



# Full wwPDB X-ray Structure Validation Report ⓘ

Dec 17, 2023 – 04:12 am GMT

PDB ID : 2YJ0  
Title : X-ray structure of chemically engineered Mycobacterium tuberculosis Dodecin  
Authors : Vinzenz, X.; Grosse, W.; Linne, U.; Meissner, B.; Essen, L.-O.  
Deposited on : 2011-05-17  
Resolution : 2.40 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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 : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

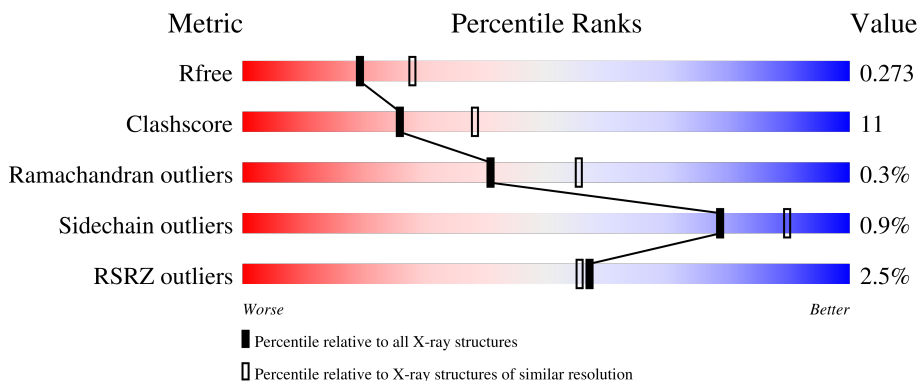
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	69	
1	B	69	
1	C	69	
1	D	69	
1	E	69	

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Mol	Chain	Length	Quality of chain
1	F	69	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	420	A	100[A]	-	-	X	-
2	420	C	100[A]	-	-	X	-
3	COA	B	200	X	-	-	-

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 4001 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.

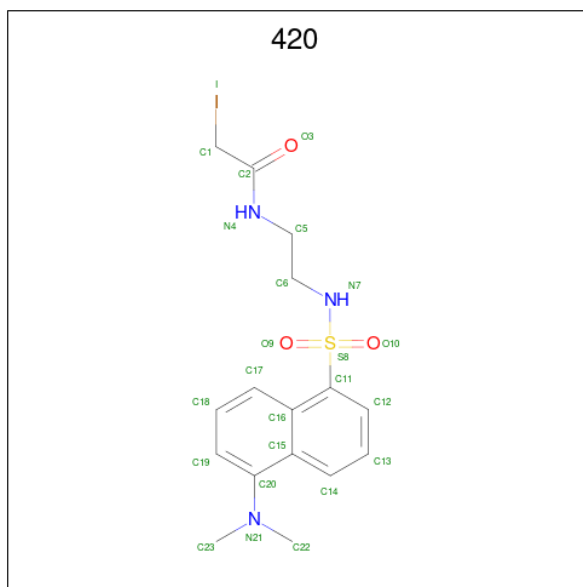
- Molecule 1 is a protein called DODECIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	68	529	332	97	96	4	0	2	0
1	B	67	525	328	97	96	4	0	2	0
1	C	68	523	327	96	96	4	0	1	0
1	D	68	528	330	97	97	4	0	2	0
1	E	68	523	327	96	96	4	0	1	0
1	F	68	528	330	97	97	4	0	2	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	59	CYS	THR	engineered mutation	UNP Q8VK10
B	59	CYS	THR	engineered mutation	UNP Q8VK10
C	59	CYS	THR	engineered mutation	UNP Q8VK10
D	59	CYS	THR	engineered mutation	UNP Q8VK10
E	59	CYS	THR	engineered mutation	UNP Q8VK10
F	59	CYS	THR	engineered mutation	UNP Q8VK10

- Molecule 2 is N-[2-({[5-(DIMETHYLAMINO)NAPHTHALEN-1-YL]SULFONYL}AMINO)ETHYL]-2-IODOACETAMIDE (three-letter code: 420) (formula: C<sub>16</sub>H<sub>20</sub>IN<sub>3</sub>O<sub>3</sub>S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
2	A	1	Total	C	N	O	S	0	1
			30	20	5	4	1		
2	B	1	Total	C	N	O	S	0	1
			33	20	5	6	2		
2	C	1	Total	C	N	O	S	0	1
			33	20	5	6	2		
2	D	1	Total	C	N	O	S	0	1
			33	20	5	6	2		
2	E	1	Total	C	N	O	S	0	1
			33	20	5	6	2		
2	F	1	Total	C	N	O	S	0	1
			33	20	5	6	2		

- Molecule 3 is COENZYME A (three-letter code: COA) (formula:  $C_{21}H_{36}N_7O_{16}P_3S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	P			S
3	A	1	48	21	7	16	3	1	0	0
3	B	1	48	21	7	16	3	1	0	0
3	C	1	48	21	7	16	3	1	0	0
3	D	1	48	21	7	16	3	1	0	0
3	E	1	48	21	7	16	3	1	0	0
3	F	1	48	21	7	16	3	1	0	0

- Molecule 4 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	
4	A	1	Total	C	N	O	P	0	1
			31	17	4	9	1		
4	B	1	Total	C	N	O	P	0	1
			31	17	4	9	1		
4	C	1	Total	C	N	O	P	0	1
			31	17	4	9	1		
4	D	1	Total	C	N	O	P	0	1
			31	17	4	9	1		
4	E	1	Total	C	N	O	P	0	1
			31	17	4	9	1		
4	F	1	Total	C	N	O	P	0	1
			31	17	4	9	1		

- Molecule 5 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Cl	0	0
			1	1		
5	B	2	Total	Cl	0	0
			2	2		

- Molecule 6 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 7 is water.

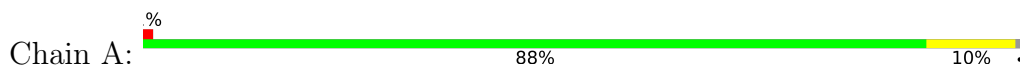
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	31	Total	O	0	0
			31	31		
7	B	26	Total	O	0	0
			26	26		
7	C	27	Total	O	0	0
			27	27		
7	D	31	Total	O	0	0
			31	31		
7	E	22	Total	O	0	0
			22	22		
7	F	26	Total	O	0	0
			26	26		



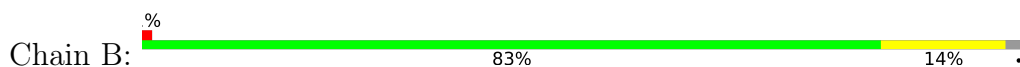
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

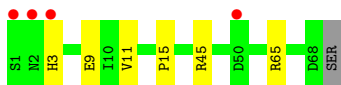
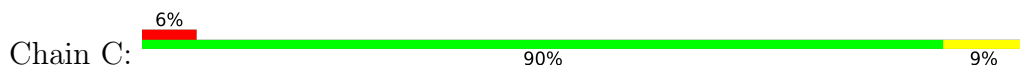
- Molecule 1: DODECIN



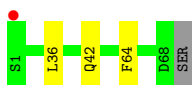
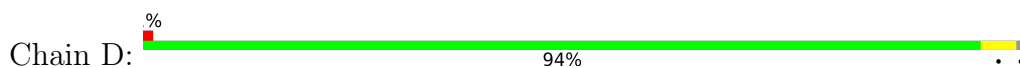
- Molecule 1: DODECIN



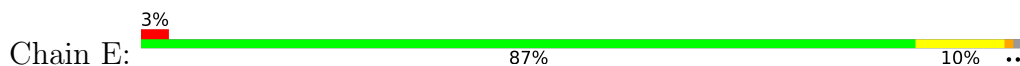
- Molecule 1: DODECIN




- Molecule 1: DODECIN



- Molecule 1: DODECIN



- Molecule 1: DODECIN

Chain F:  %  
87% 12%



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.00Å 103.80Å 82.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.40 47.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.8 (25.00-2.40) 99.8 (47.00-2.40)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.85 (at 2.39Å)	Xtrriage
Refinement program	REFMAC 5.5.0110	Depositor
R, $R_{free}$	0.217 , 0.275 0.219 , 0.273	Depositor DCC
$R_{free}$ test set	974 reflections (6.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	32.1	Xtrriage
Anisotropy	0.471	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 32.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	4001	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.26% of the height of the origin peak. No significant pseudotranslation is detected.*

<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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 420, FMN, SO4, COA, CL

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	0.31	0/544	0.56	0/734
1	B	0.31	0/537	0.53	0/726
1	C	0.29	0/535	0.53	0/723
1	D	0.33	0/543	0.57	0/734
1	E	0.30	0/535	0.52	0/723
1	F	0.33	0/543	0.58	0/734
All	All	0.31	0/3237	0.55	0/4374

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	529	0	526	8	0
1	B	525	0	509	14	0
1	C	523	0	513	10	0
1	D	528	0	519	3	0
1	E	523	0	513	8	0
1	F	528	0	519	6	0
2	A	30	0	24	13	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	33	0	15	8	0
2	C	33	0	16	10	0
2	D	33	0	16	5	0
2	E	33	0	16	15	0
2	F	33	0	24	7	0
3	A	48	0	32	1	0
3	B	48	0	32	1	0
3	C	48	0	32	0	0
3	D	48	0	32	0	0
3	E	48	0	32	0	0
3	F	48	0	32	0	0
4	A	31	0	19	1	0
4	B	31	0	9	0	0
4	C	31	0	8	2	0
4	D	31	0	19	0	0
4	E	31	0	10	0	0
4	F	31	0	13	0	0
5	A	1	0	0	0	0
5	B	2	0	0	0	0
6	D	5	0	0	0	0
6	E	5	0	0	0	0
7	A	31	0	0	1	0
7	B	26	0	0	0	0
7	C	27	0	0	0	0
7	D	31	0	0	0	0
7	E	22	0	0	0	0
7	F	26	0	0	0	0
All	All	4001	0	3480	77	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 11.

All (77) 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:F:45:ARG:HD3	2:F:100[A]:420:O9	1.46	1.16
2:A:100[B]:420:O3	2:A:100[B]:420:H62C	1.47	1.13
2:B:100[A]:420:C22	2:B:100[A]:420:H14	1.88	1.03
1:B:45:ARG:HD2	2:B:100[A]:420:O9	1.61	1.00
2:A:100[A]:420:C23	2:A:100[A]:420:H14	1.92	0.98
1:B:45:ARG:CD	2:B:100[A]:420:O9	2.12	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:LYS:CD	2:E:100[B]:420:O9	2.18	0.90
2:C:100[A]:420:H14	2:C:100[A]:420:C23	2.01	0.89
1:D:42:GLN:NE2	2:D:100[B]:420:O10	2.08	0.87
1:A:45:ARG:CD	2:A:100[A]:420:O9	2.24	0.85
1:A:45:ARG:HD2	2:A:100[A]:420:O9	1.77	0.82
2:D:100[A]:420:C23	2:D:100[A]:420:H14	2.10	0.80
2:A:100[A]:420:H14	2:A:100[A]:420:H232	1.62	0.80
2:A:100[A]:420:H14	2:A:100[A]:420:H233	1.63	0.80
1:F:45:ARG:CD	2:F:100[A]:420:O9	2.29	0.79
1:C:45:ARG:HD3	2:C:100[A]:420:O9	1.85	0.77
2:E:100[A]:420:H14	2:E:100[A]:420:C23	2.15	0.76
1:E:49:VAL:HG23	1:E:54:ALA:HB2	1.68	0.76
2:A:100[B]:420:O3	2:A:100[B]:420:C6	2.33	0.72
1:B:61:LYS:NZ	2:E:100[B]:420:O9	2.22	0.72
1:B:45:ARG:HD3	2:B:100[A]:420:O9	1.90	0.71
2:F:100[A]:420:H14	2:F:100[A]:420:C23	2.22	0.70
1:B:61:LYS:HD3	2:E:100[B]:420:O9	1.91	0.70
1:F:42:GLN:HB3	2:F:100[B]:420:H61C	1.75	0.69
1:A:45:ARG:HD3	2:A:100[A]:420:O9	1.91	0.68
1:B:61:LYS:HD2	2:E:100[B]:420:O9	1.92	0.68
2:D:100[A]:420:H7	2:D:100[A]:420:C17	2.07	0.68
3:B:200:COA:H141	3:B:200:COA:O9P	1.94	0.68
1:C:45:ARG:CD	2:C:100[A]:420:O9	2.42	0.67
2:B:100[A]:420:H7	2:B:100[A]:420:C17	2.07	0.66
4:C:300[B]:FMN:O3P	4:C:300[B]:FMN:O3'	2.09	0.66
1:D:42:GLN:HB3	2:D:100[B]:420:H61C	1.79	0.65
2:C:100[A]:420:C23	2:C:100[A]:420:C14	2.76	0.64
1:E:2:ASN:HD22	1:E:3:HIS:H	1.46	0.61
1:B:11:VAL:HG11	2:B:100[A]:420:H12	1.84	0.60
2:B:100[A]:420:C17	2:B:100[A]:420:N7	2.65	0.59
1:B:40:GLU:OE1	2:E:100[B]:420:O10	2.23	0.56
1:A:36:LEU:HD23	1:A:64:PHE:HB3	1.88	0.55
2:C:100[B]:420:C1	2:C:100[B]:420:C6	2.84	0.55
2:E:100[A]:420:O3	2:E:100[A]:420:H61C	2.06	0.54
2:B:100[A]:420:C22	2:B:100[A]:420:C14	2.63	0.53
1:C:9:GLU:OE2	2:C:100[A]:420:O3	2.27	0.53
1:F:36:LEU:HD23	1:F:64:PHE:HB3	1.92	0.52
2:E:100[A]:420:N7	2:E:100[A]:420:C17	2.73	0.52
1:A:11:VAL:HG11	2:A:100[A]:420:H12	1.92	0.51
2:F:100[A]:420:H14	2:F:100[A]:420:H233	1.91	0.51
1:C:11:VAL:HG11	2:C:100[A]:420:O10	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:15:PRO:CD	1:E:1:SER:HA	2.41	0.51
2:F:100[A]:420:H14	2:F:100[A]:420:H232	1.91	0.50
1:A:1:SER:N	7:A:2001:HOH:O	2.43	0.50
1:B:36:LEU:HD23	1:B:64:PHE:HB3	1.92	0.50
2:E:100[A]:420:C17	2:E:100[A]:420:H7	2.25	0.50
2:A:100[A]:420:C23	2:A:100[A]:420:C14	2.67	0.50
1:C:15:PRO:HD3	1:E:1:SER:HA	1.93	0.49
3:A:200:COA:H141	3:A:200:COA:O9P	2.11	0.49
2:E:100[A]:420:O3	2:E:100[A]:420:C6	2.59	0.49
1:E:45:ARG:CD	2:E:100[A]:420:O9	2.61	0.48
1:B:40:GLU:OE1	2:E:100[B]:420:N7	2.47	0.48
1:F:6:ARG:HB2	1:F:66:LEU:HD21	1.96	0.48
1:C:45:ARG:CD	2:C:100[A]:420:H62C	2.44	0.48
2:A:100[A]:420:H232	2:A:100[A]:420:C14	2.39	0.47
2:D:100[A]:420:C17	2:D:100[A]:420:N7	2.74	0.47
1:B:3:HIS:CE1	1:B:65:ARG:CZ	2.98	0.47
2:E:100[A]:420:C23	2:E:100[A]:420:C14	2.88	0.47
1:B:61:LYS:CE	2:E:100[B]:420:O9	2.63	0.46
1:E:45:ARG:HD3	2:E:100[A]:420:O9	2.15	0.46
2:A:100[A]:420:H233	2:A:100[A]:420:C14	2.39	0.46
1:E:49:VAL:CG2	1:E:54:ALA:HB2	2.40	0.46
4:A:300[B]:FMN:H1'2	4:C:300[B]:FMN:C2	2.46	0.45
1:C:45:ARG:HD2	2:C:100[A]:420:O9	2.16	0.45
1:D:36:LEU:HD23	1:D:64:PHE:HB3	1.99	0.44
1:A:45:ARG:HD2	2:A:100[A]:420:H62C	2.00	0.44
1:C:45:ARG:NE	2:C:100[A]:420:H62C	2.32	0.44
1:A:26:LEU:HD13	1:B:48:LEU:HD21	2.01	0.43
1:C:3:HIS:CG	1:C:65:ARG:HD2	2.56	0.41
1:F:43:SER:HB3	2:F:100[B]:420:H62C	2.03	0.41
1:E:36:LEU:HD23	1:E:64:PHE:HB3	2.04	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	68/69 (99%)	66 (97%)	2 (3%)	0	100	100
1	B	66/69 (96%)	66 (100%)	0	0	100	100
1	C	67/69 (97%)	67 (100%)	0	0	100	100
1	D	68/69 (99%)	68 (100%)	0	0	100	100
1	E	67/69 (97%)	65 (97%)	2 (3%)	0	100	100
1	F	68/69 (99%)	66 (97%)	0	2 (3%)	4	4
All	All	404/414 (98%)	398 (98%)	4 (1%)	2 (0%)	41	41

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	F	2[A]	ASN
1	F	2[B]	ASN

### 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	55/54 (102%)	54 (98%)	1 (2%)	59	76
1	B	54/54 (100%)	53 (98%)	1 (2%)	57	75
1	C	54/54 (100%)	54 (100%)	0	100	100
1	D	55/54 (102%)	55 (100%)	0	100	100
1	E	54/54 (100%)	53 (98%)	1 (2%)	57	75
1	F	55/54 (102%)	55 (100%)	0	100	100
All	All	327/324 (101%)	324 (99%)	3 (1%)	78	90

All (3) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	A	2	ASN
1	B	50	ASP
1	E	2	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	2	ASN

### 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](#)

Of 29 ligands modelled in this entry, 3 are monoatomic - leaving 26 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	420	D	100[A]	1	24,24,25	1.36	2 (8%)	34,34,35	2.04	6 (17%)
4	FMN	D	300[B]	-	33,33,33	1.07	2 (6%)	48,50,50	1.20	6 (12%)
2	420	B	100[A]	1	24,24,25	1.32	2 (8%)	34,34,35	1.25	4 (11%)
2	420	E	100[B]	1	7,9,25	0.99	0	6,10,35	1.22	1 (16%)
2	420	C	100[A]	1	24,24,25	1.39	2 (8%)	34,34,35	1.75	6 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	420	A	100[B]	1	6,6,25	0.69	0	5,6,35	0.97	0
3	COA	C	200	-	41,50,50	1.72	3 (7%)	52,75,75	1.06	2 (3%)
4	FMN	C	300[B]	-	33,33,33	1.08	2 (6%)	48,50,50	1.21	7 (14%)
6	SO4	D	1070	-	4,4,4	0.14	0	6,6,6	0.08	0
3	COA	F	200	-	41,50,50	1.72	3 (7%)	52,75,75	0.98	2 (3%)
4	FMN	F	300[B]	-	33,33,33	1.07	2 (6%)	48,50,50	1.20	6 (12%)
6	SO4	E	1070	-	4,4,4	0.14	0	6,6,6	0.12	0
2	420	F	100[B]	1	7,9,25	0.98	0	6,10,35	0.64	0
2	420	E	100[A]	1	24,24,25	1.39	2 (8%)	34,34,35	2.00	6 (17%)
3	COA	E	200	-	41,50,50	1.70	3 (7%)	52,75,75	1.07	2 (3%)
3	COA	B	200	-	41,50,50	1.74	3 (7%)	52,75,75	1.19	4 (7%)
2	420	A	100[A]	1	24,24,25	1.41	2 (8%)	34,34,35	1.43	4 (11%)
4	FMN	E	300[B]	-	33,33,33	1.08	2 (6%)	48,50,50	1.25	7 (14%)
4	FMN	A	300[B]	-	33,33,33	1.07	2 (6%)	48,50,50	1.20	7 (14%)
3	COA	A	200	-	41,50,50	1.71	3 (7%)	52,75,75	1.10	2 (3%)
4	FMN	B	300[B]	-	33,33,33	1.09	2 (6%)	48,50,50	1.20	7 (14%)
2	420	D	100[B]	1	7,9,25	1.14	1 (14%)	6,10,35	0.52	0
2	420	F	100[A]	1	24,24,25	1.42	2 (8%)	34,34,35	2.30	9 (26%)
3	COA	D	200	-	41,50,50	1.67	3 (7%)	52,75,75	1.09	2 (3%)
2	420	B	100[B]	1	7,9,25	0.96	0	6,10,35	0.72	0
2	420	C	100[B]	1	7,9,25	1.04	0	6,10,35	0.61	0

