



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:41 PM UTC

PDB ID : 9HG2 / pdb_00009hg2
Title : Crystal structure of M. smegmatis GMP reductase in complex with IMP and ATP.
Authors : Dolezal, M.; Pichova, I.
Deposited on : 2024-11-18
Resolution : 1.60 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 types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

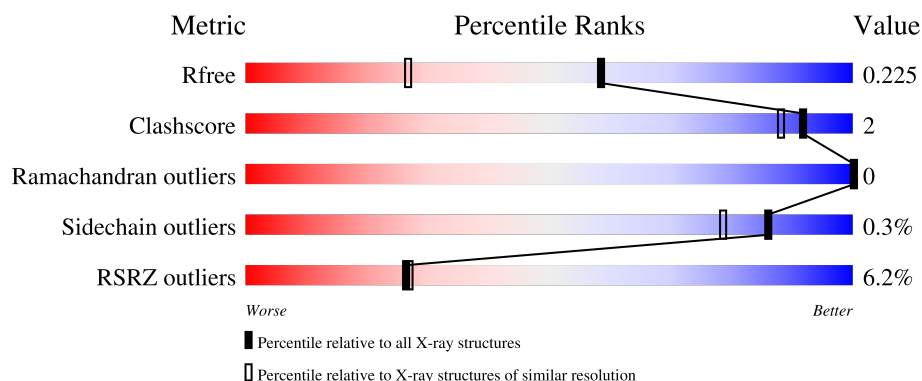
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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.



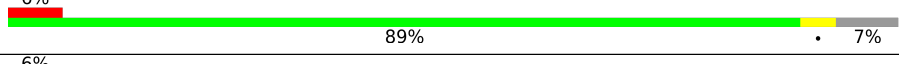
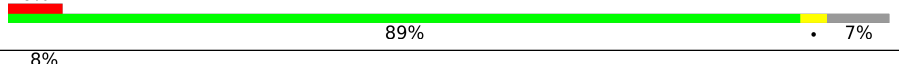
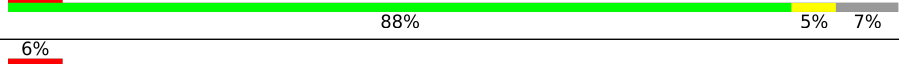
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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 ≥ 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	496	<div> <div>4%</div> <div>89%</div> <div>7%</div> </div>
1	B	496	<div> <div>5%</div> <div>90%</div> <div>7%</div> </div>
1	C	496	<div> <div>5%</div> <div>89%</div> <div>7%</div> </div>
1	D	496	<div> <div>4%</div> <div>89%</div> <div>7%</div> </div>
1	E	496	<div> <div>4%</div> <div>90%</div> <div>7%</div> </div>

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	496	
1	G	496	
1	H	496	
1	I	496	
1	J	496	
1	K	496	
1	L	496	
1	M	496	
1	N	496	
1	O	496	
1	P	496	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 57872 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 GMP reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	461	Total	C	N	O	S	0	4	0
			3359	2096	602	646	15			
1	B	460	Total	C	N	O	S	0	3	0
			3339	2083	596	646	14			
1	C	461	Total	C	N	O	S	0	4	0
			3360	2095	598	652	15			
1	D	460	Total	C	N	O	S	0	3	0
			3348	2087	597	650	14			
1	E	460	Total	C	N	O	S	0	4	0
			3352	2090	601	647	14			
1	F	461	Total	C	N	O	S	0	3	0
			3356	2092	598	651	15			
1	G	461	Total	C	N	O	S	0	3	0
			3341	2085	594	647	15			
1	H	461	Total	C	N	O	S	0	3	0
			3348	2089	597	647	15			
1	I	461	Total	C	N	O	S	0	3	0
			3345	2086	597	647	15			
1	J	461	Total	C	N	O	S	0	3	0
			3348	2090	598	645	15			
1	K	460	Total	C	N	O	S	0	3	0
			3342	2086	596	646	14			
1	L	461	Total	C	N	O	S	0	3	0
			3341	2085	594	647	15			
1	M	461	Total	C	N	O	S	0	3	0
			3339	2084	594	646	15			
1	N	460	Total	C	N	O	S	0	3	0
			3336	2082	596	644	14			
1	O	461	Total	C	N	O	S	0	3	0
			3351	2090	597	649	15			
1	P	460	Total	C	N	O	S	0	3	0
			3334	2080	594	646	14			

