2.6
1	K	188	SER	2.6
1	J	180	LEU	2.6
1	I	9	PHE	2.6
1	E	36	GLY	2.6
1	K	181	ALA	2.5
1	L	183	LEU	2.5
1	J	182	ALA	2.5
1	H	186	THR	2.5
1	I	185	ARG	2.5
1	I	44	SER	2.5
1	L	179	ALA	2.5
1	H	110	LYS	2.5
1	H	54	ASP	2.5
1	L	178	ASP	2.5
1	J	28	ALA	2.5
1	H	183	LEU	2.4
1	L	172	PHE	2.4
1	H	61	PRO	2.4
1	H	105	GLN	2.4
1	H	51	TYR	2.4
1	I	121	MET	2.4
1	J	145	ARG	2.4
1	J	114	LEU	2.4
1	K	67	MET	2.4
1	J	148	ALA	2.4
1	H	141	TYR	2.4
1	J	63	PRO	2.4
1	K	9	PHE	2.4
1	G	153	VAL	2.3
1	K	137	PHE	2.3
1	H	101	ALA	2.3
1	H	188	SER	2.3
1	L	112	VAL	2.3
1	J	70	GLN	2.3

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Mol	Chain	Res	Type	RSRZ
1	I	32	ILE	2.3
1	I	7	LEU	2.3
1	G	160	GLN	2.3
1	L	8	HIS	2.3
1	I	21	PHE	2.3
1	J	24	ALA	2.3
1	L	113	ALA	2.3
1	H	40	THR	2.3
1	J	40	THR	2.3
1	J	143	LEU	2.3
1	L	160	GLN	2.2
1	L	7	LEU	2.2
1	H	52	SER	2.2
1	K	10	VAL	2.2
1	I	13	LEU	2.2
1	K	104	PRO	2.2
1	F	120	GLY	2.2
1	G	33	ALA	2.1
1	K	106	PRO	2.1
1	L	60	PHE	2.1
1	J	184	ALA	2.1
1	L	125	ALA	2.1
1	G	39	ILE	2.1
1	J	71	ILE	2.1
1	H	138	LEU	2.1
1	G	8	HIS	2.1
1	J	50	HIS	2.1
1	I	141	TYR	2.1
1	G	188	SER	2.1
1	J	144	ASN	2.1
1	I	84	TYR	2.1
1	K	101	ALA	2.1
1	K	102	VAL	2.1
1	G	15	SER	2.1
1	I	172	PHE	2.1
1	G	104	PRO	2.1
1	H	180	LEU	2.1
1	J	170	ARG	2.1
1	K	63	PRO	2.0
1	G	19	ALA	2.0
1	L	28	ALA	2.0
1	I	8	HIS	2.0

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Mol	Chain	Res	Type	RSRZ
1	G	10	VAL	2.0
1	D	6	PRO	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 monosaccharides in this entry.