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	420	D	100[A]	1	-	14/18/18/20	0/2/2/2
4	FMN	D	300[B]	-	-	7/18/18/18	0/3/3/3
2	420	B	100[A]	1	-	12/18/18/20	0/2/2/2
2	420	E	100[B]	1	-	1/5/7/20	-
2	420	C	100[A]	1	-	10/18/18/20	0/2/2/2
2	420	A	100[B]	1	-	2/4/4/20	-
3	COA	C	200	-	-	10/44/64/64	0/3/3/3
4	FMN	C	300[B]	-	-	10/18/18/18	0/3/3/3
3	COA	F	200	-	-	11/44/64/64	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	FMN	F	300[B]	-	-	11/18/18/18	0/3/3/3
2	420	F	100[B]	1	-	1/5/7/20	-
2	420	E	100[A]	1	-	15/18/18/20	0/2/2/2
3	COA	E	200	-	-	11/44/64/64	0/3/3/3
3	COA	B	200	-	1/1/11/13	12/44/64/64	0/3/3/3
2	420	A	100[A]	1	-	8/18/18/20	0/2/2/2
4	FMN	E	300[B]	-	-	8/18/18/18	0/3/3/3
4	FMN	A	300[B]	-	-	13/18/18/18	0/3/3/3
3	COA	A	200	-	-	9/44/64/64	0/3/3/3
4	FMN	B	300[B]	-	-	12/18/18/18	0/3/3/3
2	420	D	100[B]	1	-	1/5/7/20	-
2	420	F	100[A]	1	-	8/18/18/20	0/2/2/2
3	COA	D	200	-	-	11/44/64/64	0/3/3/3
2	420	B	100[B]	1	-	1/5/7/20	-
2	420	C	100[B]	1	-	4/5/7/20	-

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	200	COA	O9P-C9P	9.16	1.41	1.23
3	A	200	COA	O9P-C9P	9.01	1.41	1.23
3	F	200	COA	O9P-C9P	8.99	1.41	1.23
3	C	200	COA	O9P-C9P	8.98	1.41	1.23
3	E	200	COA	O9P-C9P	8.97	1.41	1.23
3	D	200	COA	O9P-C9P	8.77	1.40	1.23
2	A	100[A]	420	S8-N7	4.86	1.69	1.61
2	E	100[A]	420	S8-N7	4.84	1.69	1.61
2	F	100[A]	420	S8-N7	4.81	1.69	1.61
2	C	100[A]	420	S8-N7	4.76	1.68	1.61
2	D	100[A]	420	S8-N7	4.66	1.68	1.61
2	B	100[A]	420	S8-N7	4.52	1.68	1.61
3	B	200	COA	C2A-N3A	4.17	1.38	1.32
3	C	200	COA	C2A-N3A	4.15	1.38	1.32
3	A	200	COA	C2A-N3A	4.13	1.38	1.32
3	F	200	COA	C2A-N3A	4.13	1.38	1.32
4	C	300[B]	FMN	C4A-N5	4.01	1.38	1.30
4	A	300[B]	FMN	C4A-N5	4.01	1.38	1.30
4	B	300[B]	FMN	C4A-N5	4.00	1.38	1.30
4	D	300[B]	FMN	C4A-N5	3.99	1.38	1.30
4	F	300[B]	FMN	C4A-N5	3.95	1.38	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	200	COA	C2A-N3A	3.86	1.38	1.32
4	E	300[B]	FMN	C4A-N5	3.83	1.38	1.30
3	E	200	COA	C2A-N3A	3.80	1.38	1.32
3	F	200	COA	C2A-N1A	2.75	1.39	1.33
3	C	200	COA	C2A-N1A	2.73	1.39	1.33
3	D	200	COA	C2A-N1A	2.61	1.38	1.33
3	A	200	COA	C2A-N1A	2.59	1.38	1.33
3	B	200	COA	C2A-N1A	2.53	1.38	1.33
3	E	200	COA	C2A-N1A	2.53	1.38	1.33
4	E	300[B]	FMN	C10-N1	2.53	1.38	1.33
4	C	300[B]	FMN	C10-N1	2.52	1.38	1.33
4	B	300[B]	FMN	C10-N1	2.52	1.38	1.33
4	D	300[B]	FMN	C10-N1	2.47	1.38	1.33
4	A	300[B]	FMN	C10-N1	2.45	1.38	1.33
4	F	300[B]	FMN	C10-N1	2.45	1.38	1.33
2	F	100[A]	420	C12-C11	2.40	1.40	1.37
2	A	100[A]	420	C12-C11	2.40	1.40	1.37
2	E	100[A]	420	C12-C11	2.36	1.40	1.37
2	C	100[A]	420	C12-C11	2.34	1.40	1.37
2	B	100[A]	420	C12-C11	2.22	1.40	1.37
2	D	100[A]	420	C12-C11	2.10	1.39	1.37
2	D	100[B]	420	S8-N7	2.03	1.69	1.65

All (90) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	100[A]	420	C15-C20-N21	9.03	131.22	118.23
2	D	100[A]	420	C15-C20-N21	8.11	129.90	118.23
2	E	100[A]	420	C15-C20-N21	8.10	129.89	118.23
2	C	100[A]	420	C15-C20-N21	6.65	127.80	118.23
3	E	200	COA	N3A-C2A-N1A	-5.87	119.50	128.68
3	A	200	COA	N3A-C2A-N1A	-5.50	120.08	128.68
3	B	200	COA	N3A-C2A-N1A	-5.47	120.13	128.68
3	C	200	COA	N3A-C2A-N1A	-5.43	120.19	128.68
3	F	200	COA	N3A-C2A-N1A	-5.41	120.23	128.68
3	D	200	COA	N3A-C2A-N1A	-5.36	120.30	128.68
2	A	100[A]	420	C15-C20-N21	5.12	125.60	118.23
2	F	100[A]	420	C19-C20-N21	-4.98	114.17	121.68
2	D	100[A]	420	C19-C20-N21	-4.73	114.55	121.68
2	E	100[A]	420	C19-C20-N21	-4.64	114.69	121.68
2	B	100[A]	420	C15-C20-N21	3.97	123.94	118.23
2	C	100[A]	420	C19-C20-N21	-3.84	115.89	121.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	100[A]	420	C20-C15-C16	3.56	123.17	119.33
3	A	200	COA	P2A-O3A-P1A	-3.54	120.68	132.83
3	B	200	COA	P2A-O3A-P1A	-3.47	120.93	132.83
4	E	300[B]	FMN	C4-N3-C2	-3.27	119.59	125.64
4	C	300[B]	FMN	C4-N3-C2	-3.08	119.94	125.64
2	A	100[A]	420	C19-C20-N21	-3.08	117.04	121.68
4	F	300[B]	FMN	C4-N3-C2	-3.07	119.96	125.64
4	A	300[B]	FMN	C4-N3-C2	-3.05	120.01	125.64
4	B	300[B]	FMN	C4-N3-C2	-3.02	120.05	125.64
4	D	300[B]	FMN	C4-N3-C2	-3.02	120.07	125.64
2	C	100[A]	420	C20-C15-C16	2.99	122.56	119.33
2	F	100[A]	420	C19-C20-C15	-2.98	114.61	119.94
3	C	200	COA	P2A-O3A-P1A	-2.96	122.67	132.83
2	F	100[A]	420	C20-C15-C16	2.83	122.38	119.33
2	B	100[A]	420	C19-C20-N21	-2.75	117.53	121.68
4	F	300[B]	FMN	C4A-C4-N3	2.73	120.12	113.19
3	E	200	COA	P2A-O3A-P1A	-2.71	123.53	132.83
4	E	300[B]	FMN	C4A-C4-N3	2.69	120.02	113.19
4	A	300[B]	FMN	C4A-C10-N10	2.68	120.40	116.48
2	F	100[A]	420	C17-C16-C11	-2.66	120.11	123.60
4	B	300[B]	FMN	C4A-C4-N3	2.66	119.94	113.19
4	D	300[B]	FMN	C4A-C4-N3	2.64	119.90	113.19
4	C	300[B]	FMN	C4A-C4-N3	2.64	119.89	113.19
4	A	300[B]	FMN	C4A-C4-N3	2.61	119.82	113.19
2	F	100[A]	420	C11-C16-C15	2.59	121.20	117.94
4	E	300[B]	FMN	C4A-C10-N10	2.59	120.26	116.48
4	C	300[B]	FMN	C4A-C10-N10	2.58	120.26	116.48
2	C	100[A]	420	C5-N4-C2	2.56	126.49	122.56
2	F	100[A]	420	C5-N4-C2	-2.55	118.63	122.56
2	E	100[A]	420	C19-C20-C15	-2.53	115.43	119.94
4	E	300[B]	FMN	O4-C4-C4A	-2.52	119.92	126.60
4	B	300[B]	FMN	C4A-C10-N10	2.50	120.14	116.48
4	B	300[B]	FMN	O4-C4-C4A	-2.50	119.97	126.60
4	A	300[B]	FMN	C10-C4A-N5	-2.49	119.56	124.86
4	D	300[B]	FMN	C4A-C10-N10	2.49	120.12	116.48
4	F	300[B]	FMN	O4-C4-C4A	-2.48	120.02	126.60
4	F	300[B]	FMN	C4A-C10-N10	2.47	120.10	116.48
2	E	100[A]	420	C20-C15-C16	2.47	122.00	119.33
3	F	200	COA	P2A-O3A-P1A	-2.47	124.36	132.83
2	D	100[A]	420	C19-C20-C15	-2.46	115.55	119.94
3	D	200	COA	P2A-O3A-P1A	-2.45	124.42	132.83
4	C	300[B]	FMN	O4-C4-C4A	-2.44	120.13	126.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	300[B]	FMN	O4-C4-C4A	-2.43	120.16	126.60
4	E	300[B]	FMN	C10-C4A-N5	-2.40	119.76	124.86
4	C	300[B]	FMN	C10-C4A-N5	-2.39	119.80	124.86
4	D	300[B]	FMN	C10-C4A-N5	-2.37	119.82	124.86
4	A	300[B]	FMN	O4-C4-C4A	-2.34	120.38	126.60
2	D	100[A]	420	C23-N21-C20	2.34	121.08	114.12
3	B	200	COA	CEP-CBP-CAP	2.34	112.87	108.82
2	F	100[A]	420	C23-N21-C20	2.33	121.06	114.12
2	E	100[B]	420	C1-C2-N4	-2.32	111.97	116.09
4	B	300[B]	FMN	C10-C4A-N5	-2.29	120.00	124.86
4	F	300[B]	FMN	C10-C4A-N5	-2.29	120.01	124.86
4	E	300[B]	FMN	C4A-C10-N1	-2.26	119.48	124.73
2	E	100[A]	420	C23-N21-C20	2.24	120.78	114.12
4	E	300[B]	FMN	C9A-C5A-N5	-2.23	120.01	122.43
2	A	100[A]	420	C20-C15-C16	2.23	121.73	119.33
4	A	300[B]	FMN	C9A-C5A-N5	-2.18	120.06	122.43
2	B	100[A]	420	C1-C2-N4	-2.18	112.23	116.09
4	C	300[B]	FMN	C9A-C5A-N5	-2.17	120.08	122.43
4	D	300[B]	FMN	C9A-C5A-N5	-2.16	120.08	122.43
4	F	300[B]	FMN	C9A-C5A-N5	-2.15	120.09	122.43
2	D	100[A]	420	C16-C11-S8	2.14	124.13	121.58
4	C	300[B]	FMN	C4A-C10-N1	-2.14	119.77	124.73
2	F	100[A]	420	C18-C19-C20	2.13	123.63	120.66
4	B	300[B]	FMN	C9A-C5A-N5	-2.10	120.15	122.43
3	B	200	COA	C7P-C6P-C5P	-2.09	108.88	112.36
2	B	100[A]	420	O10-S8-O9	2.08	122.10	119.55
4	B	300[B]	FMN	C4A-C10-N1	-2.04	119.99	124.73
2	C	100[A]	420	C23-N21-C20	2.04	120.18	114.12
2	C	100[A]	420	C19-C20-C15	-2.03	116.31	119.94
2	E	100[A]	420	C11-C16-C15	2.03	120.50	117.94
2	A	100[A]	420	C22-N21-C20	2.02	120.13	114.12
4	A	300[B]	FMN	C4A-C10-N1	-2.02	120.04	124.73

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	B	200	COA	CAP

All (202) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	100[B]	420	C6-C5-N4-C2

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Mol	Chain	Res	Type	Atoms
2	B	100[A]	420	C12-C11-S8-N7
2	B	100[A]	420	C16-C11-S8-N7
2	B	100[A]	420	C16-C11-S8-O9
2	B	100[A]	420	C12-C11-S8-O10
2	B	100[A]	420	C16-C11-S8-O10
2	D	100[A]	420	C6-N7-S8-C11
2	D	100[A]	420	C12-C11-S8-N7
2	D	100[A]	420	C16-C11-S8-N7
2	D	100[A]	420	C12-C11-S8-O9
2	D	100[A]	420	C16-C11-S8-O9
2	D	100[A]	420	C12-C11-S8-O10
2	D	100[A]	420	C16-C11-S8-O10
2	E	100[A]	420	C6-C5-N4-C2
2	E	100[A]	420	C12-C11-S8-N7
2	E	100[A]	420	C16-C11-S8-N7
2	E	100[A]	420	C12-C11-S8-O9
2	E	100[A]	420	C16-C11-S8-O9
2	E	100[A]	420	C12-C11-S8-O10
2	E	100[A]	420	C16-C11-S8-O10
3	A	200	COA	CCP-O6A-P2A-O5A
3	A	200	COA	C6P-C5P-N4P-C3P
3	A	200	COA	S1P-C2P-C3P-N4P
3	B	200	COA	P1A-O3A-P2A-O6A
3	B	200	COA	O9P-C9P-CAP-OAP
3	C	200	COA	C3B-O3B-P3B-O9A
3	C	200	COA	CCP-O6A-P2A-O5A
3	D	200	COA	CAP-C9P-N8P-C7P
3	D	200	COA	O9P-C9P-N8P-C7P
3	E	200	COA	CCP-O6A-P2A-O5A
3	F	200	COA	CCP-O6A-P2A-O5A
3	F	200	COA	CAP-CBP-CCP-O6A
4	A	300[B]	FMN	C1'-C2'-C3'-O3'
4	A	300[B]	FMN	C1'-C2'-C3'-C4'
4	A	300[B]	FMN	O2'-C2'-C3'-O3'
4	A	300[B]	FMN	O2'-C2'-C3'-C4'
4	A	300[B]	FMN	C5'-O5'-P-O1P
4	A	300[B]	FMN	C5'-O5'-P-O2P
4	A	300[B]	FMN	C5'-O5'-P-O3P
4	B	300[B]	FMN	N10-C1'-C2'-O2'
4	B	300[B]	FMN	N10-C1'-C2'-C3'
4	B	300[B]	FMN	C1'-C2'-C3'-O3'
4	B	300[B]	FMN	C1'-C2'-C3'-C4'

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Mol	Chain	Res	Type	Atoms
4	B	300[B]	FMN	O2'-C2'-C3'-O3'
4	B	300[B]	FMN	O2'-C2'-C3'-C4'
4	B	300[B]	FMN	O3'-C3'-C4'-C5'
4	B	300[B]	FMN	O4'-C4'-C5'-O5'
4	C	300[B]	FMN	C1'-C2'-C3'-O3'
4	C	300[B]	FMN	C1'-C2'-C3'-C4'
4	C	300[B]	FMN	O2'-C2'-C3'-O3'
4	C	300[B]	FMN	O2'-C2'-C3'-C4'
4	C	300[B]	FMN	C3'-C4'-C5'-O5'
4	C	300[B]	FMN	O4'-C4'-C5'-O5'
4	C	300[B]	FMN	C5'-O5'-P-O1P
4	C	300[B]	FMN	C5'-O5'-P-O2P
4	C	300[B]	FMN	C5'-O5'-P-O3P
4	D	300[B]	FMN	C1'-C2'-C3'-O3'
4	D	300[B]	FMN	C1'-C2'-C3'-C4'
4	D	300[B]	FMN	O2'-C2'-C3'-O3'
4	D	300[B]	FMN	O2'-C2'-C3'-C4'
4	E	300[B]	FMN	C1'-C2'-C3'-O3'
4	E	300[B]	FMN	C1'-C2'-C3'-C4'
4	E	300[B]	FMN	O2'-C2'-C3'-O3'
4	E	300[B]	FMN	O2'-C2'-C3'-C4'
4	E	300[B]	FMN	C5'-O5'-P-O1P
4	E	300[B]	FMN	C5'-O5'-P-O2P
4	E	300[B]	FMN	C5'-O5'-P-O3P
4	F	300[B]	FMN	N10-C1'-C2'-O2'
4	F	300[B]	FMN	N10-C1'-C2'-C3'
4	F	300[B]	FMN	C1'-C2'-C3'-O3'
4	F	300[B]	FMN	C1'-C2'-C3'-C4'
4	F	300[B]	FMN	O2'-C2'-C3'-O3'
4	F	300[B]	FMN	O2'-C2'-C3'-C4'
4	F	300[B]	FMN	C3'-C4'-C5'-O5'
4	F	300[B]	FMN	O4'-C4'-C5'-O5'
4	F	300[B]	FMN	C5'-O5'-P-O1P
4	F	300[B]	FMN	C5'-O5'-P-O2P
4	F	300[B]	FMN	C5'-O5'-P-O3P
3	A	200	COA	O5P-C5P-N4P-C3P
2	A	100[A]	420	C19-C20-N21-C23
2	B	100[A]	420	C19-C20-N21-C22
2	C	100[A]	420	C19-C20-N21-C23
2	D	100[A]	420	C19-C20-N21-C23
2	E	100[A]	420	C19-C20-N21-C23
2	F	100[A]	420	C19-C20-N21-C23

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Mol	Chain	Res	Type	Atoms
2	A	100[A]	420	C15-C20-N21-C23
2	B	100[A]	420	C15-C20-N21-C22
2	C	100[A]	420	C15-C20-N21-C23
2	D	100[A]	420	C15-C20-N21-C23
2	E	100[A]	420	C15-C20-N21-C23
2	F	100[A]	420	C15-C20-N21-C23
2	D	100[A]	420	C6-N7-S8-O9
2	A	100[A]	420	C6-N7-S8-O10
2	C	100[A]	420	C6-N7-S8-O10
2	B	100[A]	420	C6-N7-S8-C11
3	D	200	COA	C6P-C7P-N8P-C9P
2	B	100[A]	420	N4-C5-C6-N7
2	C	100[B]	420	N4-C5-C6-N7
2	D	100[B]	420	N4-C5-C6-N7
2	E	100[A]	420	N4-C5-C6-N7
2	A	100[A]	420	C6-N7-S8-C11
2	C	100[A]	420	C6-N7-S8-C11
4	B	300[B]	FMN	C2'-C3'-C4'-C5'
2	C	100[A]	420	C1-C2-N4-C5
2	C	100[A]	420	O3-C2-N4-C5
2	C	100[B]	420	C1-C2-N4-C5
2	C	100[B]	420	O3-C2-N4-C5
2	D	100[A]	420	C1-C2-N4-C5
2	D	100[A]	420	O3-C2-N4-C5
2	C	100[A]	420	N4-C5-C6-N7
2	F	100[A]	420	N4-C5-C6-N7
4	B	300[B]	FMN	C2'-C3'-C4'-O4'
2	E	100[A]	420	C6-N7-S8-O9
2	B	100[A]	420	C12-C11-S8-O9
2	F	100[A]	420	C6-N7-S8-O10
2	B	100[B]	420	N4-C5-C6-N7
4	B	300[B]	FMN	O3'-C3'-C4'-O4'
2	E	100[A]	420	C6-N7-S8-O10
2	F	100[A]	420	C6-N7-S8-O9
4	C	300[B]	FMN	C4'-C5'-O5'-P
3	F	200	COA	C6P-C5P-N4P-C3P
3	F	200	COA	O5P-C5P-N4P-C3P
2	C	100[B]	420	C6-C5-N4-C2
2	F	100[B]	420	N4-C5-C6-N7
2	E	100[A]	420	C6-N7-S8-C11
2	C	100[A]	420	C19-C20-N21-C22
2	A	100[A]	420	C19-C20-N21-C22