There are 288 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	VAL	-	insertion	UNP A0QYE8
A	480	THR	-	expression tag	UNP A0QYE8
A	481	ALA	-	expression tag	UNP A0QYE8
A	482	ALA	-	expression tag	UNP A0QYE8
A	483	ALA	-	expression tag	UNP A0QYE8
A	484	LYS	-	expression tag	UNP A0QYE8
A	485	GLU	-	expression tag	UNP A0QYE8
A	486	ASP	-	expression tag	UNP A0QYE8
A	487	LEU	-	expression tag	UNP A0QYE8
A	488	GLU	-	expression tag	UNP A0QYE8
A	489	HIS	-	expression tag	UNP A0QYE8
A	490	HIS	-	expression tag	UNP A0QYE8
A	491	HIS	-	expression tag	UNP A0QYE8
A	492	HIS	-	expression tag	UNP A0QYE8
A	493	HIS	-	expression tag	UNP A0QYE8
A	494	HIS	-	expression tag	UNP A0QYE8
A	495	HIS	-	expression tag	UNP A0QYE8
A	496	HIS	-	expression tag	UNP A0QYE8
B	2	VAL	-	insertion	UNP A0QYE8
B	480	THR	-	expression tag	UNP A0QYE8
B	481	ALA	-	expression tag	UNP A0QYE8
B	482	ALA	-	expression tag	UNP A0QYE8
B	483	ALA	-	expression tag	UNP A0QYE8
B	484	LYS	-	expression tag	UNP A0QYE8
B	485	GLU	-	expression tag	UNP A0QYE8
B	486	ASP	-	expression tag	UNP A0QYE8
B	487	LEU	-	expression tag	UNP A0QYE8
B	488	GLU	-	expression tag	UNP A0QYE8
B	489	HIS	-	expression tag	UNP A0QYE8
B	490	HIS	-	expression tag	UNP A0QYE8
B	491	HIS	-	expression tag	UNP A0QYE8
B	492	HIS	-	expression tag	UNP A0QYE8
B	493	HIS	-	expression tag	UNP A0QYE8
B	494	HIS	-	expression tag	UNP A0QYE8
B	495	HIS	-	expression tag	UNP A0QYE8
B	496	HIS	-	expression tag	UNP A0QYE8
C	2	VAL	-	insertion	UNP A0QYE8
C	480	THR	-	expression tag	UNP A0QYE8
C	481	ALA	-	expression tag	UNP A0QYE8
C	482	ALA	-	expression tag	UNP A0QYE8
C	483	ALA	-	expression tag	UNP A0QYE8
C	484	LYS	-	expression tag	UNP A0QYE8

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	485	GLU	-	expression tag	UNP A0QYE8
C	486	ASP	-	expression tag	UNP A0QYE8
C	487	LEU	-	expression tag	UNP A0QYE8
C	488	GLU	-	expression tag	UNP A0QYE8
C	489	HIS	-	expression tag	UNP A0QYE8
C	490	HIS	-	expression tag	UNP A0QYE8
C	491	HIS	-	expression tag	UNP A0QYE8
C	492	HIS	-	expression tag	UNP A0QYE8
C	493	HIS	-	expression tag	UNP A0QYE8
C	494	HIS	-	expression tag	UNP A0QYE8
C	495	HIS	-	expression tag	UNP A0QYE8
C	496	HIS	-	expression tag	UNP A0QYE8
D	2	VAL	-	insertion	UNP A0QYE8
D	480	THR	-	expression tag	UNP A0QYE8
D	481	ALA	-	expression tag	UNP A0QYE8
D	482	ALA	-	expression tag	UNP A0QYE8
D	483	ALA	-	expression tag	UNP A0QYE8
D	484	LYS	-	expression tag	UNP A0QYE8
D	485	GLU	-	expression tag	UNP A0QYE8
D	486	ASP	-	expression tag	UNP A0QYE8
D	487	LEU	-	expression tag	UNP A0QYE8
D	488	GLU	-	expression tag	UNP A0QYE8
D	489	HIS	-	expression tag	UNP A0QYE8
D	490	HIS	-	expression tag	UNP A0QYE8
D	491	HIS	-	expression tag	UNP A0QYE8
D	492	HIS	-	expression tag	UNP A0QYE8
D	493	HIS	-	expression tag	UNP A0QYE8
D	494	HIS	-	expression tag	UNP A0QYE8
D	495	HIS	-	expression tag	UNP A0QYE8
D	496	HIS	-	expression tag	UNP A0QYE8
E	2	VAL	-	insertion	UNP A0QYE8
E	480	THR	-	expression tag	UNP A0QYE8
E	481	ALA	-	expression tag	UNP A0QYE8
E	482	ALA	-	expression tag	UNP A0QYE8
E	483	ALA	-	expression tag	UNP A0QYE8
E	484	LYS	-	expression tag	UNP A0QYE8
E	485	GLU	-	expression tag	UNP A0QYE8
E	486	ASP	-	expression tag	UNP A0QYE8
E	487	LEU	-	expression tag	UNP A0QYE8
E	488	GLU	-	expression tag	UNP A0QYE8
E	489	HIS	-	expression tag	UNP A0QYE8
E	490	HIS	-	expression tag	UNP A0QYE8