6.4 Ligands [i](#)

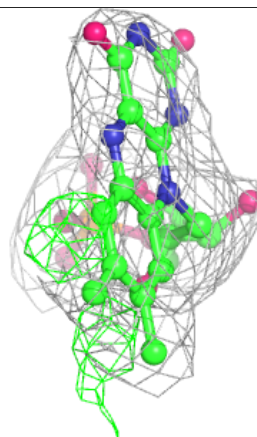
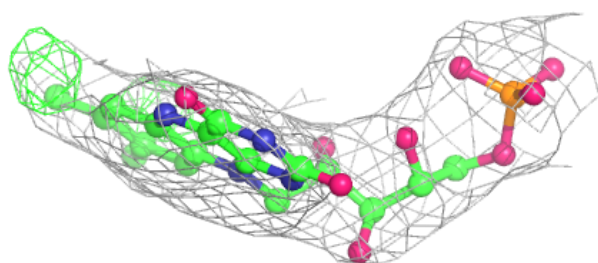
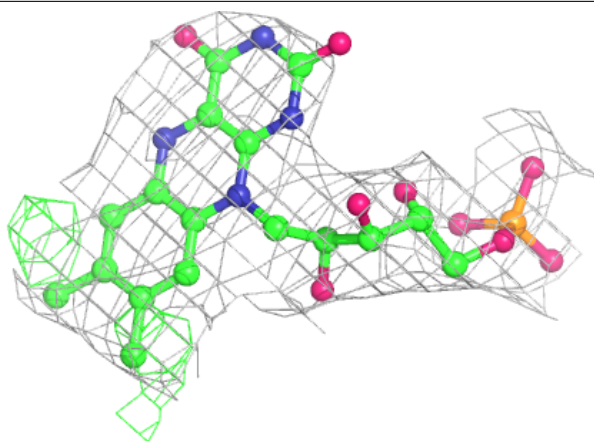
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	FMN	G	201	31/31	0.78	0.21	62,72,77,79	0
2	FMN	I	201	31/31	0.79	0.19	97,98,121,121	0
2	FMN	J	201	31/31	0.83	0.19	60,65,112,112	0
2	FMN	H	201	31/31	0.90	0.21	49,66,74,77	0
2	FMN	L	201	31/31	0.90	0.17	61,67,75,80	0
2	FMN	K	201	31/31	0.91	0.16	60,73,91,92	0
2	FMN	E	201	31/31	0.94	0.14	33,43,47,49	0
2	FMN	C	201	31/31	0.96	0.14	25,30,32,34	0
2	FMN	B	201	31/31	0.96	0.16	20,24,30,31	0
2	FMN	F	201	31/31	0.96	0.16	25,28,30,32	0
2	FMN	D	201	31/31	0.96	0.13	25,31,37,40	0
2	FMN	A	201	31/31	0.97	0.14	23,25,29,34	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.

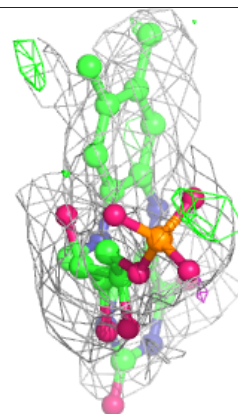
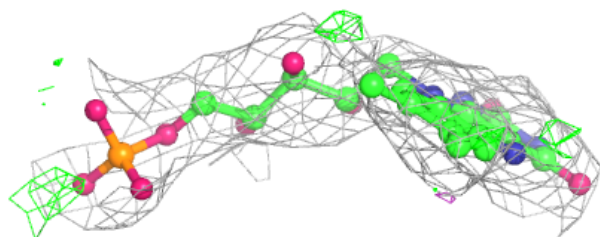
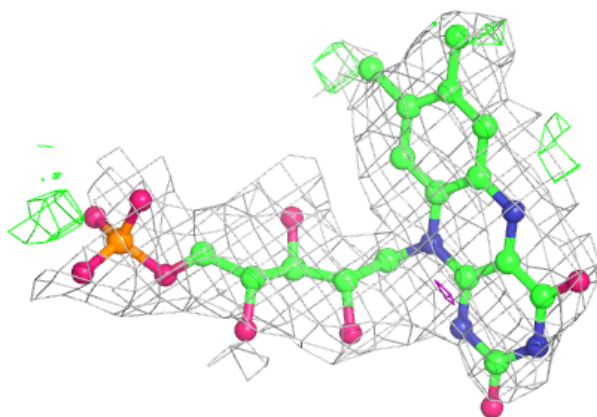
Electron density around FMN G 201:

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



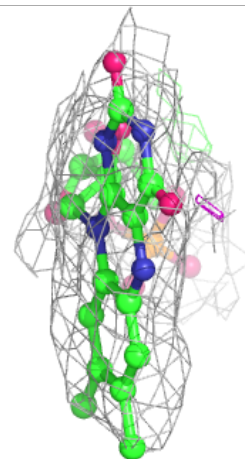
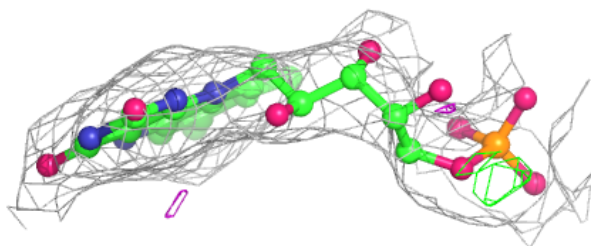
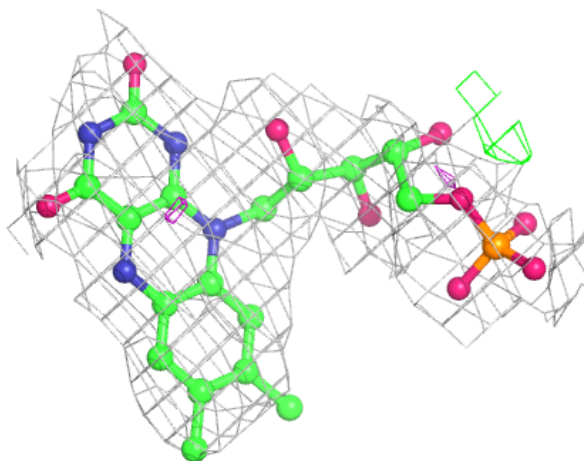
Electron density around FMN I 201:

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



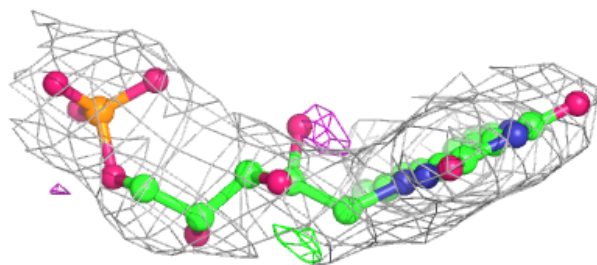
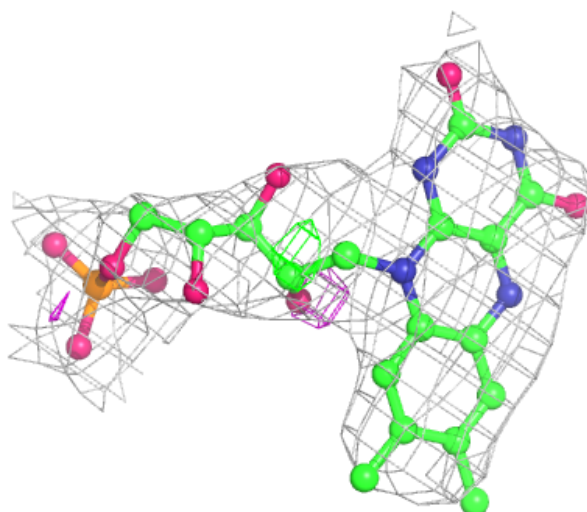
Electron density around FMN J 201:

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



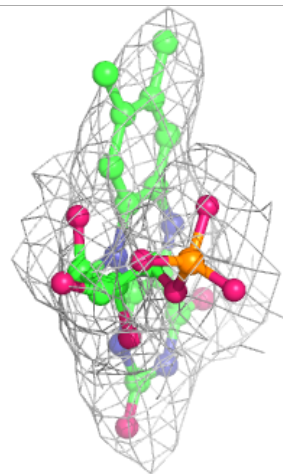
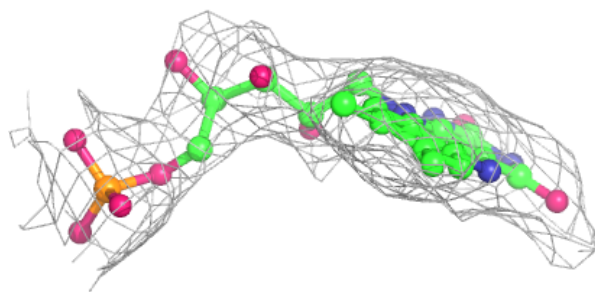
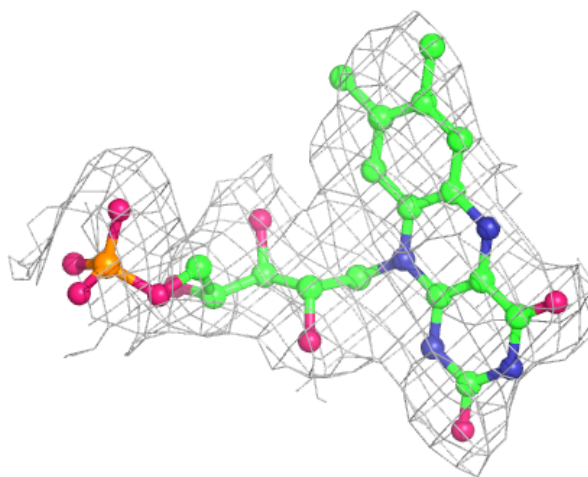
Electron density around FMN H 201:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