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Mol	Chain	Res	Type	Atoms
2	B	100[A]	420	C19-C20-N21-C23
2	D	100[A]	420	C19-C20-N21-C22
2	E	100[A]	420	C19-C20-N21-C22
2	F	100[A]	420	C19-C20-N21-C22
3	F	200	COA	CDP-CBP-CCP-O6A
3	F	200	COA	CEP-CBP-CCP-O6A
2	A	100[A]	420	C15-C20-N21-C22
2	C	100[A]	420	C15-C20-N21-C22
2	E	100[A]	420	C15-C20-N21-C22
2	F	100[A]	420	C6-N7-S8-C11
2	B	100[A]	420	C15-C20-N21-C23
2	D	100[A]	420	C15-C20-N21-C22
2	F	100[A]	420	C15-C20-N21-C22
2	A	100[B]	420	N4-C5-C6-N7
3	A	200	COA	P1A-O3A-P2A-O6A
3	C	200	COA	P1A-O3A-P2A-O6A
3	D	200	COA	P1A-O3A-P2A-O6A
3	E	200	COA	P1A-O3A-P2A-O6A
3	F	200	COA	P1A-O3A-P2A-O6A
3	B	200	COA	N8P-C9P-CAP-OAP
3	A	200	COA	CCP-O6A-P2A-O3A
3	B	200	COA	CCP-O6A-P2A-O3A
3	C	200	COA	CCP-O6A-P2A-O3A
3	D	200	COA	C3B-O3B-P3B-O8A
3	E	200	COA	CCP-O6A-P2A-O3A
3	F	200	COA	CCP-O6A-P2A-O3A
3	B	200	COA	C4B-C5B-O5B-P1A
3	A	200	COA	CCP-O6A-P2A-O4A
3	B	200	COA	CCP-O6A-P2A-O4A
3	B	200	COA	CCP-O6A-P2A-O5A
3	C	200	COA	CCP-O6A-P2A-O4A
3	E	200	COA	CCP-O6A-P2A-O4A
3	F	200	COA	CCP-O6A-P2A-O4A
4	A	300[B]	FMN	N10-C1'-C2'-O2'
3	C	200	COA	CAP-CBP-CCP-O6A
3	E	200	COA	CAP-CBP-CCP-O6A
2	E	100[B]	420	N4-C5-C6-N7
3	C	200	COA	CDP-CBP-CCP-O6A
3	D	200	COA	C4B-C5B-O5B-P1A
3	D	200	COA	P1A-O3A-P2A-O4A
4	A	300[B]	FMN	C2'-C3'-C4'-O4'
3	A	200	COA	C4B-C5B-O5B-P1A

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Mol	Chain	Res	Type	Atoms
3	C	200	COA	C4B-C5B-O5B-P1A
3	E	200	COA	C4B-C5B-O5B-P1A
2	A	100[A]	420	N4-C5-C6-N7
3	B	200	COA	C5P-C6P-C7P-N8P
3	D	200	COA	C5P-C6P-C7P-N8P
3	B	200	COA	O9P-C9P-CAP-CBP
4	A	300[B]	FMN	C4'-C5'-O5'-P
2	C	100[A]	420	C6-N7-S8-O9
4	D	300[B]	FMN	O3'-C3'-C4'-C5'
3	C	200	COA	CEP-CBP-CCP-O6A
3	E	200	COA	CDP-CBP-CCP-O6A
3	E	200	COA	CEP-CBP-CCP-O6A
3	B	200	COA	N8P-C9P-CAP-CBP
3	F	200	COA	C4B-C5B-O5B-P1A
4	D	300[B]	FMN	O3'-C3'-C4'-O4'
4	E	300[B]	FMN	C4'-C5'-O5'-P
3	E	200	COA	C3B-O3B-P3B-O7A
4	B	300[B]	FMN	C5'-O5'-P-O2P
3	E	200	COA	O4B-C4B-C5B-O5B
3	D	200	COA	C3B-O3B-P3B-O9A
3	E	200	COA	C3B-O3B-P3B-O8A
4	A	300[B]	FMN	O3'-C3'-C4'-O4'
3	B	200	COA	C6P-C7P-N8P-C9P
3	C	200	COA	O4B-C4B-C5B-O5B
3	F	200	COA	O4B-C4B-C5B-O5B
4	A	300[B]	FMN	C2'-C3'-C4'-C5'
2	A	100[A]	420	C6-N7-S8-O9
4	D	300[B]	FMN	C2'-C3'-C4'-O4'
3	D	200	COA	CCP-O6A-P2A-O4A
3	A	200	COA	O4B-C4B-C5B-O5B
3	B	200	COA	O4B-C4B-C5B-O5B
3	D	200	COA	O4B-C4B-C5B-O5B
4	A	300[B]	FMN	N10-C1'-C2'-C3'

There are no ring outliers.

15 monomers are involved in 62 short contacts:

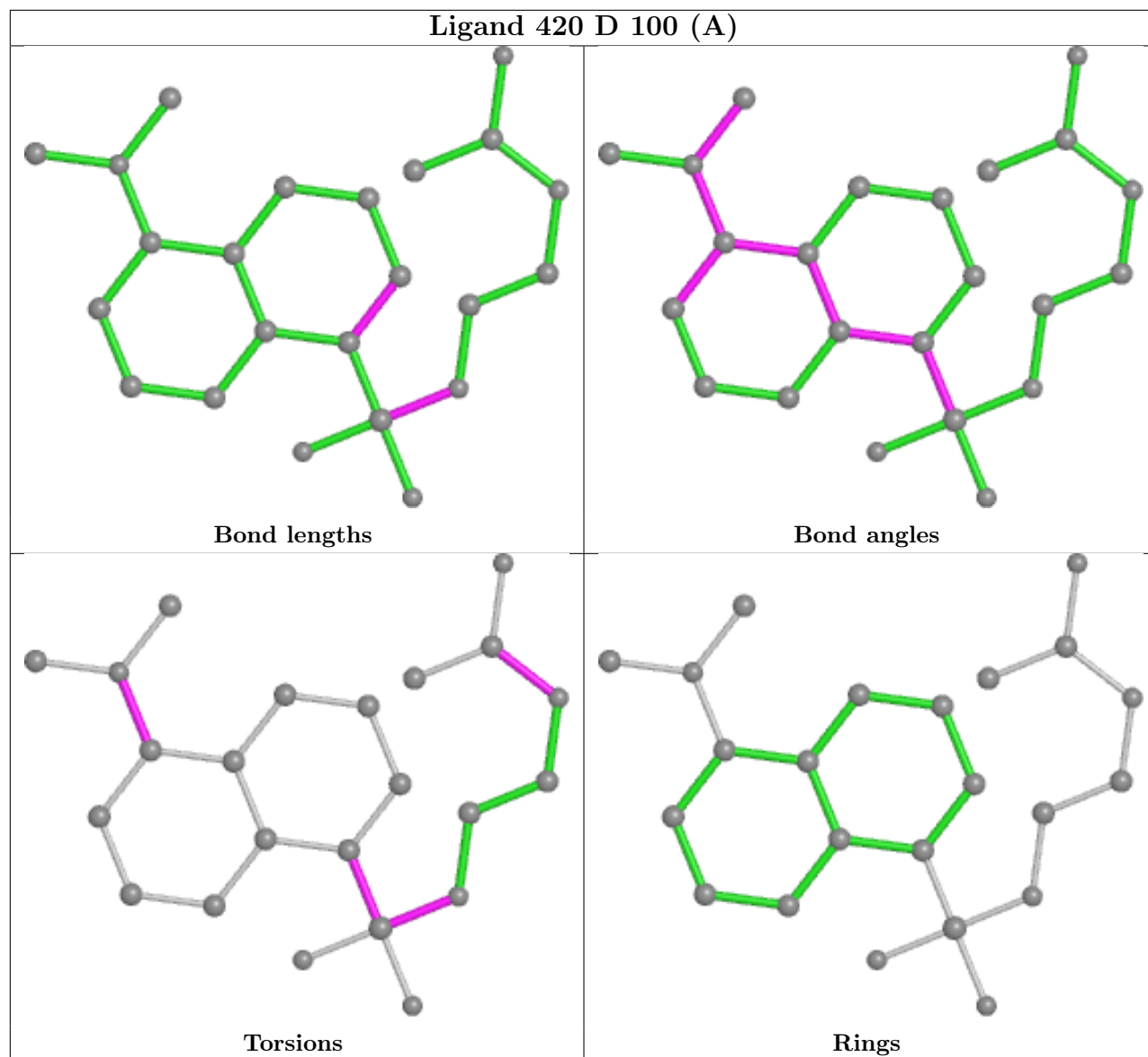
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	100[A]	420	3	0
2	B	100[A]	420	8	0
2	E	100[B]	420	7	0
2	C	100[A]	420	9	0

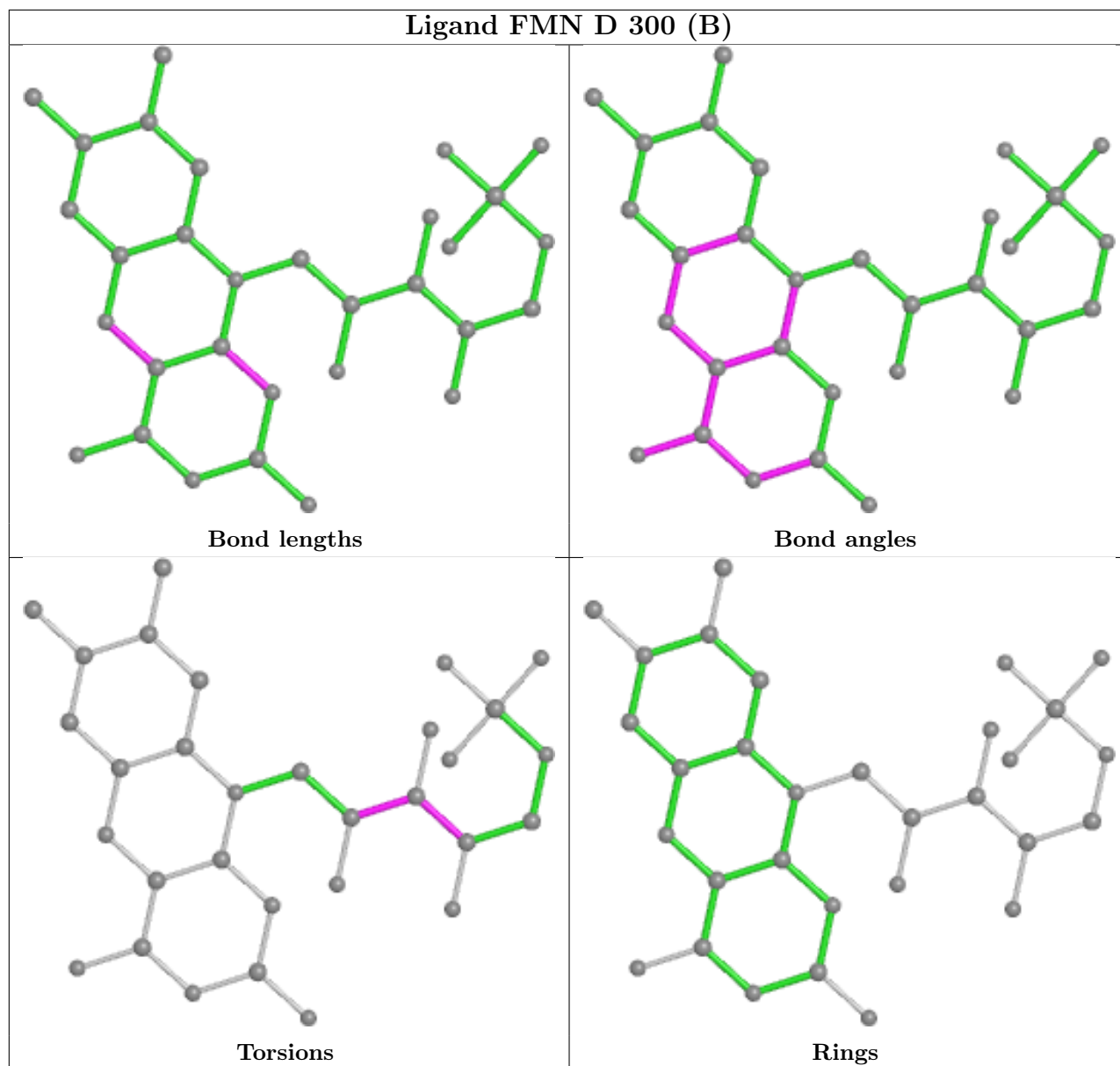
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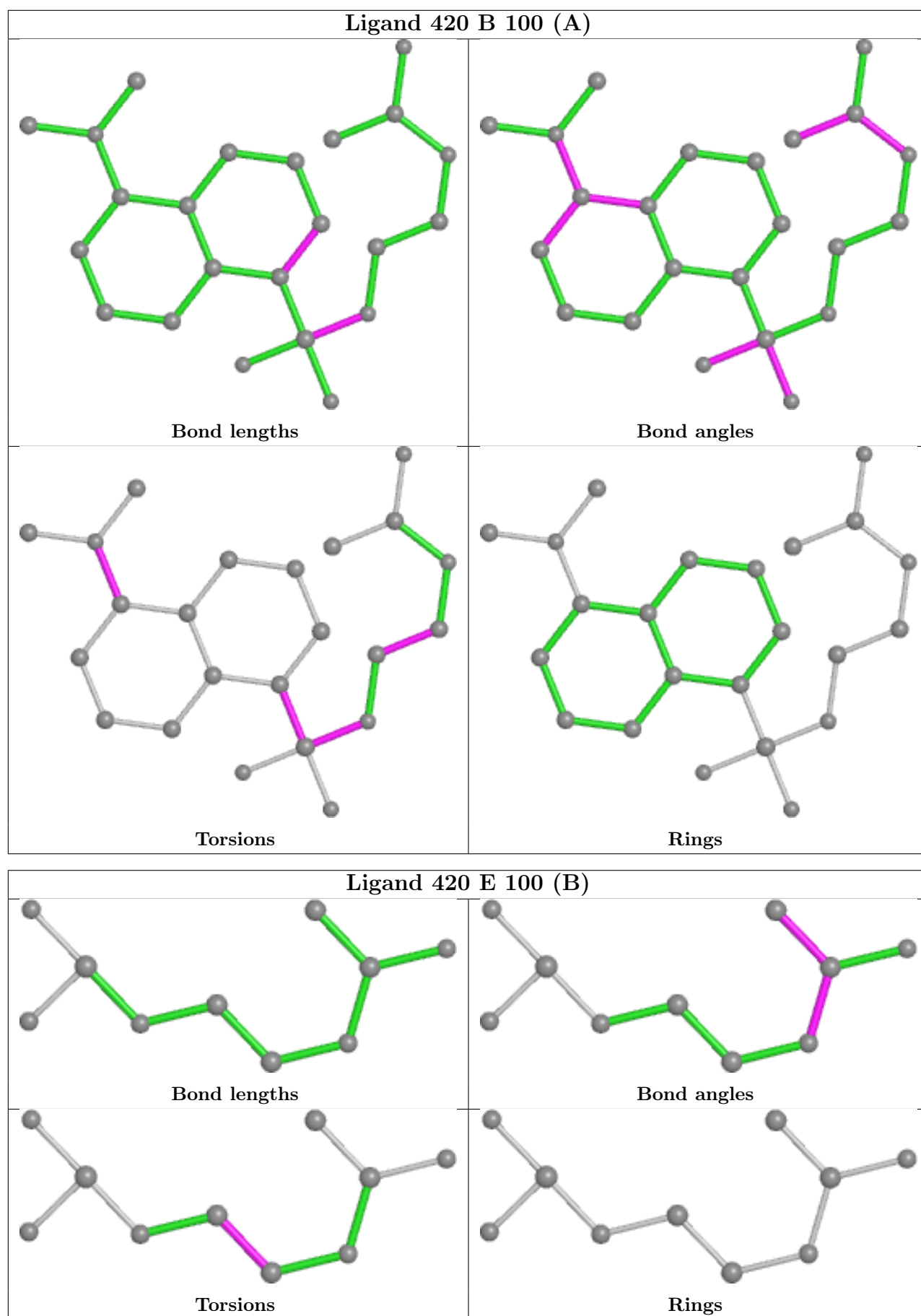
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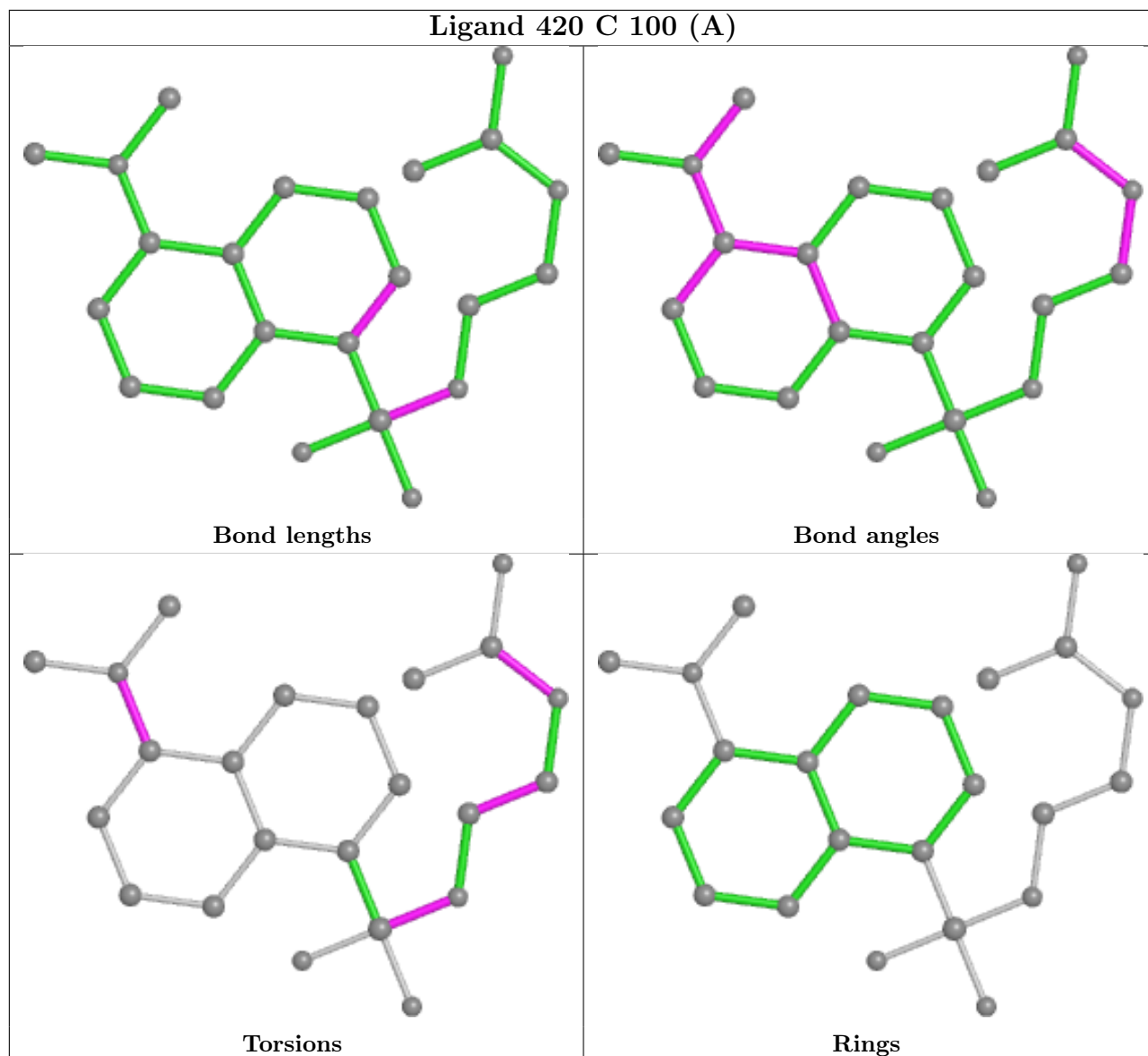
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	100[B]	420	2	0
4	C	300[B]	FMN	2	0
2	F	100[B]	420	2	0
2	E	100[A]	420	8	0
3	B	200	COA	1	0
2	A	100[A]	420	11	0
4	A	300[B]	FMN	1	0
3	A	200	COA	1	0
2	D	100[B]	420	2	0
2	F	100[A]	420	5	0
2	C	100[B]	420	1	0

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.