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
E	491	HIS	-	expression tag	UNP A0QYE8
E	492	HIS	-	expression tag	UNP A0QYE8
E	493	HIS	-	expression tag	UNP A0QYE8
E	494	HIS	-	expression tag	UNP A0QYE8
E	495	HIS	-	expression tag	UNP A0QYE8
E	496	HIS	-	expression tag	UNP A0QYE8
F	2	VAL	-	insertion	UNP A0QYE8
F	480	THR	-	expression tag	UNP A0QYE8
F	481	ALA	-	expression tag	UNP A0QYE8
F	482	ALA	-	expression tag	UNP A0QYE8
F	483	ALA	-	expression tag	UNP A0QYE8
F	484	LYS	-	expression tag	UNP A0QYE8
F	485	GLU	-	expression tag	UNP A0QYE8
F	486	ASP	-	expression tag	UNP A0QYE8
F	487	LEU	-	expression tag	UNP A0QYE8
F	488	GLU	-	expression tag	UNP A0QYE8
F	489	HIS	-	expression tag	UNP A0QYE8
F	490	HIS	-	expression tag	UNP A0QYE8
F	491	HIS	-	expression tag	UNP A0QYE8
F	492	HIS	-	expression tag	UNP A0QYE8
F	493	HIS	-	expression tag	UNP A0QYE8
F	494	HIS	-	expression tag	UNP A0QYE8
F	495	HIS	-	expression tag	UNP A0QYE8
F	496	HIS	-	expression tag	UNP A0QYE8
G	2	VAL	-	insertion	UNP A0QYE8
G	480	THR	-	expression tag	UNP A0QYE8
G	481	ALA	-	expression tag	UNP A0QYE8
G	482	ALA	-	expression tag	UNP A0QYE8
G	483	ALA	-	expression tag	UNP A0QYE8
G	484	LYS	-	expression tag	UNP A0QYE8
G	485	GLU	-	expression tag	UNP A0QYE8
G	486	ASP	-	expression tag	UNP A0QYE8
G	487	LEU	-	expression tag	UNP A0QYE8
G	488	GLU	-	expression tag	UNP A0QYE8
G	489	HIS	-	expression tag	UNP A0QYE8
G	490	HIS	-	expression tag	UNP A0QYE8
G	491	HIS	-	expression tag	UNP A0QYE8
G	492	HIS	-	expression tag	UNP A0QYE8
G	493	HIS	-	expression tag	UNP A0QYE8
G	494	HIS	-	expression tag	UNP A0QYE8
G	495	HIS	-	expression tag	UNP A0QYE8
G	496	HIS	-	expression tag	UNP A0QYE8

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
H	2	VAL	-	insertion	UNP A0QYE8
H	480	THR	-	expression tag	UNP A0QYE8
H	481	ALA	-	expression tag	UNP A0QYE8
H	482	ALA	-	expression tag	UNP A0QYE8
H	483	ALA	-	expression tag	UNP A0QYE8
H	484	LYS	-	expression tag	UNP A0QYE8
H	485	GLU	-	expression tag	UNP A0QYE8
H	486	ASP	-	expression tag	UNP A0QYE8
H	487	LEU	-	expression tag	UNP A0QYE8
H	488	GLU	-	expression tag	UNP A0QYE8
H	489	HIS	-	expression tag	UNP A0QYE8
H	490	HIS	-	expression tag	UNP A0QYE8
H	491	HIS	-	expression tag	UNP A0QYE8
H	492	HIS	-	expression tag	UNP A0QYE8
H	493	HIS	-	expression tag	UNP A0QYE8
H	494	HIS	-	expression tag	UNP A0QYE8
H	495	HIS	-	expression tag	UNP A0QYE8
H	496	HIS	-	expression tag	UNP A0QYE8
I	2	VAL	-	insertion	UNP A0QYE8
I	480	THR	-	expression tag	UNP A0QYE8
I	481	ALA	-	expression tag	UNP A0QYE8
I	482	ALA	-	expression tag	UNP A0QYE8
I	483	ALA	-	expression tag	UNP A0QYE8
I	484	LYS	-	expression tag	UNP A0QYE8
I	485	GLU	-	expression tag	UNP A0QYE8
I	486	ASP	-	expression tag	UNP A0QYE8
I	487	LEU	-	expression tag	UNP A0QYE8
I	488	GLU	-	expression tag	UNP A0QYE8
I	489	HIS	-	expression tag	UNP A0QYE8
I	490	HIS	-	expression tag	UNP A0QYE8
I	491	HIS	-	expression tag	UNP A0QYE8
I	492	HIS	-	expression tag	UNP A0QYE8
I	493	HIS	-	expression tag	UNP A0QYE8
I	494	HIS	-	expression tag	UNP A0QYE8
I	495	HIS	-	expression tag	UNP A0QYE8
I	496	HIS	-	expression tag	UNP A0QYE8
J	2	VAL	-	insertion	UNP A0QYE8
J	480	THR	-	expression tag	UNP A0QYE8
J	481	ALA	-	expression tag	UNP A0QYE8
J	482	ALA	-	expression tag	UNP A0QYE8
J	483	ALA	-	expression tag	UNP A0QYE8
J	484	LYS	-	expression tag	UNP A0QYE8