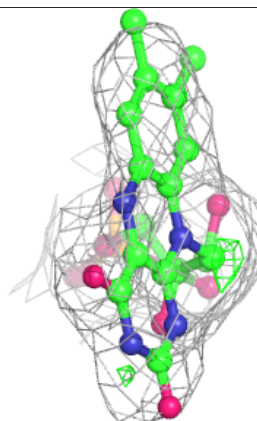
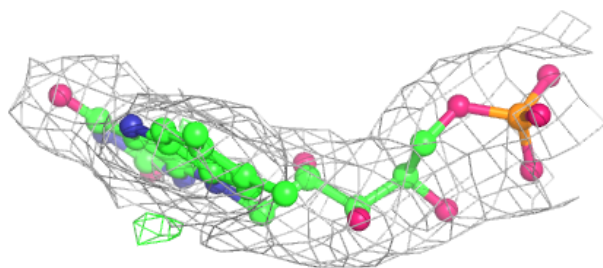
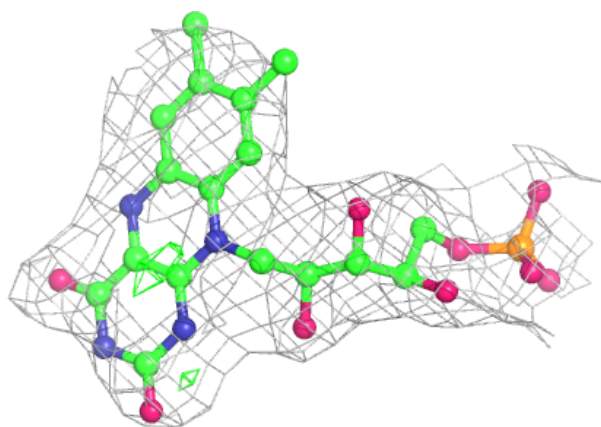
Electron density around FMN L 201:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



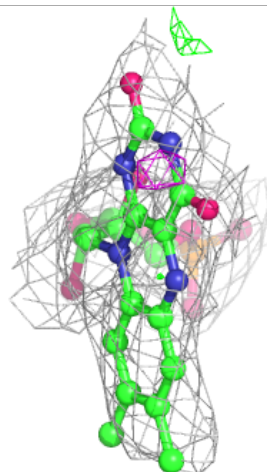
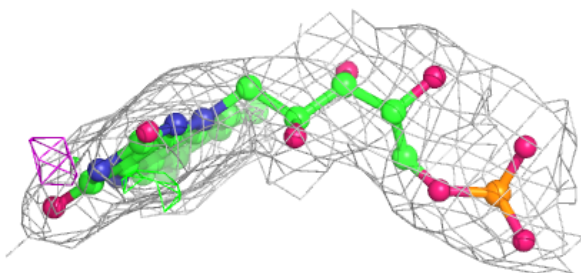
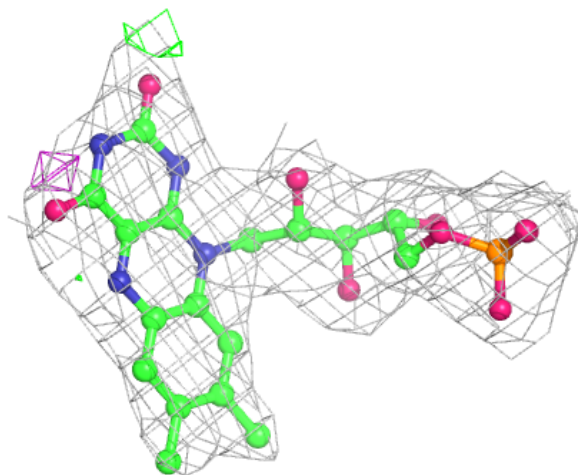
Electron density around FMN K 201:

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



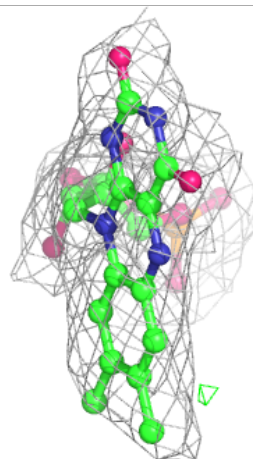
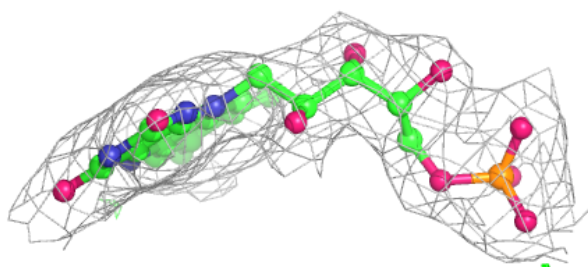
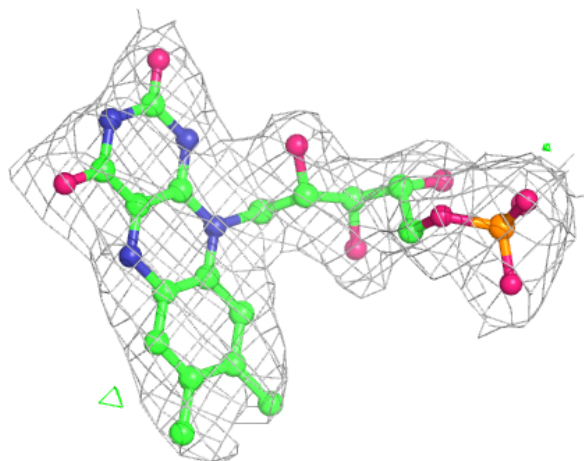
Electron density around FMN E 201:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



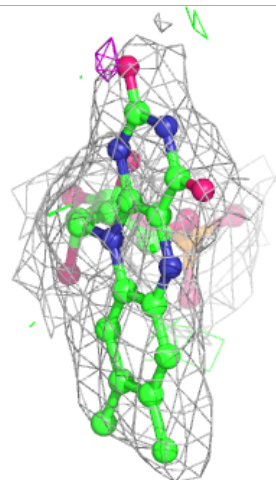
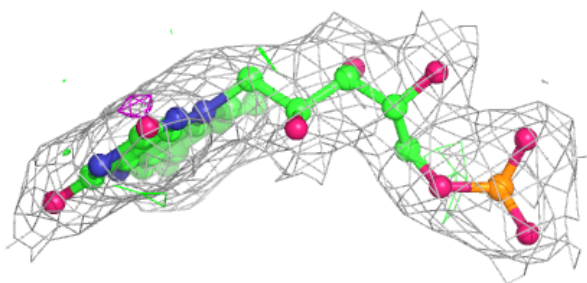
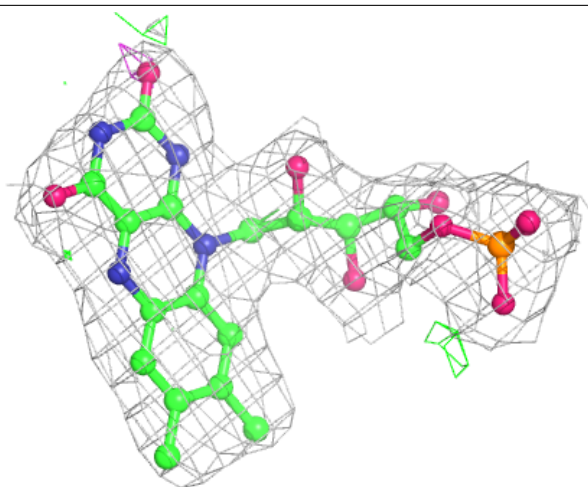
Electron density around FMN C 201:

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



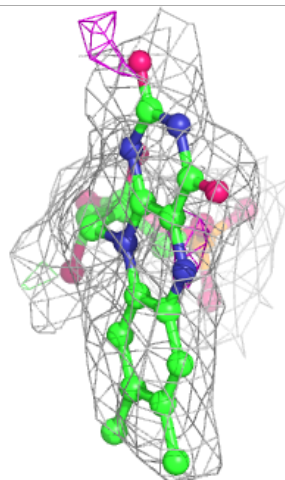
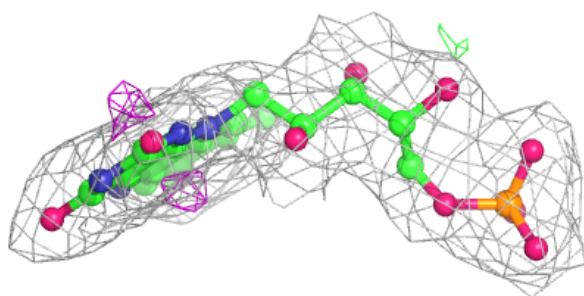
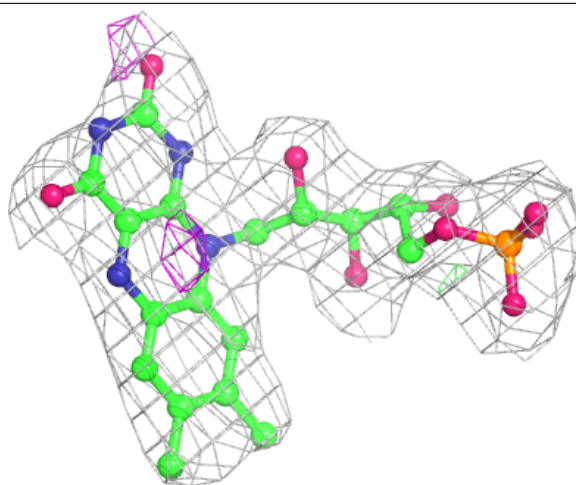
Electron density around FMN B 201:

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



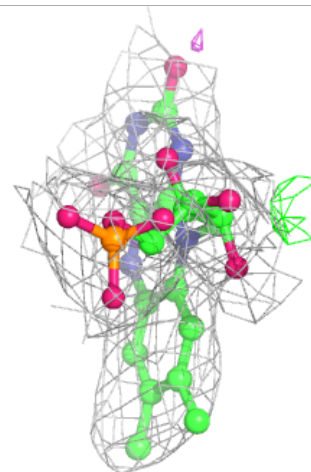
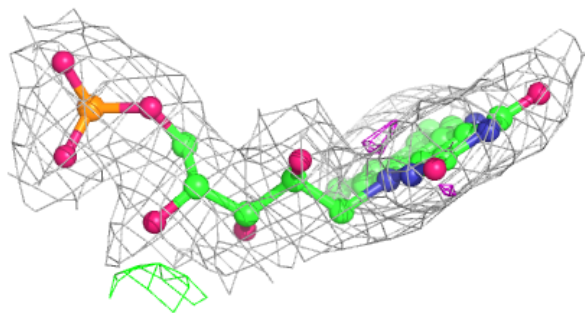
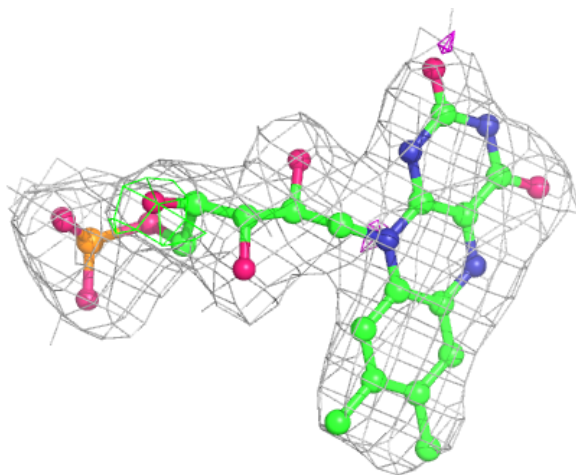
Electron density around FMN F 201:

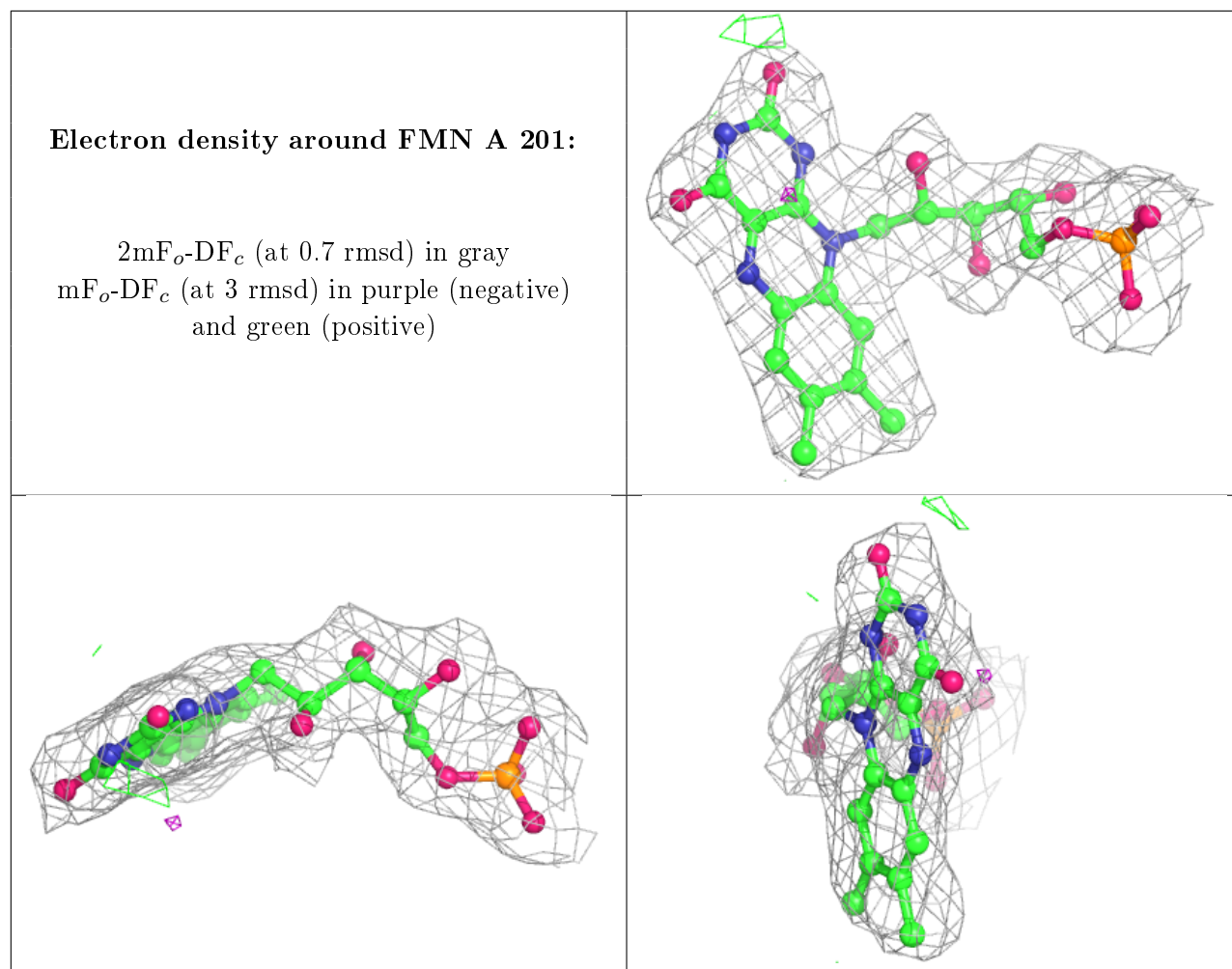
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around FMN D 201:

$2mF_o-DF_c$ (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.