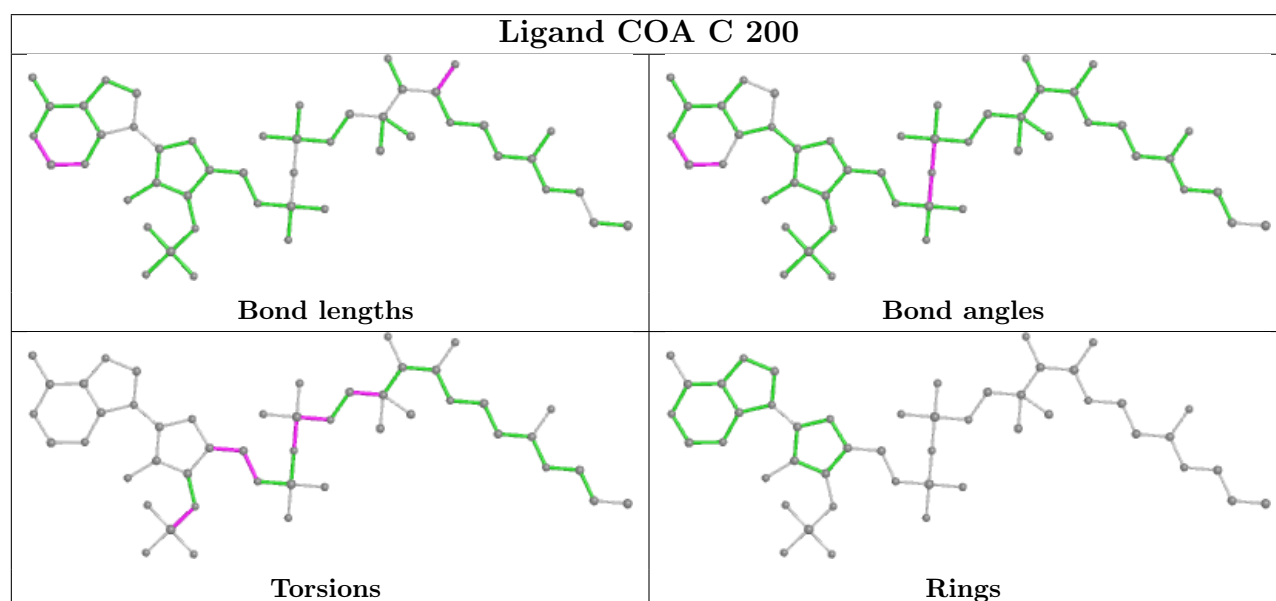
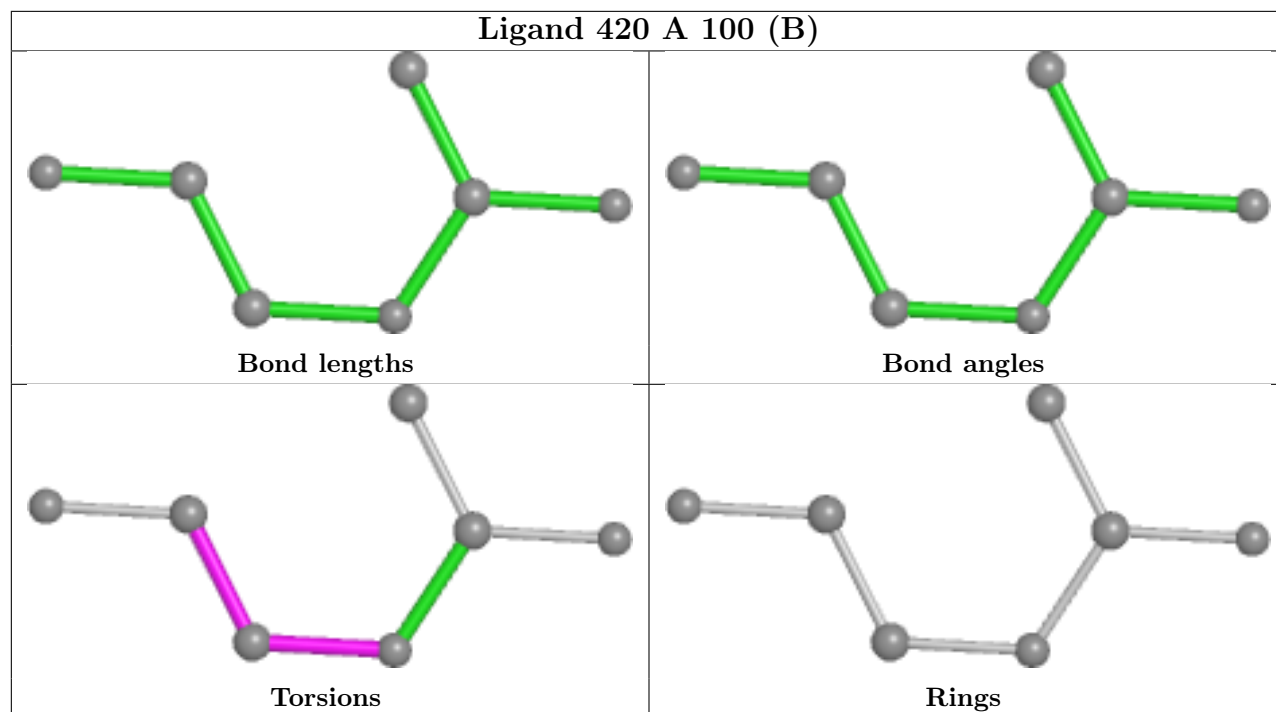


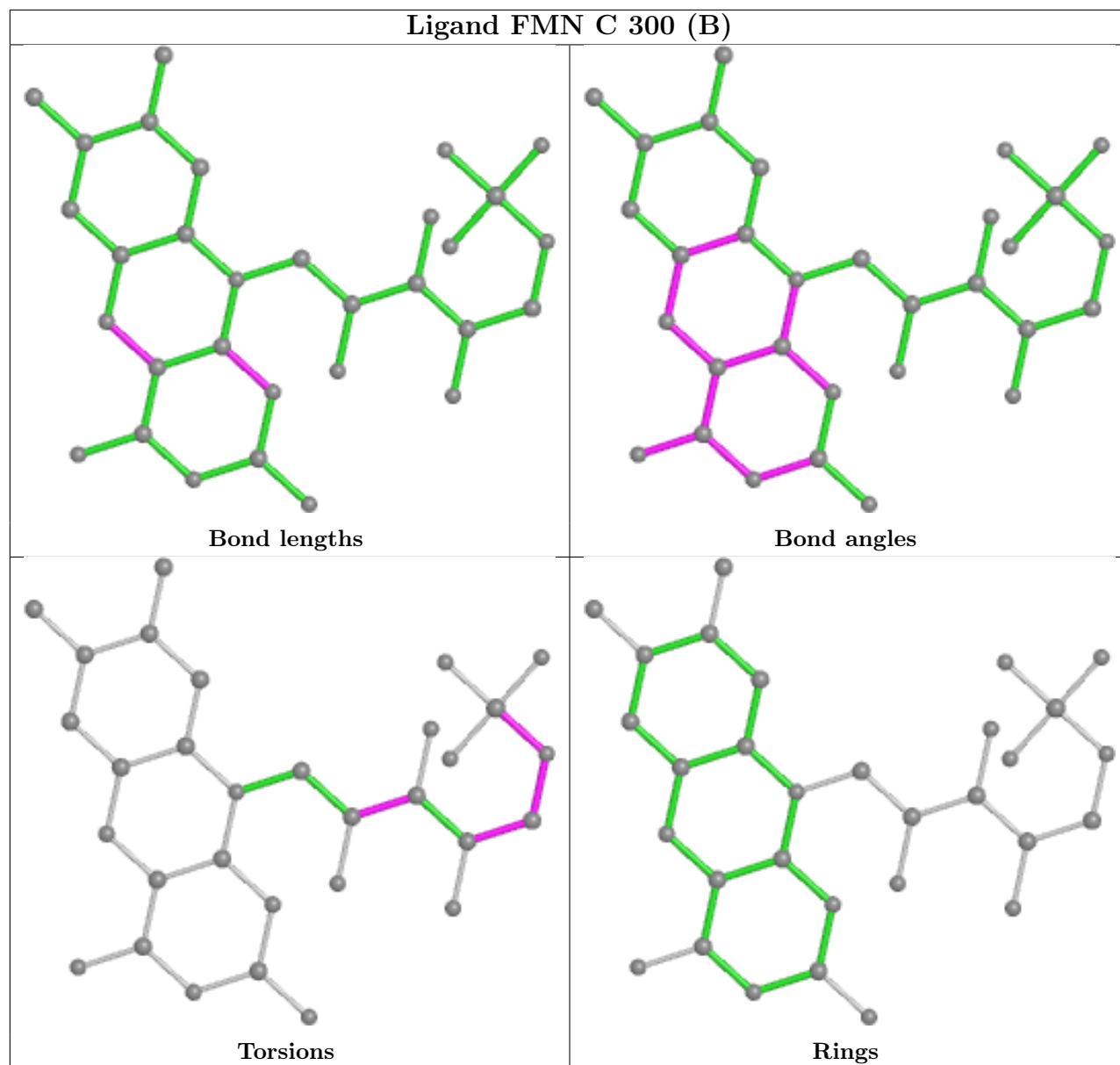


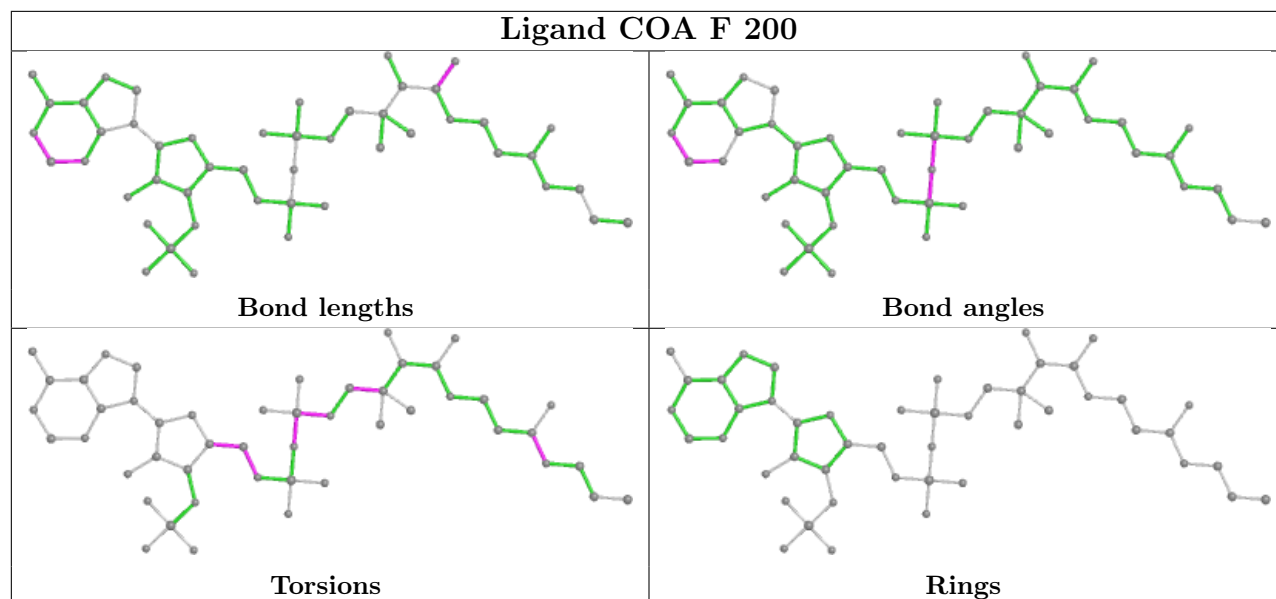


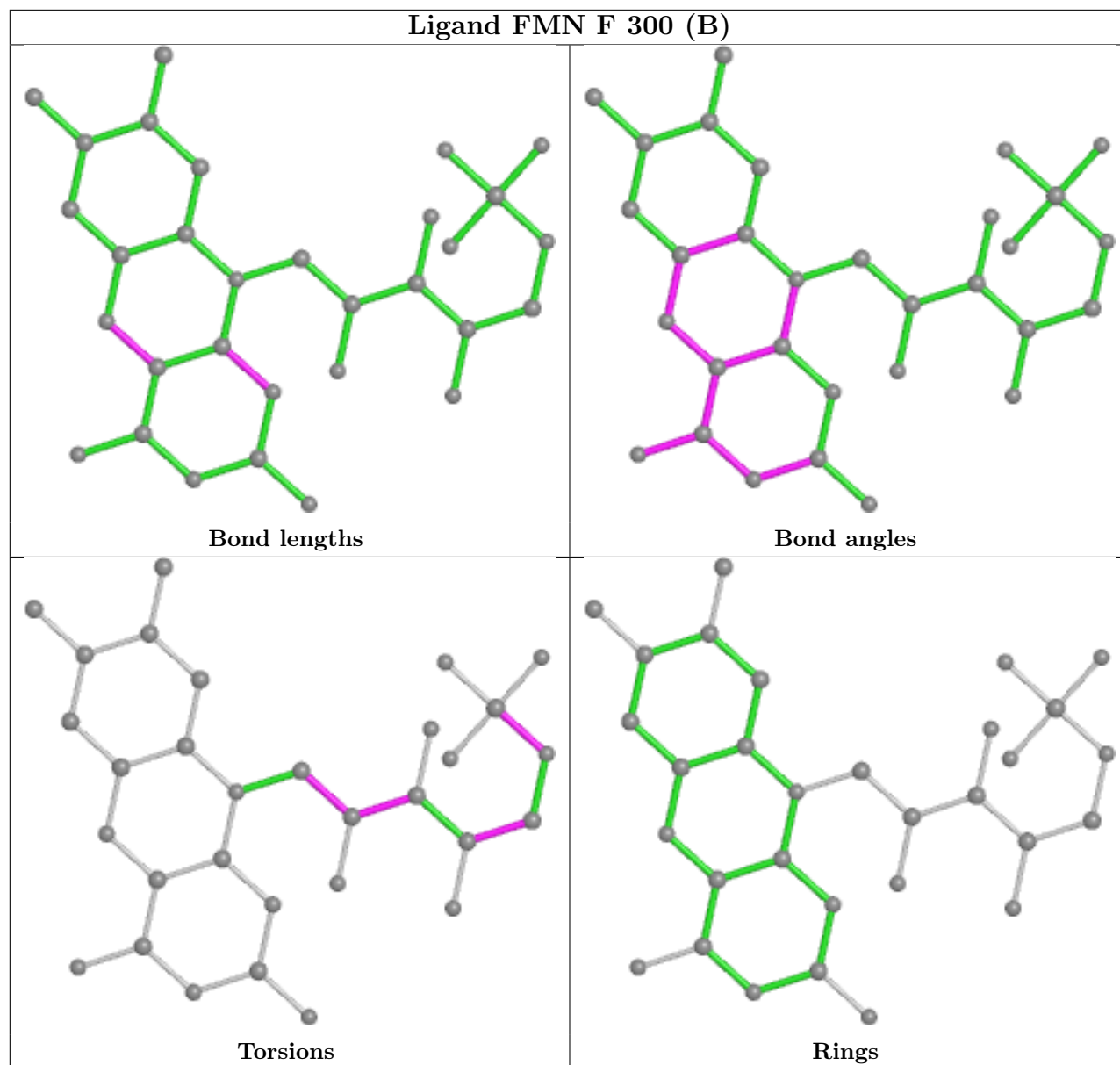


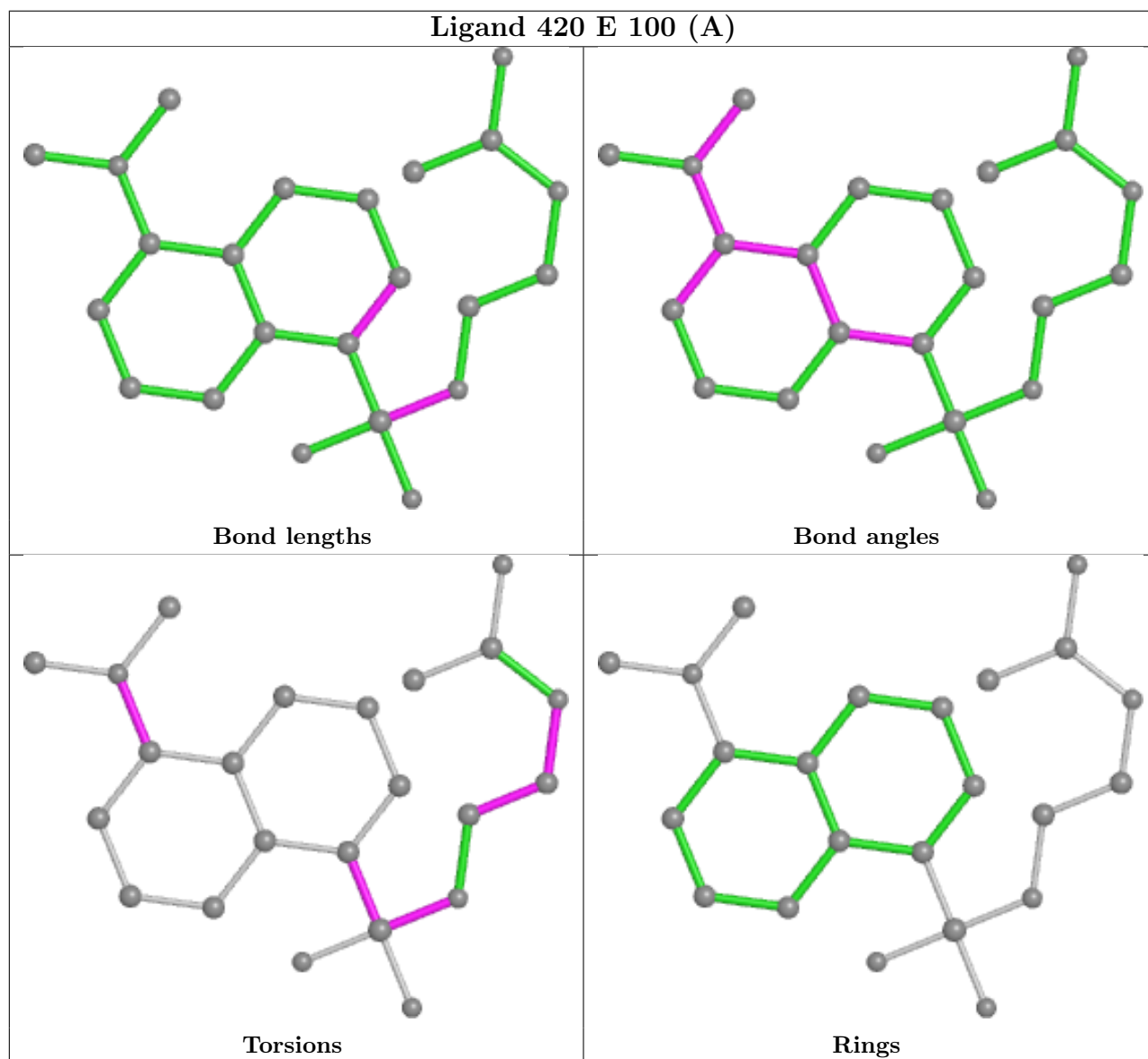
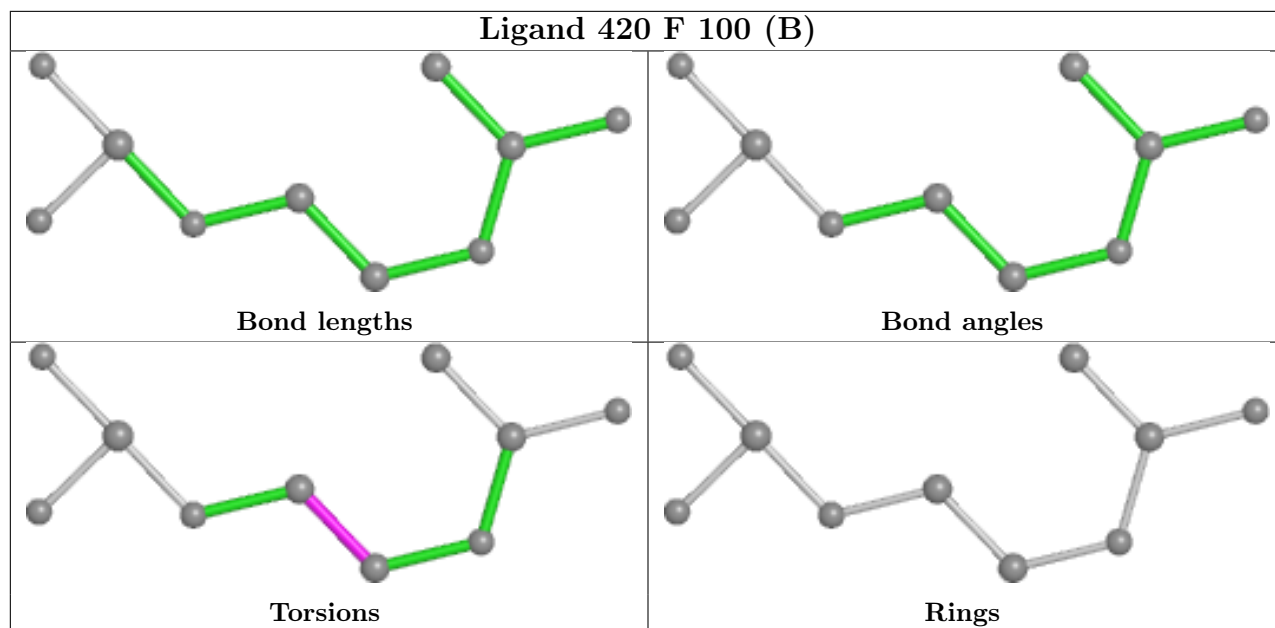