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
J	485	GLU	-	expression tag	UNP A0QYE8
J	486	ASP	-	expression tag	UNP A0QYE8
J	487	LEU	-	expression tag	UNP A0QYE8
J	488	GLU	-	expression tag	UNP A0QYE8
J	489	HIS	-	expression tag	UNP A0QYE8
J	490	HIS	-	expression tag	UNP A0QYE8
J	491	HIS	-	expression tag	UNP A0QYE8
J	492	HIS	-	expression tag	UNP A0QYE8
J	493	HIS	-	expression tag	UNP A0QYE8
J	494	HIS	-	expression tag	UNP A0QYE8
J	495	HIS	-	expression tag	UNP A0QYE8
J	496	HIS	-	expression tag	UNP A0QYE8
K	2	VAL	-	insertion	UNP A0QYE8
K	480	THR	-	expression tag	UNP A0QYE8
K	481	ALA	-	expression tag	UNP A0QYE8
K	482	ALA	-	expression tag	UNP A0QYE8
K	483	ALA	-	expression tag	UNP A0QYE8
K	484	LYS	-	expression tag	UNP A0QYE8
K	485	GLU	-	expression tag	UNP A0QYE8
K	486	ASP	-	expression tag	UNP A0QYE8
K	487	LEU	-	expression tag	UNP A0QYE8
K	488	GLU	-	expression tag	UNP A0QYE8
K	489	HIS	-	expression tag	UNP A0QYE8
K	490	HIS	-	expression tag	UNP A0QYE8
K	491	HIS	-	expression tag	UNP A0QYE8
K	492	HIS	-	expression tag	UNP A0QYE8
K	493	HIS	-	expression tag	UNP A0QYE8
K	494	HIS	-	expression tag	UNP A0QYE8
K	495	HIS	-	expression tag	UNP A0QYE8
K	496	HIS	-	expression tag	UNP A0QYE8
L	2	VAL	-	insertion	UNP A0QYE8
L	480	THR	-	expression tag	UNP A0QYE8
L	481	ALA	-	expression tag	UNP A0QYE8
L	482	ALA	-	expression tag	UNP A0QYE8
L	483	ALA	-	expression tag	UNP A0QYE8
L	484	LYS	-	expression tag	UNP A0QYE8
L	485	GLU	-	expression tag	UNP A0QYE8
L	486	ASP	-	expression tag	UNP A0QYE8
L	487	LEU	-	expression tag	UNP A0QYE8
L	488	GLU	-	expression tag	UNP A0QYE8
L	489	HIS	-	expression tag	UNP A0QYE8
L	490	HIS	-	expression tag	UNP A0QYE8

Continued on next page...

Continued from previous page...