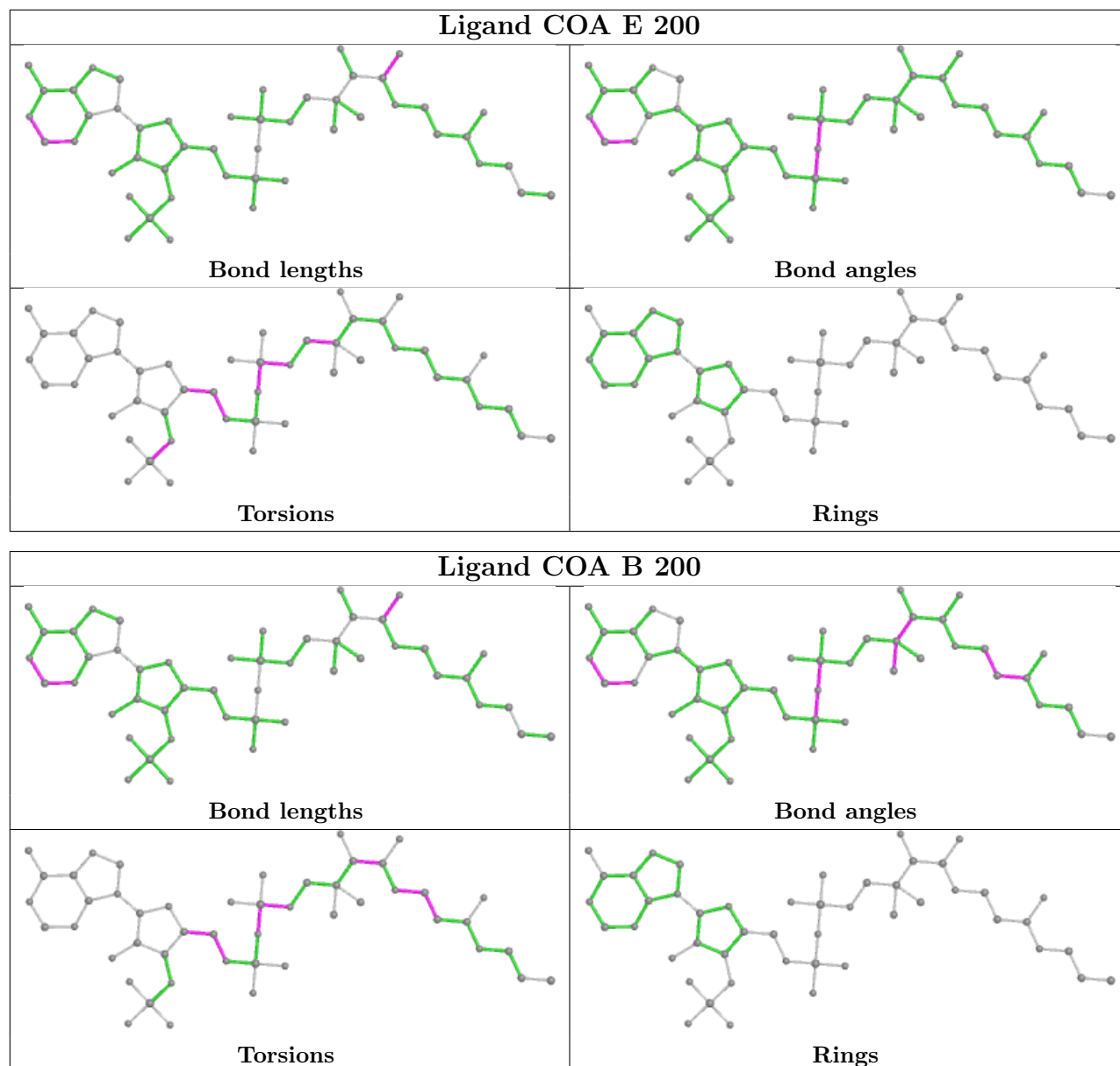


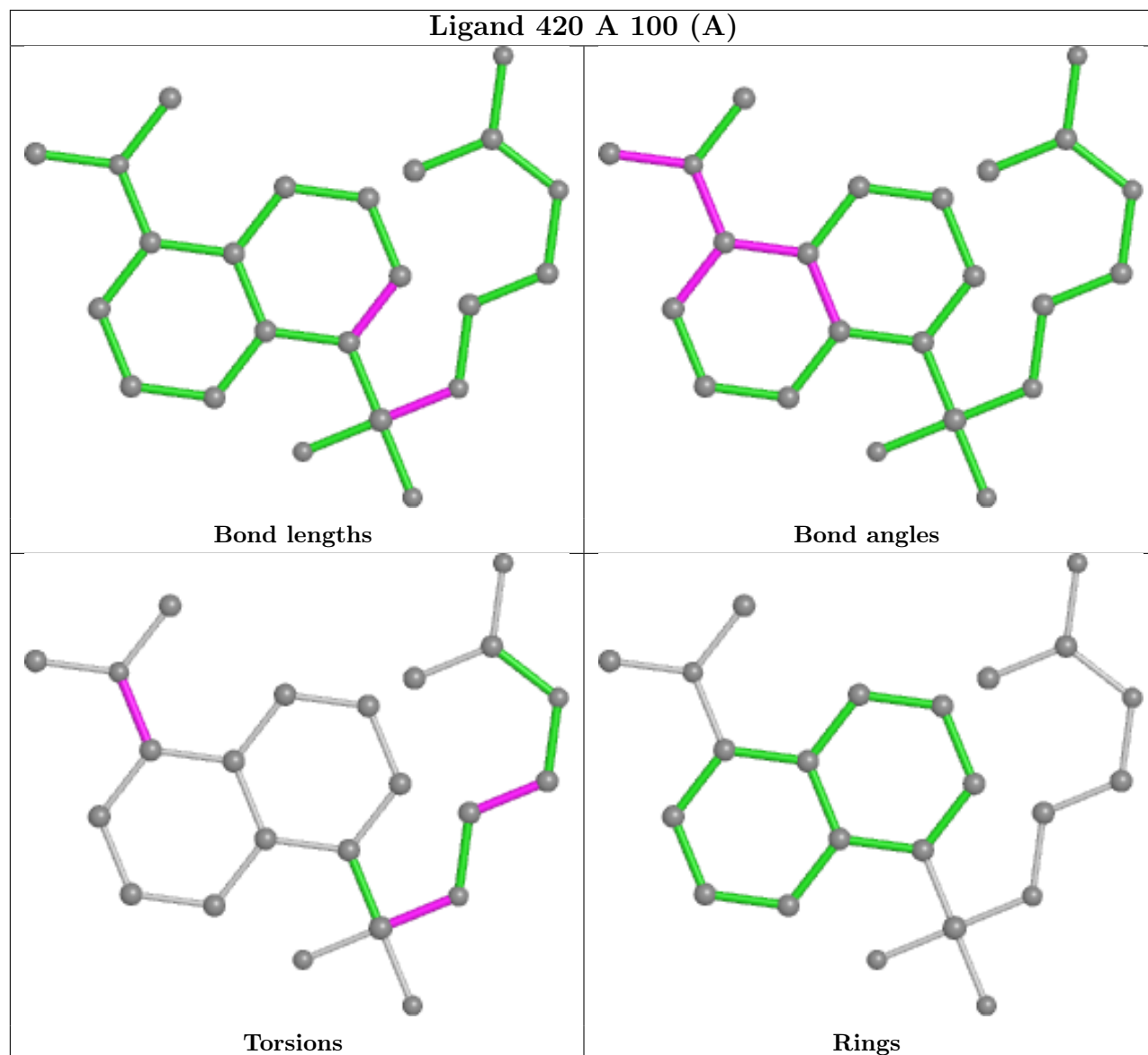


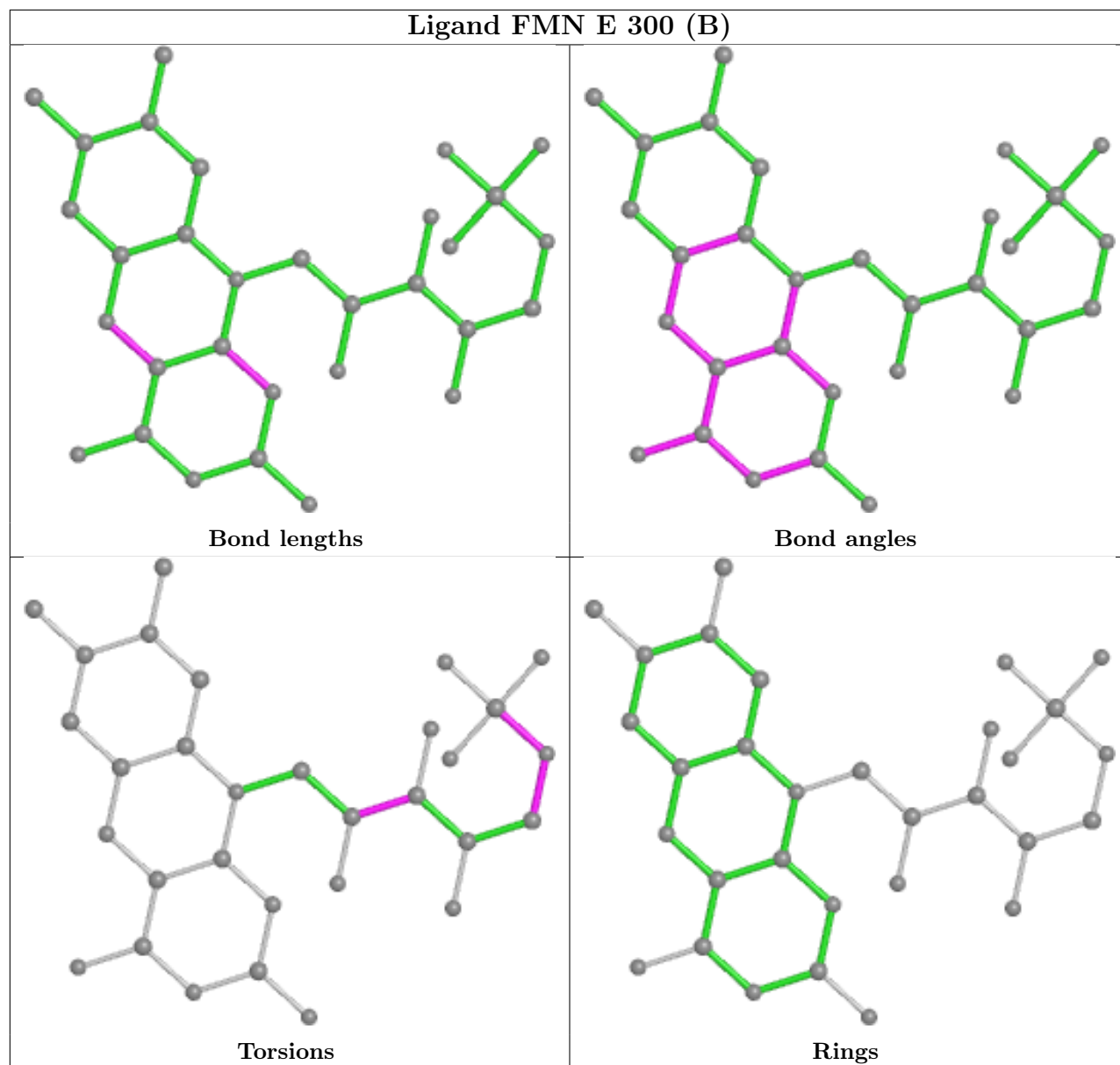




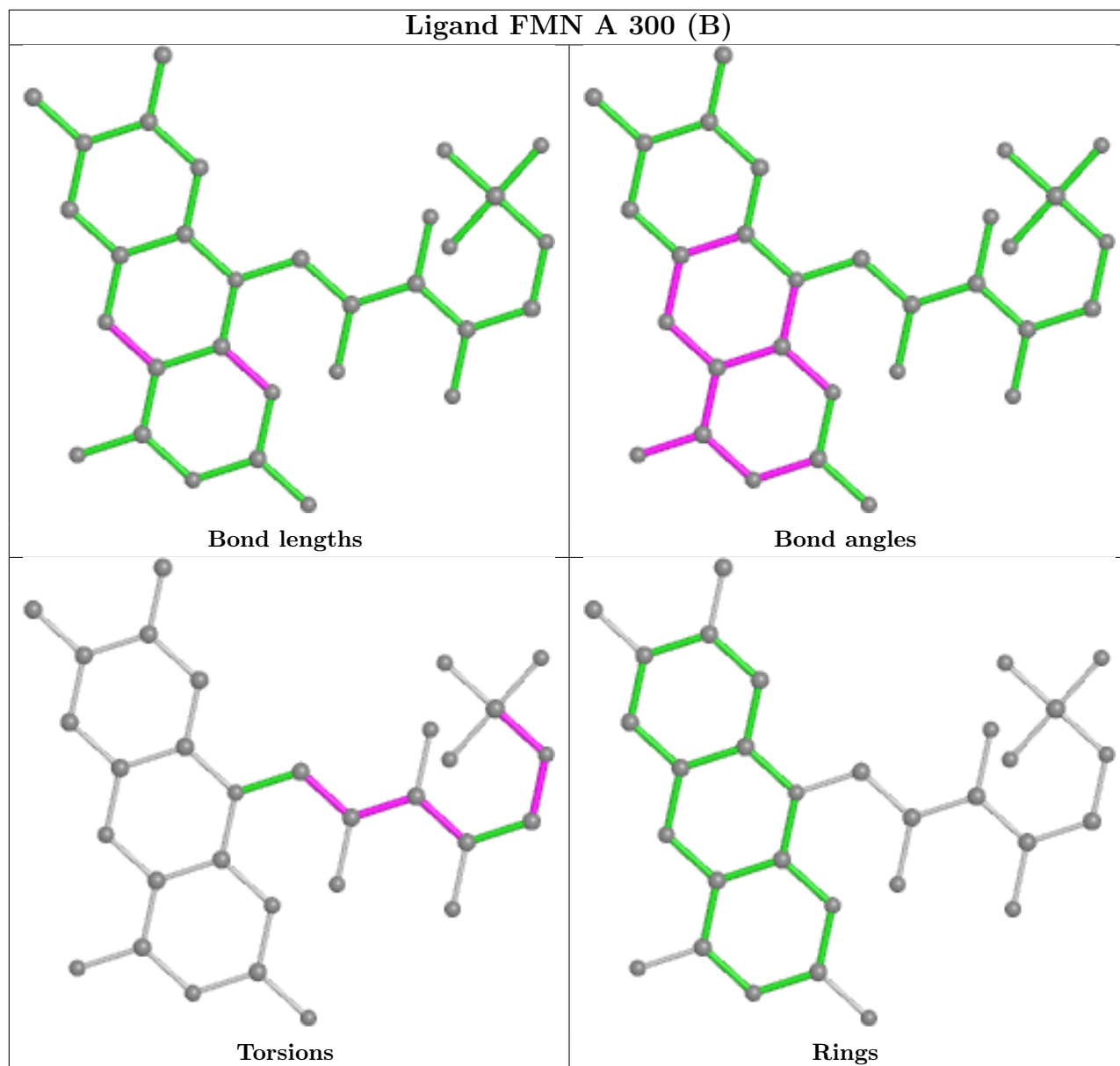


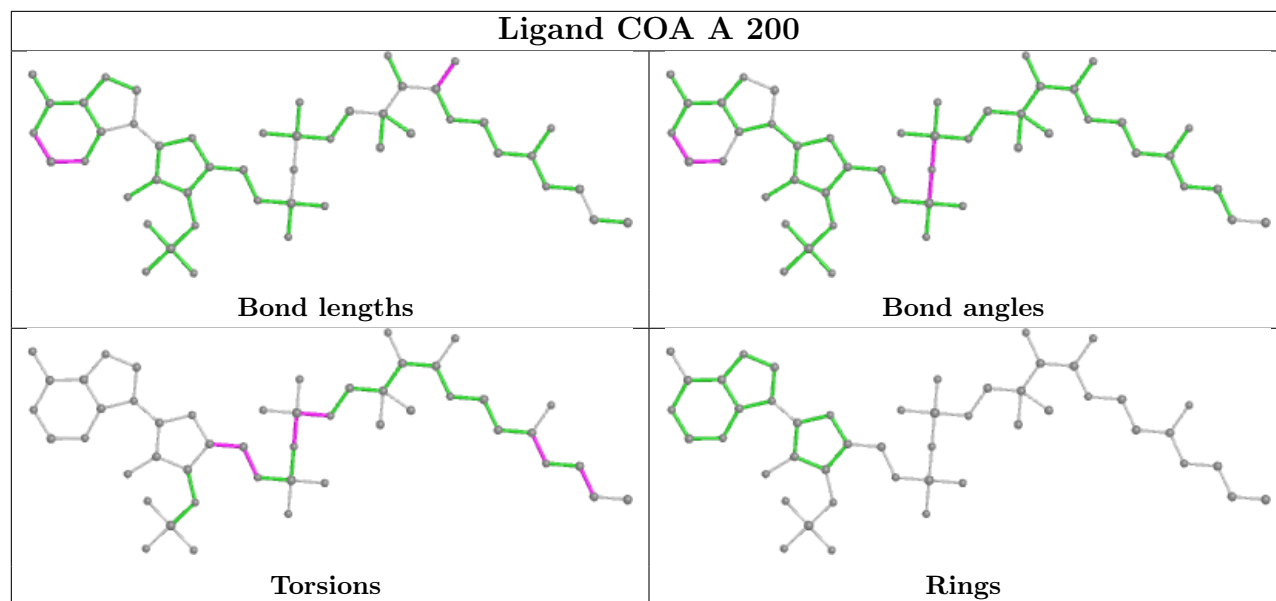


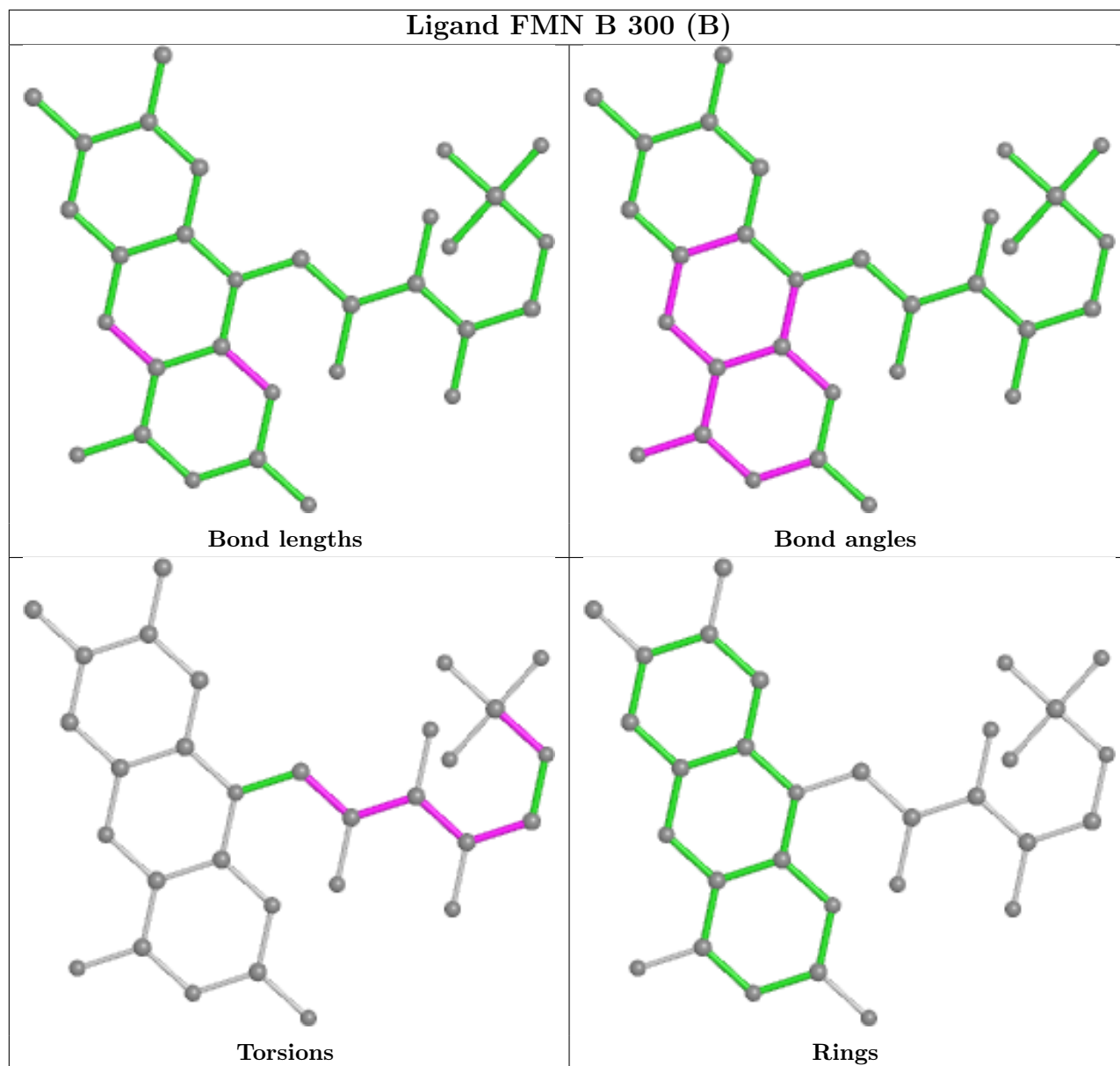


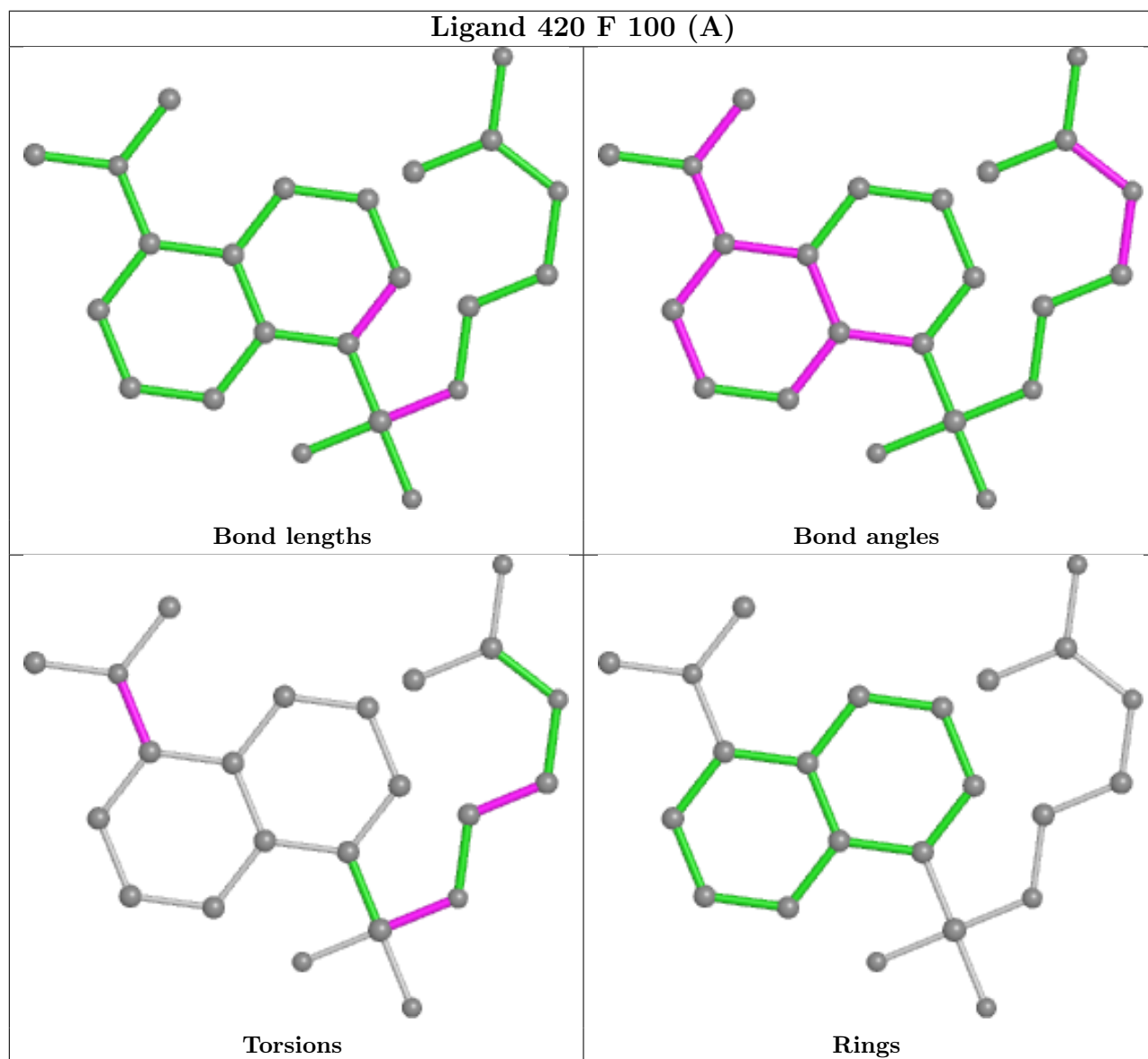
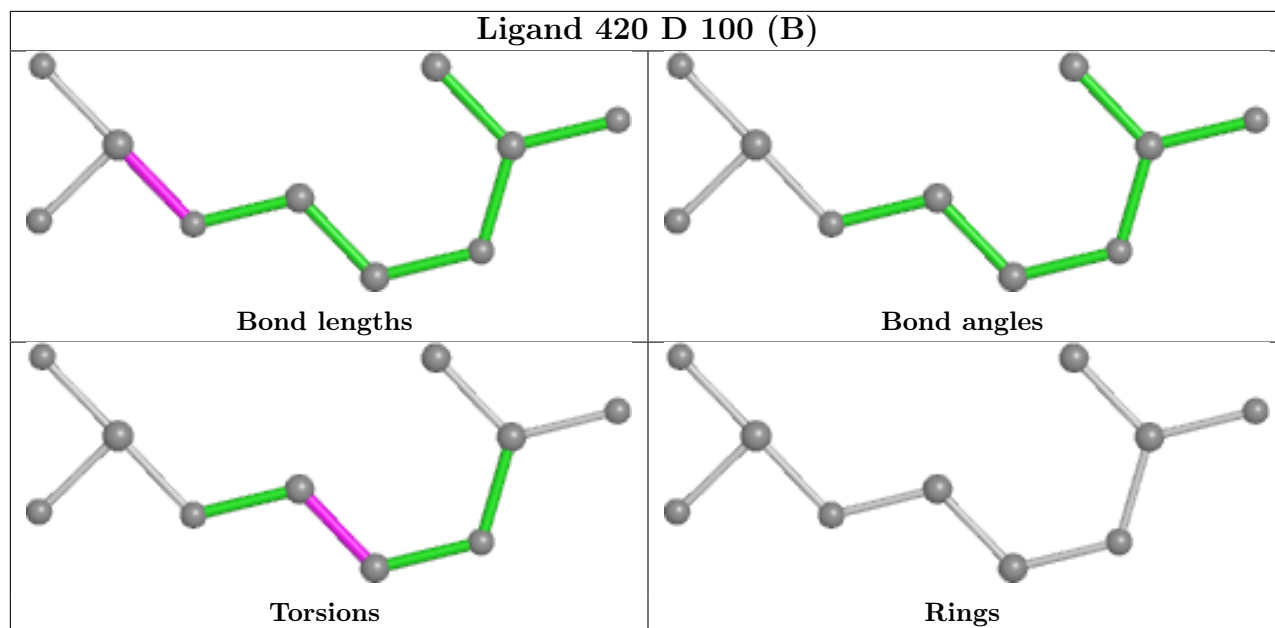


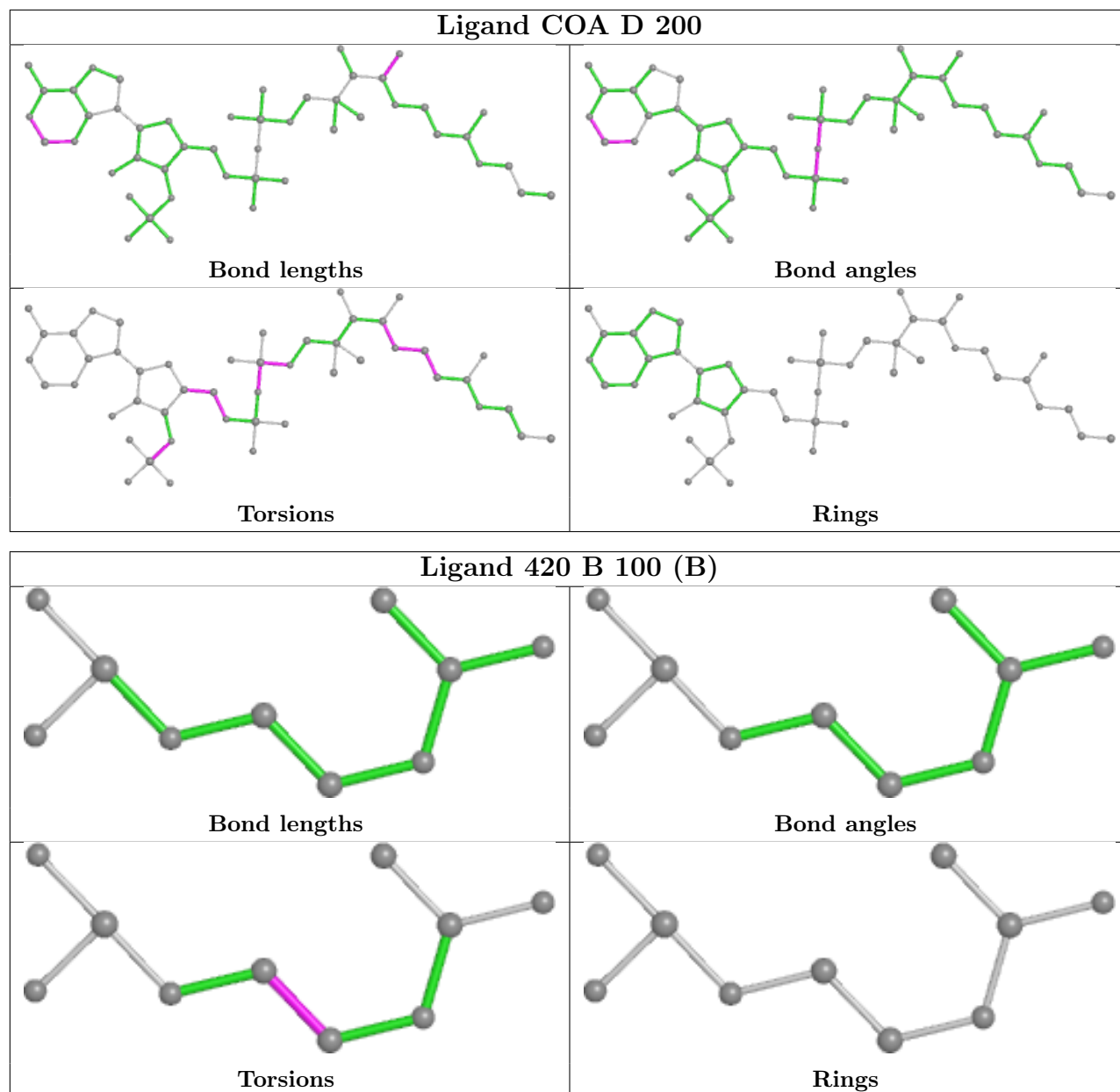


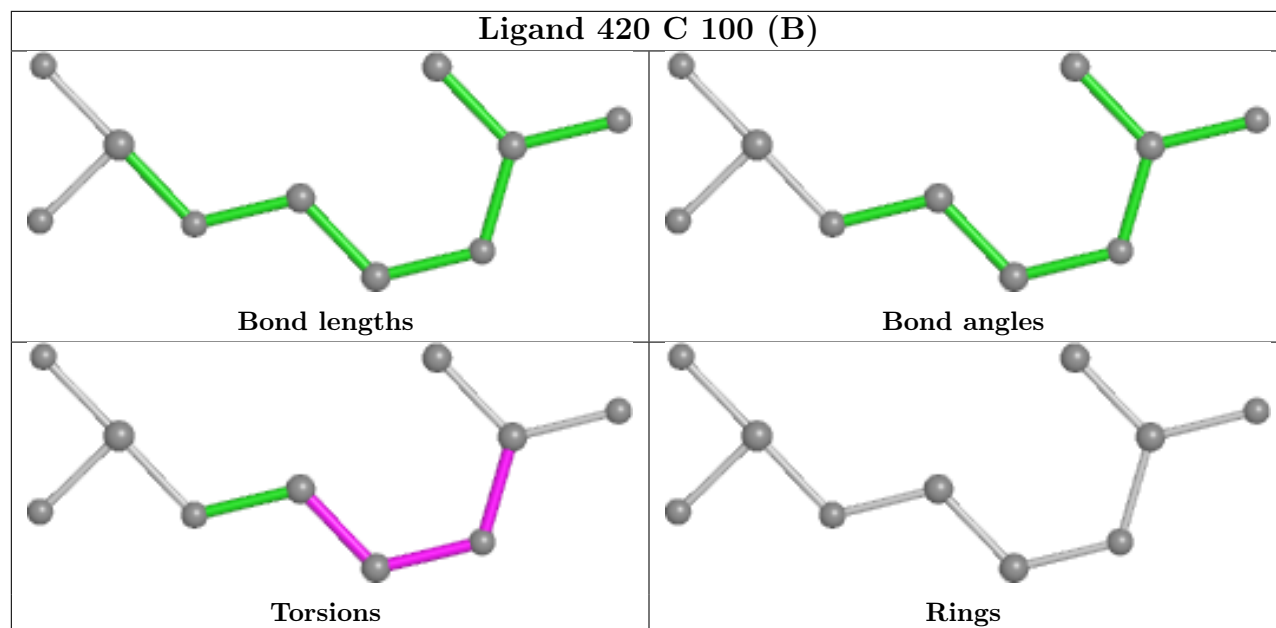












## 5.7 Other polymers [i](#)

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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	68/69 (98%)	-0.14	1 (1%) 73 72	18, 24, 39, 51	0
1	B	67/69 (97%)	-0.29	1 (1%) 73 72	18, 24, 38, 47	0
1	C	68/69 (98%)	-0.07	4 (5%) 22 21	20, 26, 38, 46	0
1	D	68/69 (98%)	-0.18	1 (1%) 73 72	18, 24, 35, 43	0
1	E	68/69 (98%)	-0.04	2 (2%) 51 50	19, 28, 45, 63	0
1	F	68/69 (98%)	-0.04	1 (1%) 73 72	19, 27, 44, 53	0
All	All	407/414 (98%)	-0.13	10 (2%) 57 55	18, 25, 42, 63	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	1	SER	6.4
1	A	1	SER	4.5
1	D	1	SER	4.2
1	C	1	SER	4.1
1	C	2	ASN	3.6
1	F	1	SER	3.5
1	E	3	HIS	2.5
1	C	50	ASP	2.4
1	B	2[A]	ASN	2.4
1	C	3	HIS	2.3

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

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

### 6.3 Carbohydrates [i](#)

There are no 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, 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(Å <sup>2</sup> )	Q<0.9
4	FMN	C	300[B]	31/31	0.66	0.35	23,25,26,26	31
4	FMN	A	300[B]	31/31	0.67	0.35	24,26,26,26	31
2	420	D	100[A]	23/24	0.72	0.35	20,31,32,32	23