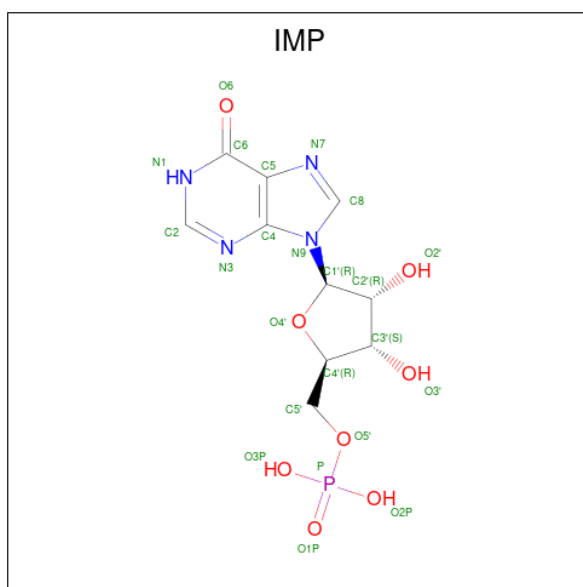
Chain	Residue	Modelled	Actual	Comment	Reference
L	491	HIS	-	expression tag	UNP A0QYE8
L	492	HIS	-	expression tag	UNP A0QYE8
L	493	HIS	-	expression tag	UNP A0QYE8
L	494	HIS	-	expression tag	UNP A0QYE8
L	495	HIS	-	expression tag	UNP A0QYE8
L	496	HIS	-	expression tag	UNP A0QYE8
M	2	VAL	-	insertion	UNP A0QYE8
M	480	THR	-	expression tag	UNP A0QYE8
M	481	ALA	-	expression tag	UNP A0QYE8
M	482	ALA	-	expression tag	UNP A0QYE8
M	483	ALA	-	expression tag	UNP A0QYE8
M	484	LYS	-	expression tag	UNP A0QYE8
M	485	GLU	-	expression tag	UNP A0QYE8
M	486	ASP	-	expression tag	UNP A0QYE8
M	487	LEU	-	expression tag	UNP A0QYE8
M	488	GLU	-	expression tag	UNP A0QYE8
M	489	HIS	-	expression tag	UNP A0QYE8
M	490	HIS	-	expression tag	UNP A0QYE8
M	491	HIS	-	expression tag	UNP A0QYE8
M	492	HIS	-	expression tag	UNP A0QYE8
M	493	HIS	-	expression tag	UNP A0QYE8
M	494	HIS	-	expression tag	UNP A0QYE8
M	495	HIS	-	expression tag	UNP A0QYE8
M	496	HIS	-	expression tag	UNP A0QYE8
N	2	VAL	-	insertion	UNP A0QYE8
N	480	THR	-	expression tag	UNP A0QYE8
N	481	ALA	-	expression tag	UNP A0QYE8
N	482	ALA	-	expression tag	UNP A0QYE8
N	483	ALA	-	expression tag	UNP A0QYE8
N	484	LYS	-	expression tag	UNP A0QYE8
N	485	GLU	-	expression tag	UNP A0QYE8
N	486	ASP	-	expression tag	UNP A0QYE8
N	487	LEU	-	expression tag	UNP A0QYE8
N	488	GLU	-	expression tag	UNP A0QYE8
N	489	HIS	-	expression tag	UNP A0QYE8
N	490	HIS	-	expression tag	UNP A0QYE8
N	491	HIS	-	expression tag	UNP A0QYE8
N	492	HIS	-	expression tag	UNP A0QYE8
N	493	HIS	-	expression tag	UNP A0QYE8
N	494	HIS	-	expression tag	UNP A0QYE8
N	495	HIS	-	expression tag	UNP A0QYE8
N	496	HIS	-	expression tag	UNP A0QYE8

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
O	2	VAL	-	insertion	UNP A0QYE8
O	480	THR	-	expression tag	UNP A0QYE8
O	481	ALA	-	expression tag	UNP A0QYE8
O	482	ALA	-	expression tag	UNP A0QYE8
O	483	ALA	-	expression tag	UNP A0QYE8
O	484	LYS	-	expression tag	UNP A0QYE8
O	485	GLU	-	expression tag	UNP A0QYE8
O	486	ASP	-	expression tag	UNP A0QYE8
O	487	LEU	-	expression tag	UNP A0QYE8
O	488	GLU	-	expression tag	UNP A0QYE8
O	489	HIS	-	expression tag	UNP A0QYE8
O	490	HIS	-	expression tag	UNP A0QYE8
O	491	HIS	-	expression tag	UNP A0QYE8
O	492	HIS	-	expression tag	UNP A0QYE8
O	493	HIS	-	expression tag	UNP A0QYE8
O	494	HIS	-	expression tag	UNP A0QYE8
O	495	HIS	-	expression tag	UNP A0QYE8
O	496	HIS	-	expression tag	UNP A0QYE8
P	2	VAL	-	insertion	UNP A0QYE8
P	480	THR	-	expression tag	UNP A0QYE8
P	481	ALA	-	expression tag	UNP A0QYE8
P	482	ALA	-	expression tag	UNP A0QYE8
P	483	ALA	-	expression tag	UNP A0QYE8
P	484	LYS	-	expression tag	UNP A0QYE8
P	485	GLU	-	expression tag	UNP A0QYE8
P	486	ASP	-	expression tag	UNP A0QYE8
P	487	LEU	-	expression tag	UNP A0QYE8
P	488	GLU	-	expression tag	UNP A0QYE8
P	489	HIS	-	expression tag	UNP A0QYE8
P	490	HIS	-	expression tag	UNP A0QYE8
P	491	HIS	-	expression tag	UNP A0QYE8
P	492	HIS	-	expression tag	UNP A0QYE8
P	493	HIS	-	expression tag	UNP A0QYE8
P	494	HIS	-	expression tag	UNP A0QYE8
P	495	HIS	-	expression tag	UNP A0QYE8
P	496	HIS	-	expression tag	UNP A0QYE8