2	420	D	100[B]	10/24	0.72	0.35	16,19,21,21	10
2	420	B	100[A]	23/24	0.74	0.36	18,30,32,33	23
4	FMN	B	300[B]	31/31	0.74	0.31	22,26,27,27	31
2	420	B	100[B]	10/24	0.74	0.36	16,20,22,22	10
4	FMN	E	300[B]	31/31	0.74	0.34	25,26,27,27	31
4	FMN	D	300[B]	31/31	0.76	0.33	25,26,26,27	31
4	FMN	F	300[B]	31/31	0.76	0.28	23,25,26,27	31
2	420	C	100[A]	23/24	0.78	0.37	19,32,33,33	23
2	420	C	100[B]	10/24	0.78	0.37	16,19,23,23	10
2	420	E	100[A]	23/24	0.79	0.32	19,31,32,32	23
2	420	E	100[B]	10/24	0.79	0.32	16,20,23,23	10
2	420	F	100[A]	23/24	0.81	0.30	19,29,29,30	23
2	420	F	100[B]	10/24	0.81	0.30	15,18,20,20	10
2	420	A	100[A]	23/24	0.83	0.28	18,27,29,30	23
2	420	A	100[B]	7/24	0.83	0.28	14,16,16,16	7
3	COA	C	200	48/48	0.93	0.14	23,27,30,31	0
3	COA	F	200	48/48	0.93	0.14	26,27,33,33	0
3	COA	A	200	48/48	0.94	0.13	24,27,31,32	0
3	COA	B	200	48/48	0.94	0.12	24,25,29,30	0
3	COA	E	200	48/48	0.95	0.12	24,26,31,32	0
5	CL	B	1071	1/1	0.95	0.05	33,33,33,33	0
5	CL	B	1070	1/1	0.96	0.09	38,38,38,38	0
3	COA	D	200	48/48	0.96	0.11	24,25,31,32	0
5	CL	A	1070	1/1	0.97	0.16	40,40,40,40	0
6	SO4	D	1070	5/5	0.98	0.13	34,34,34,34	0
6	SO4	E	1070	5/5	0.98	0.09	26,27,27,27	0

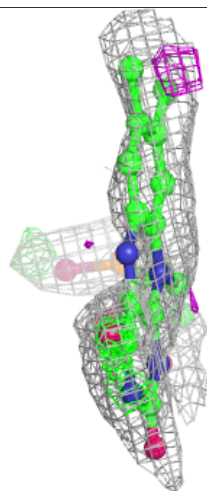
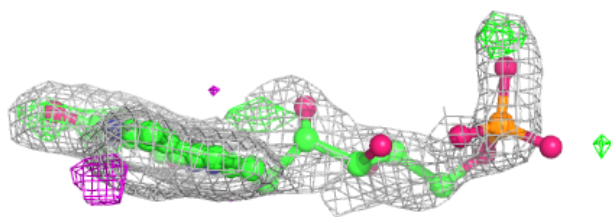
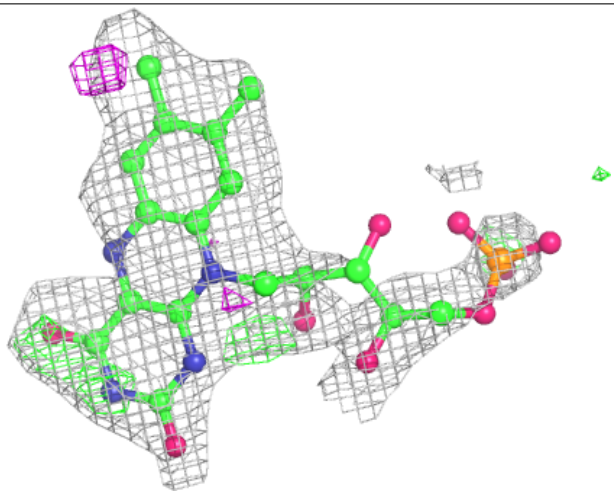
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers



as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

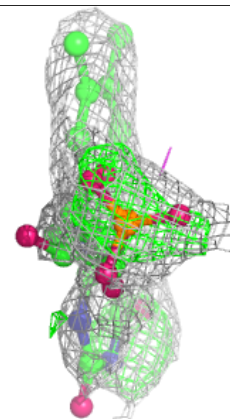
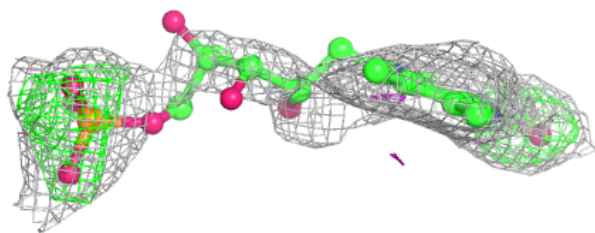
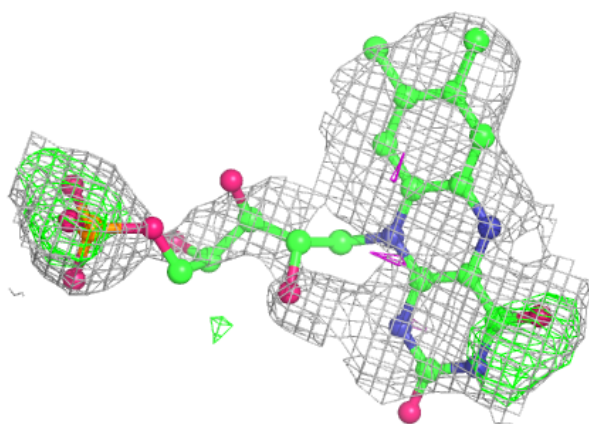
**Electron density around FMN C 300 (B):**