- Molecule 2 is INOSINIC ACID (CCD ID: IMP) (formula: $C_{10}H_{13}N_4O_8P$) (labeled as "Ligand of Interest" by depositor).



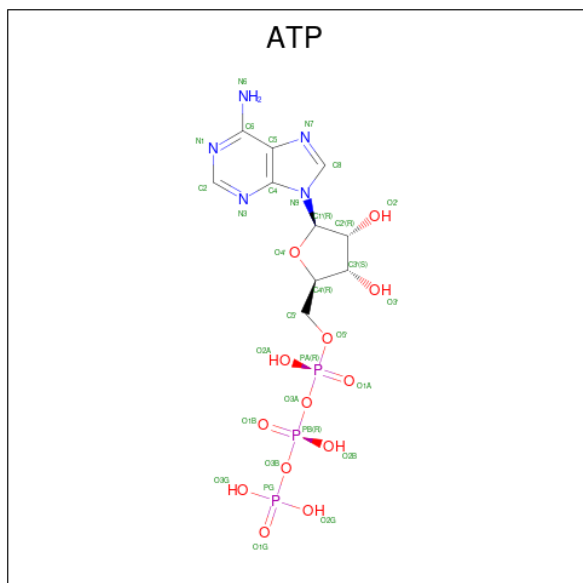
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	B	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	C	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	D	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	E	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	F	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	G	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	H	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	I	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	J	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	K	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	L	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	M	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	N	1	Total	C	N	O	P	0	0
			23	10	4	8	1		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	O	1	Total	C	N	O	P	0	0
			23	10	4	8	1		
2	P	1	Total	C	N	O	P	0	0
			23	10	4	8	1		

- Molecule 3 is ADENOSINE-5'-TRIPHOSPHATE (CCD ID: ATP) (formula: $C_{10}H_{16}N_5O_{13}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	B	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	C	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	D	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	E	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	F	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	G	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	H	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	I	1	Total	C	N	O	P	0	1
			62	20	10	26	6		

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	J	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	K	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	L	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	M	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	N	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	O	1	Total	C	N	O	P	0	1
			62	20	10	26	6		
3	P	1	Total	C	N	O	P	0	1
			62	20	10	26	6		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	213	Total	O	0	0
			213	213		
4	B	189	Total	O	0	0
			189	189		
4	C	211	Total	O	0	0
			211	211		
4	D	251	Total	O	0	0
			251	251		
4	E	241	Total	O	0	0
			241	241		
4	F	235	Total	O	0	0
			235	235		
4	G	236	Total	O	0	0
			236	236		
4	H	244	Total	O	0	0
			244	244		
4	I	116	Total	O	0	0
			116	116		
4	J	212	Total	O	0	0
			212	212		
4	K	200	Total	O	0	0
			200	200		
4	L	134	Total	O	0	0
			134	134		

Continued on next page...

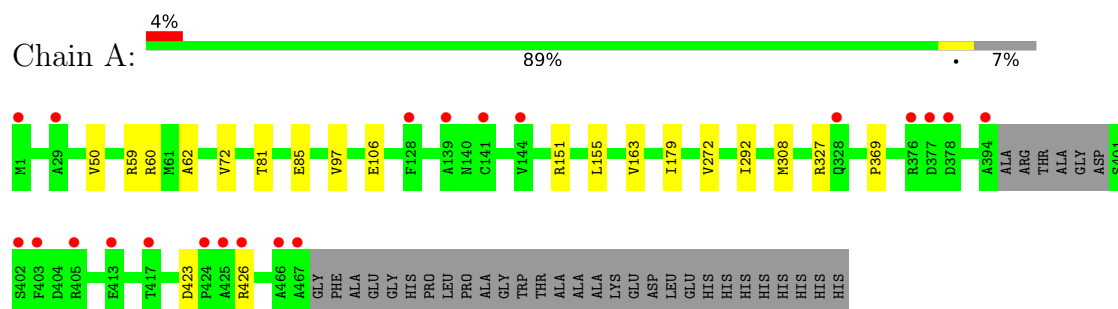
Continued from previous page...

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	M	119	Total 119	O 119	0	0
4	N	149	Total 149	O 149	0	0
4	O	120	Total 120	O 120	0	0
4	P	103	Total 103	O 103	0	0

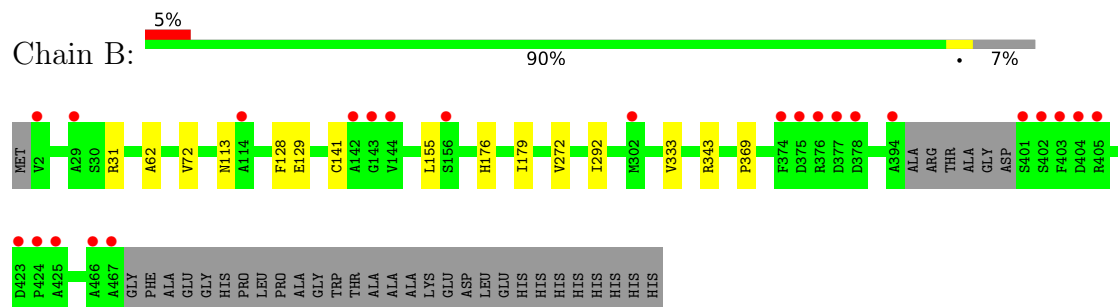
3 Residue-property plots [i](#)

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.

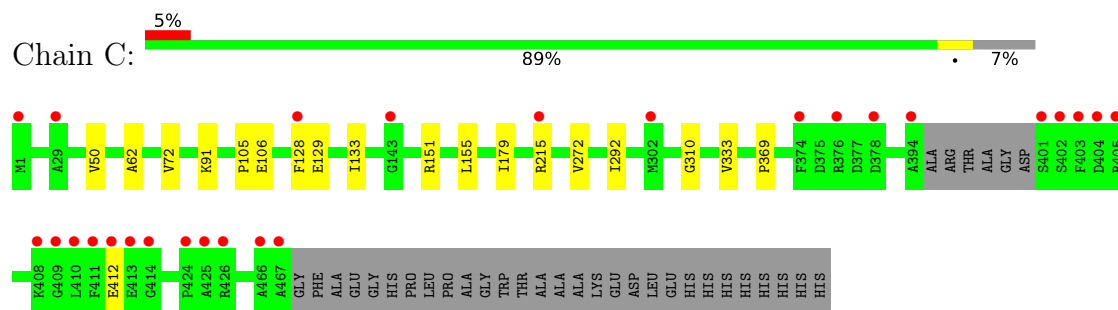
• Molecule 1: GMP reductase



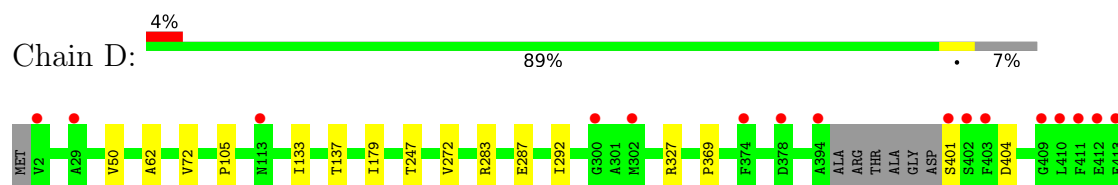
• Molecule 1: GMP reductase

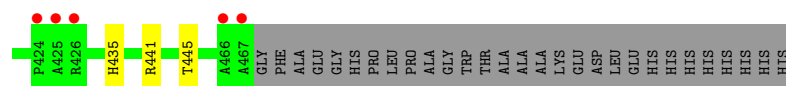


• Molecule 1: GMP reductase

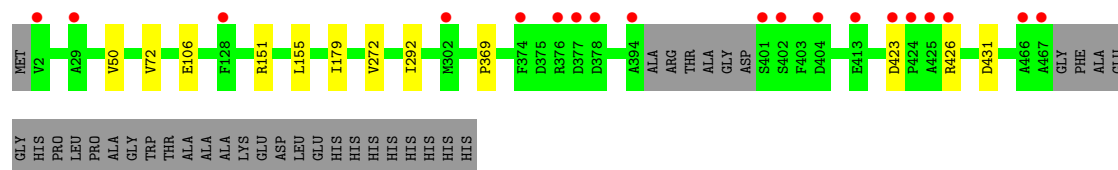
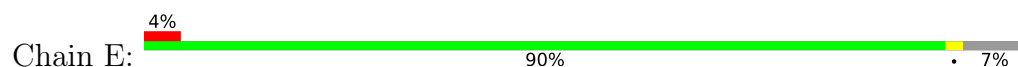