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

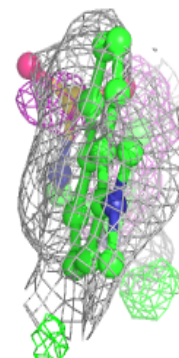
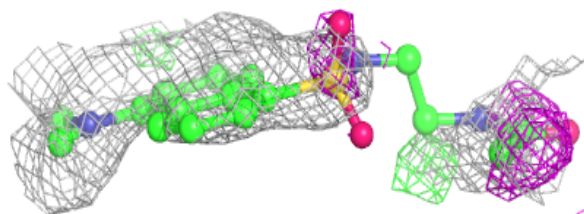
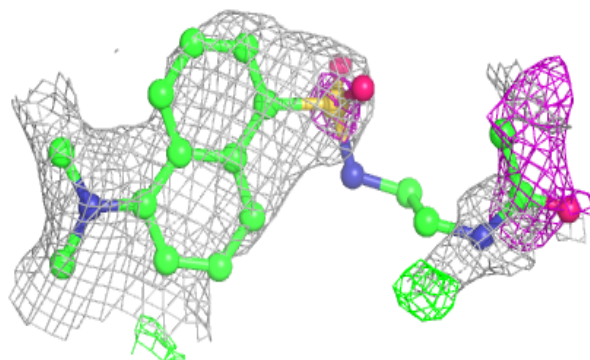


**Electron density around FMN A 300 (B):**

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

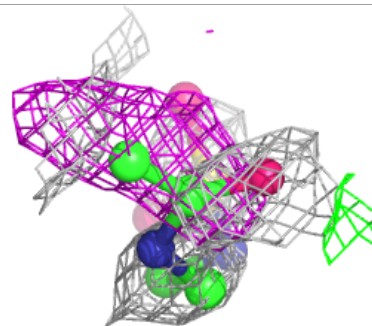
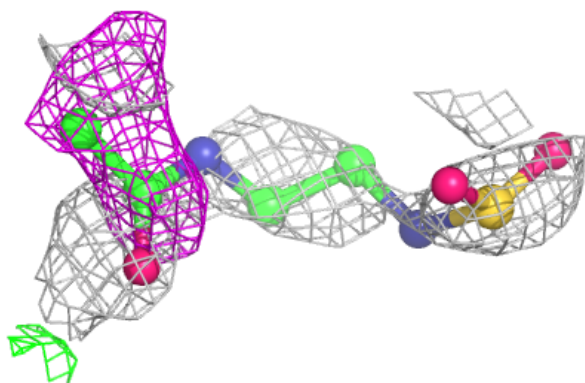
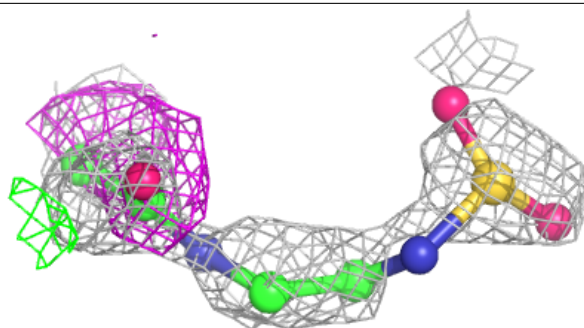
**Electron density around 420 D 100 (A):**

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

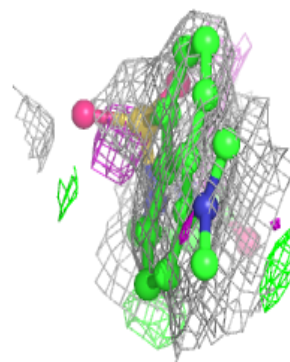
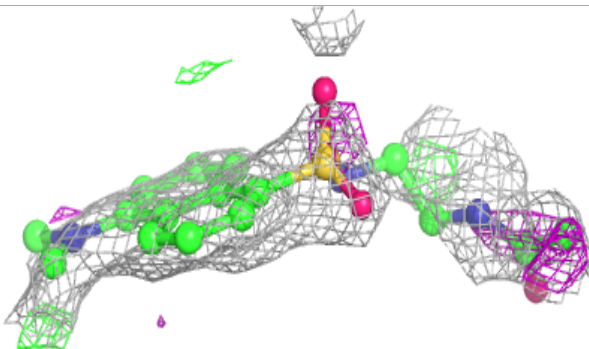
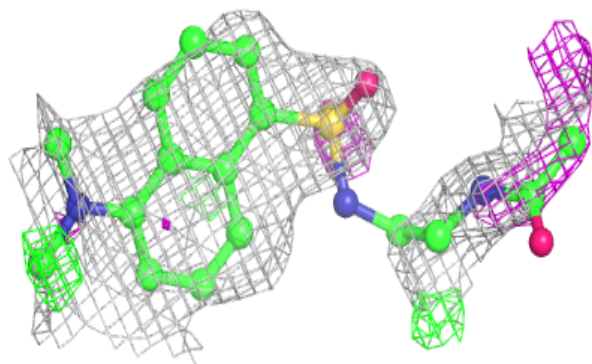


**Electron density around 420 D 100 (B):**

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

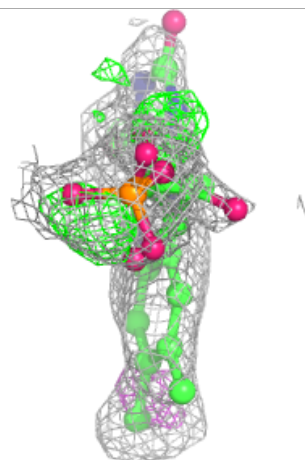
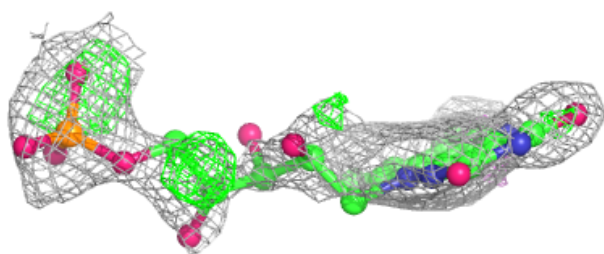
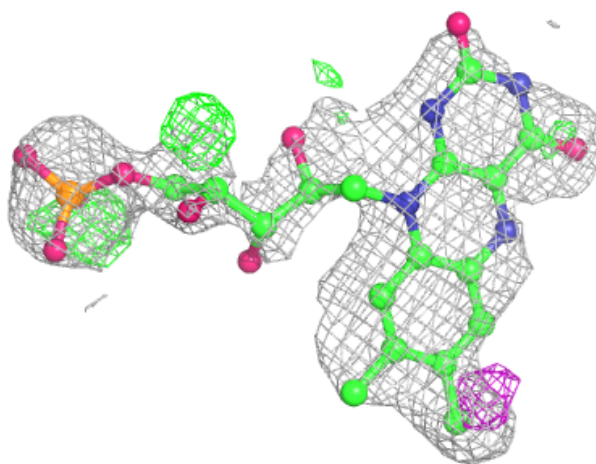
**Electron density around 420 B 100 (A):**

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



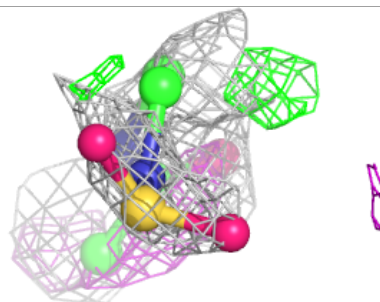
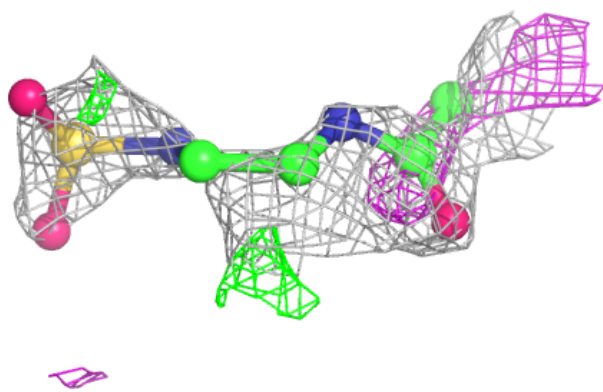
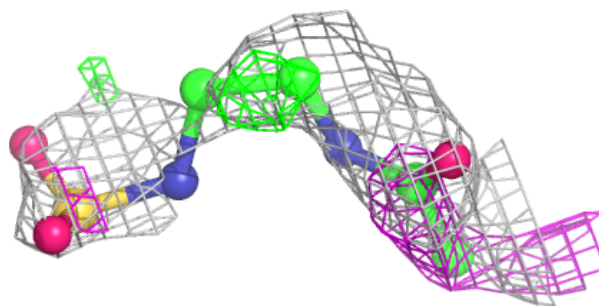
**Electron density around FMN B 300 (B):**

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



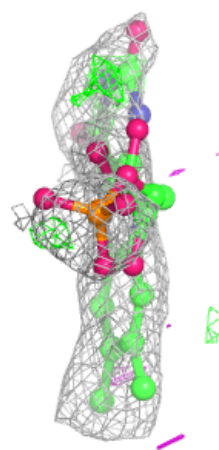
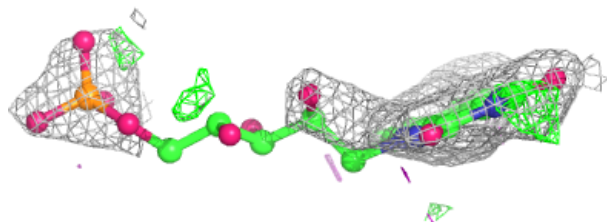
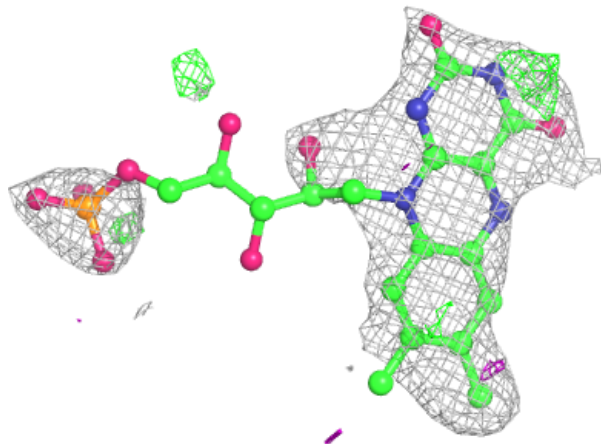
**Electron density around 420 B 100 (B):**

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



**Electron density around FMN E 300 (B):**

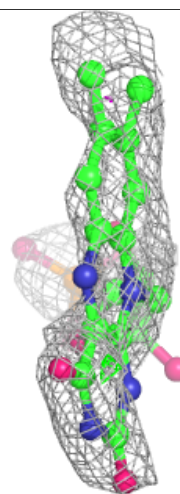
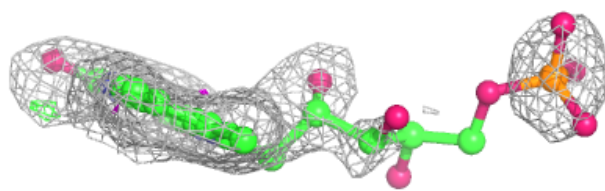
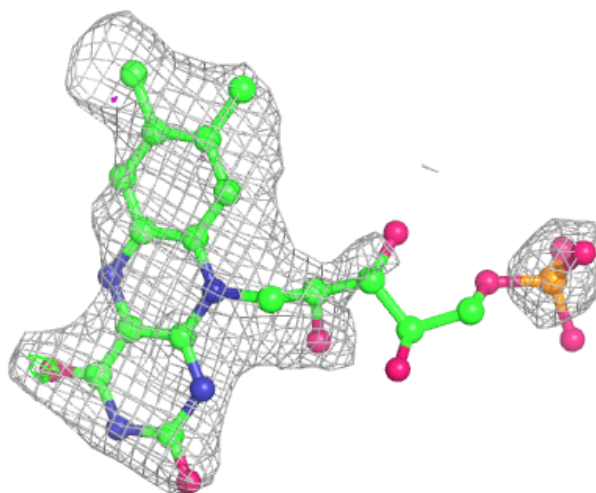
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





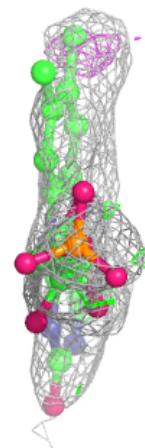
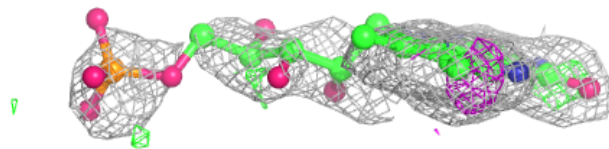
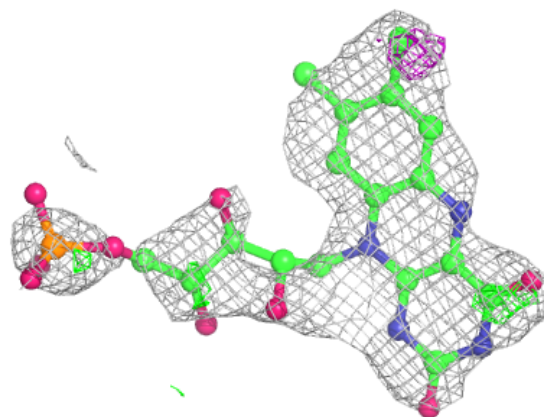
**Electron density around FMN D 300 (B):**

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



**Electron density around FMN F 300 (B):**

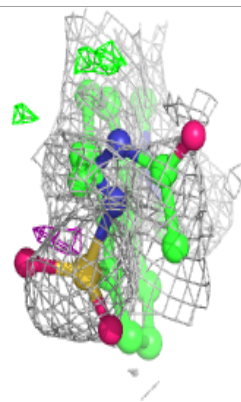
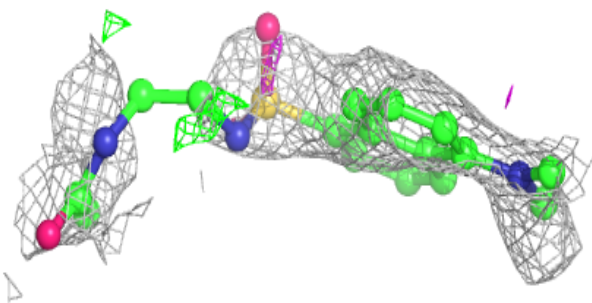
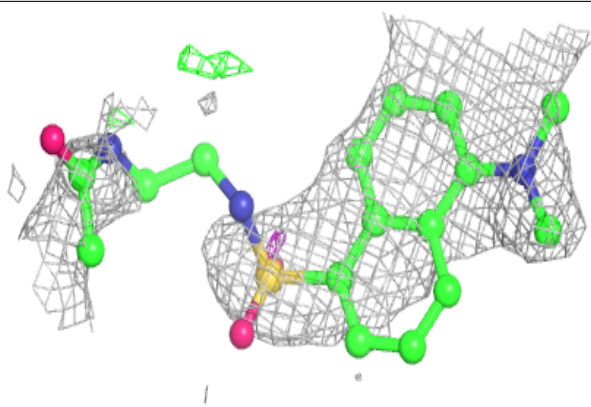
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



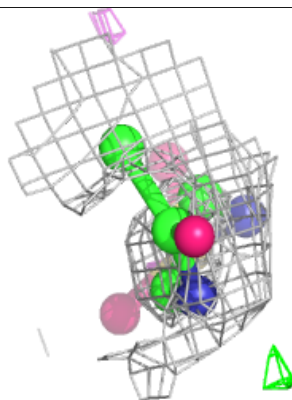
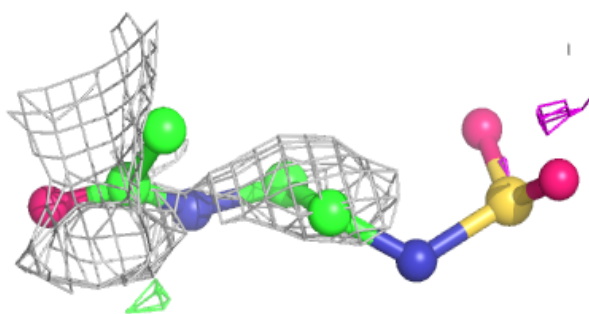
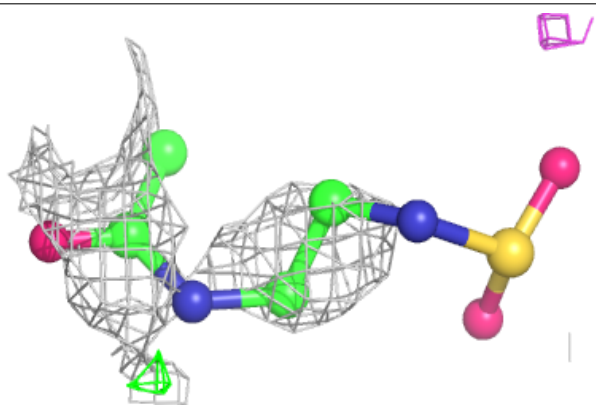


**Electron density around 420 C 100 (A):**

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

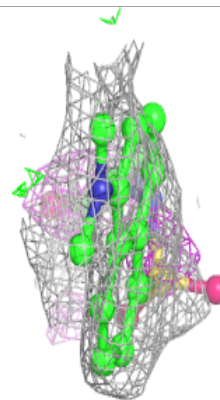
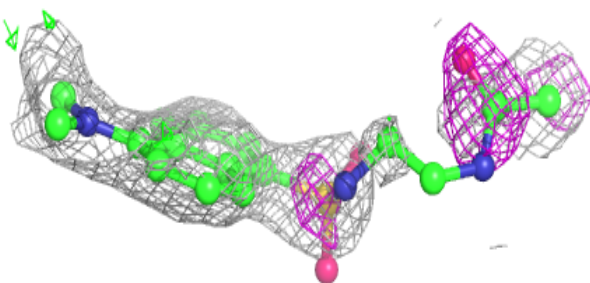
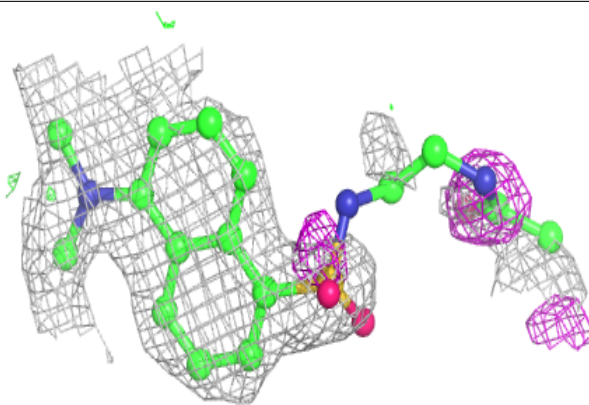
**Electron density around 420 C 100 (B):**

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

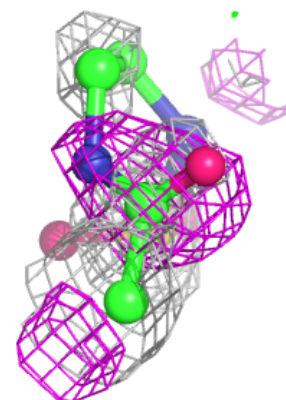
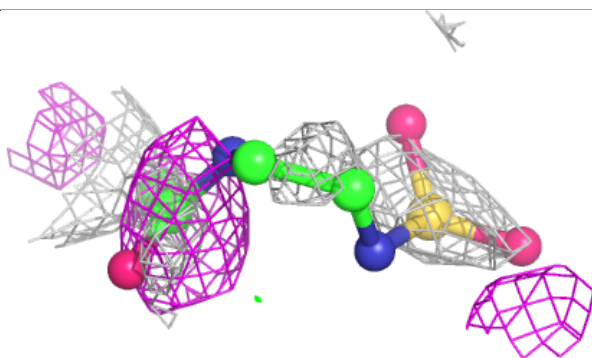
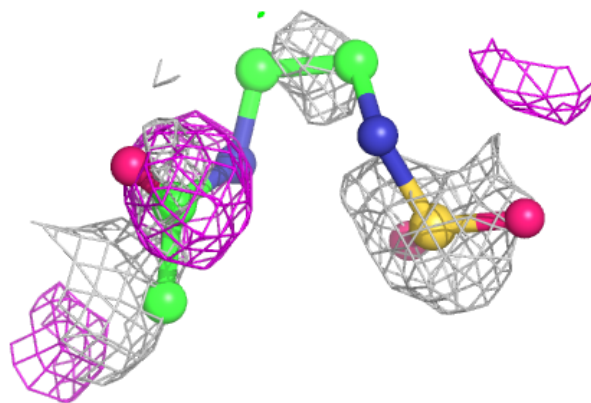


**Electron density around 420 E 100 (A):**

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

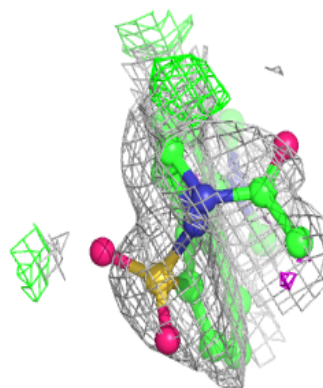
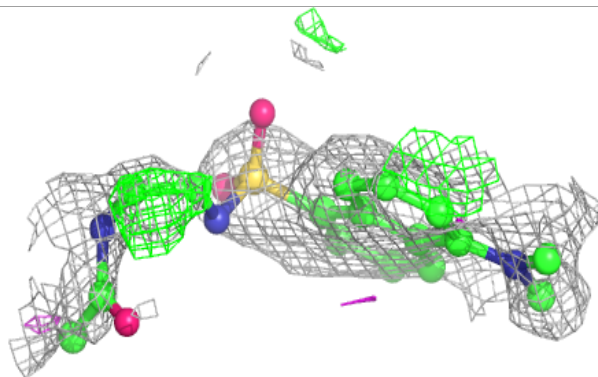
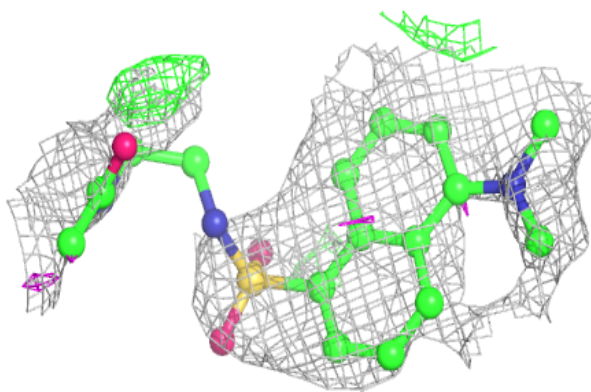
**Electron density around 420 E 100 (B):**

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

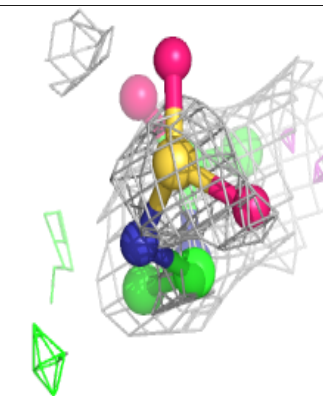
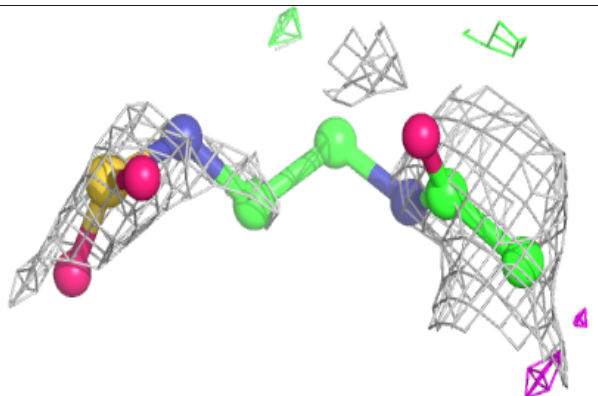
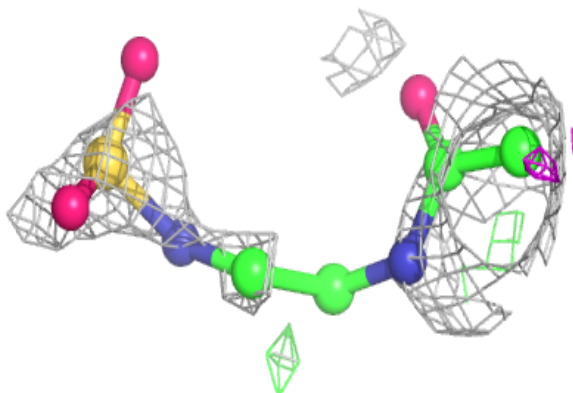


**Electron density around 420 F 100 (A):**

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

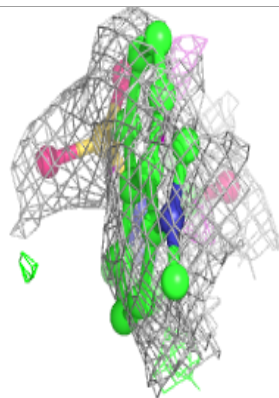
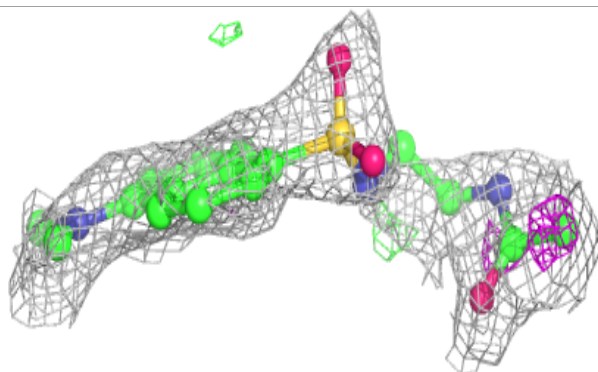
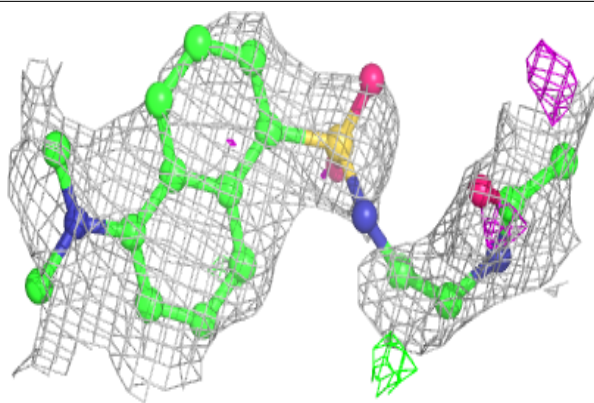
**Electron density around 420 F 100 (B):**

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

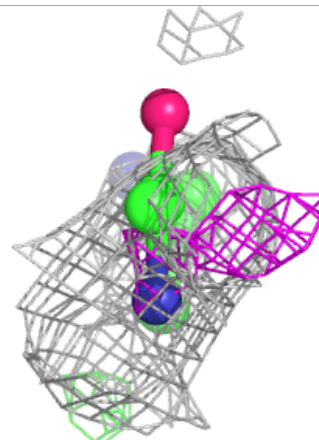
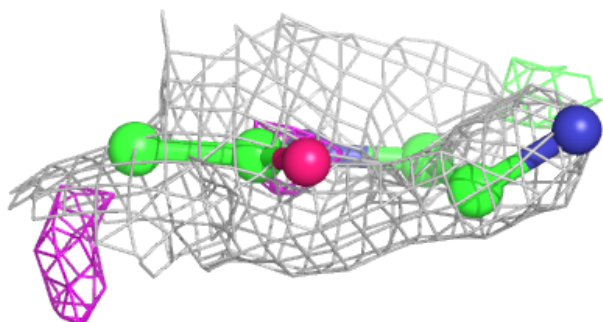
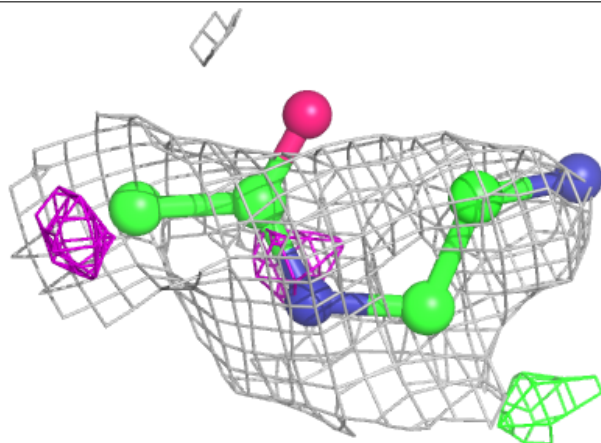


**Electron density around 420 A 100 (A):**

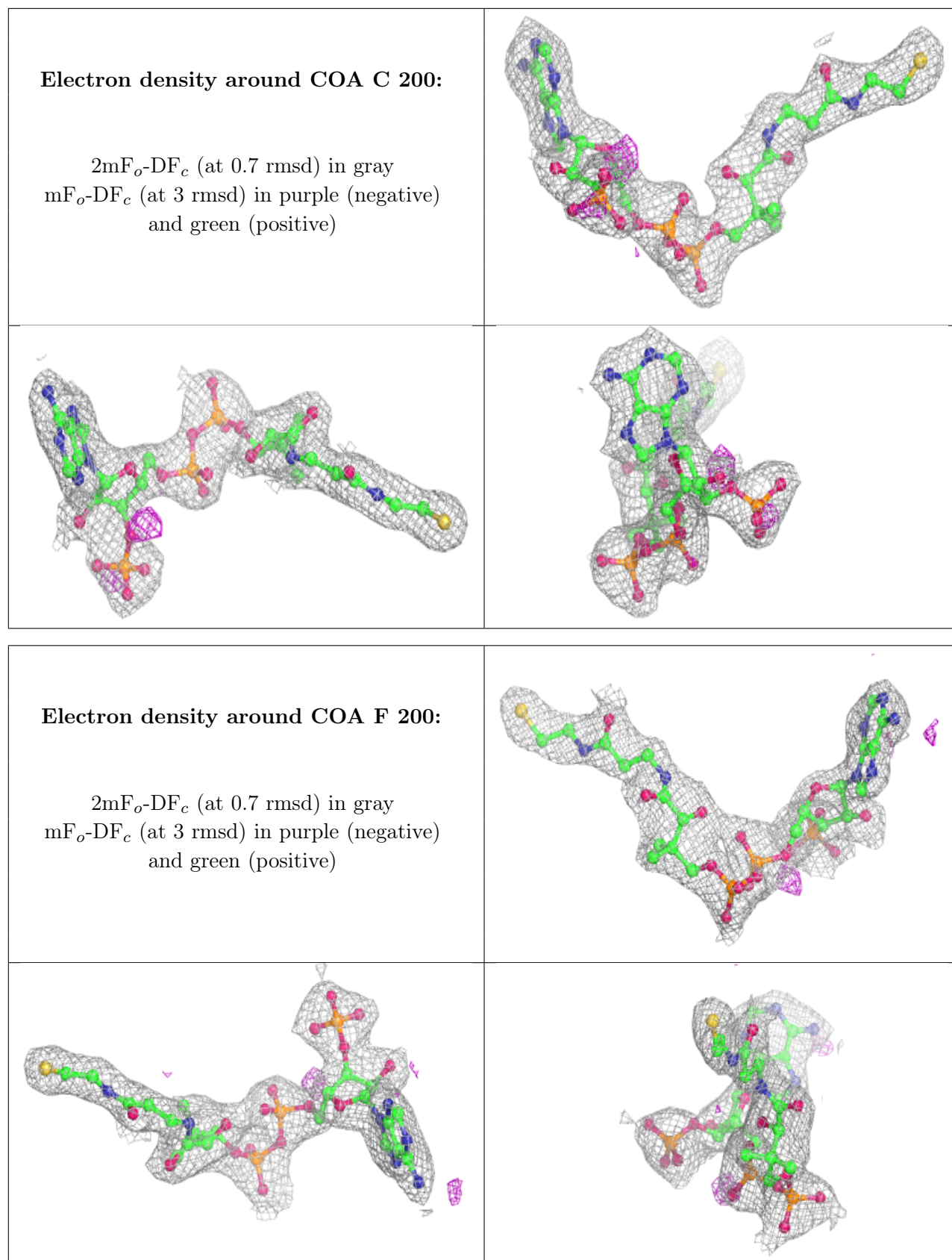
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

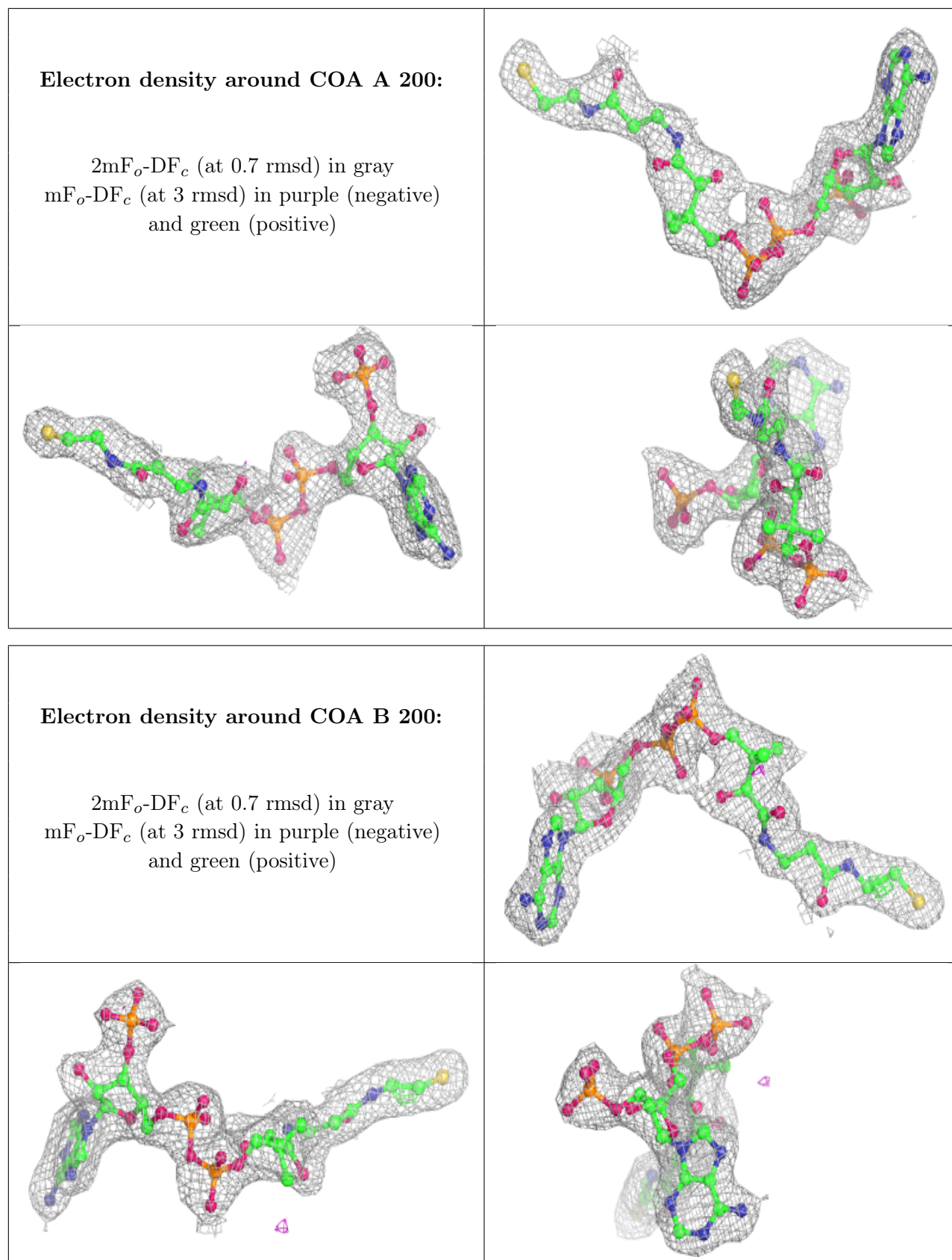
**Electron density around 420 A 100 (B):**

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



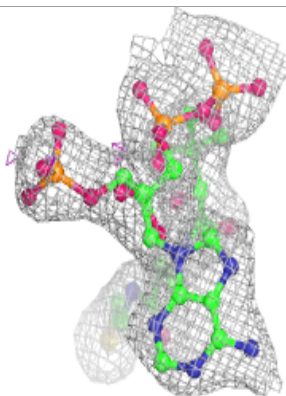
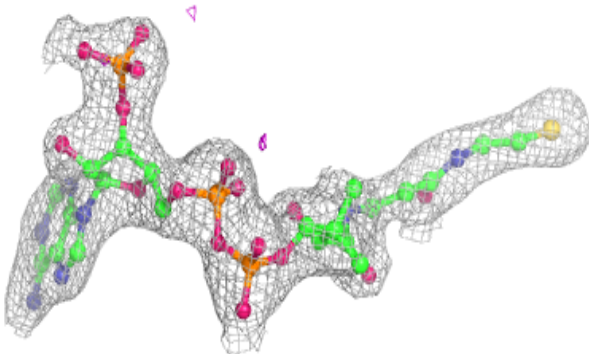
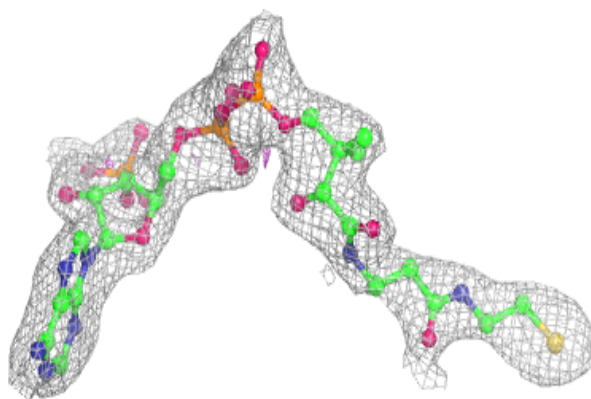




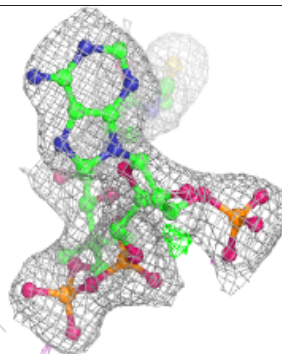
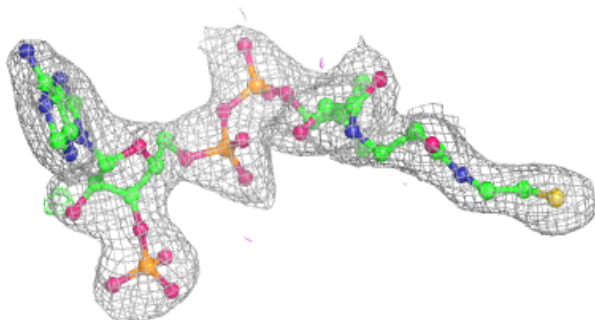
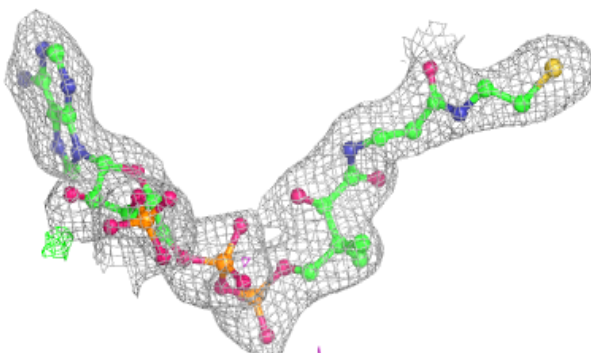


**Electron density around COA E 200:**

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

**Electron density around COA D 200:**

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



## 6.5 Other polymers [i](#)

There are no such residues in this entry.