• Molecule 1: GMP reductase

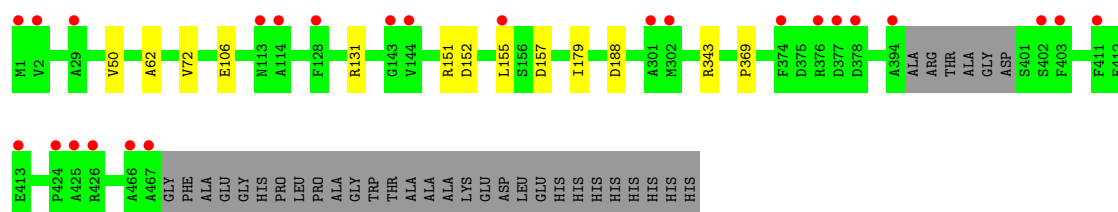
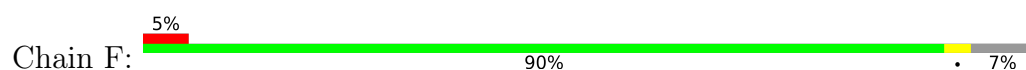




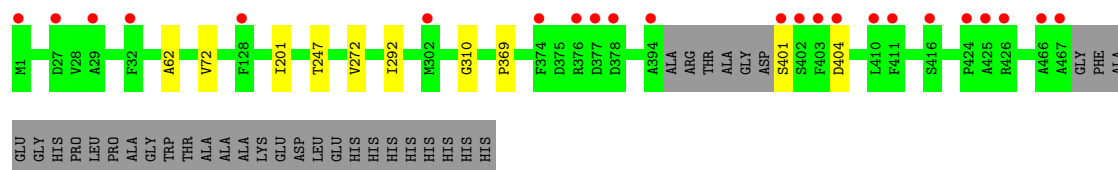
• Molecule 1: GMP reductase



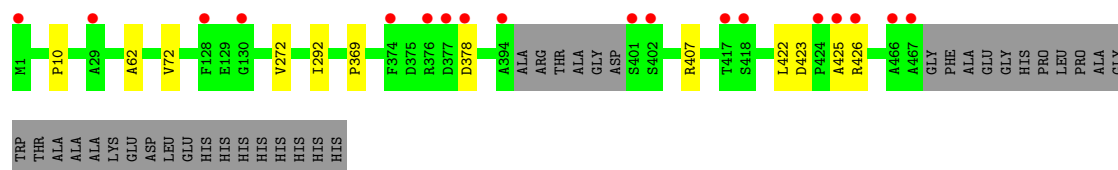
• Molecule 1: GMP reductase



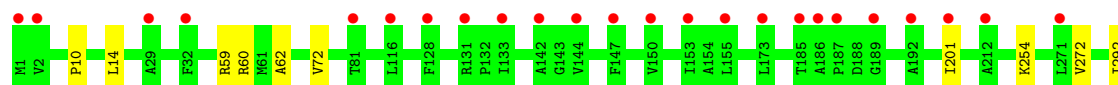
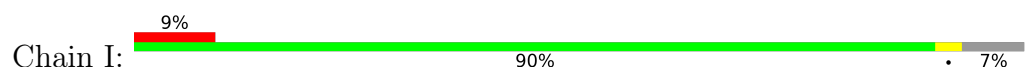
• Molecule 1: GMP reductase



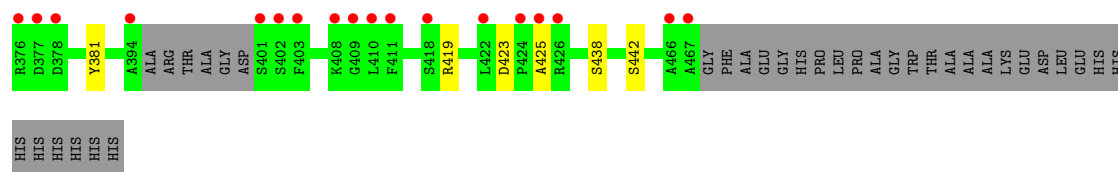
• Molecule 1: GMP reductase



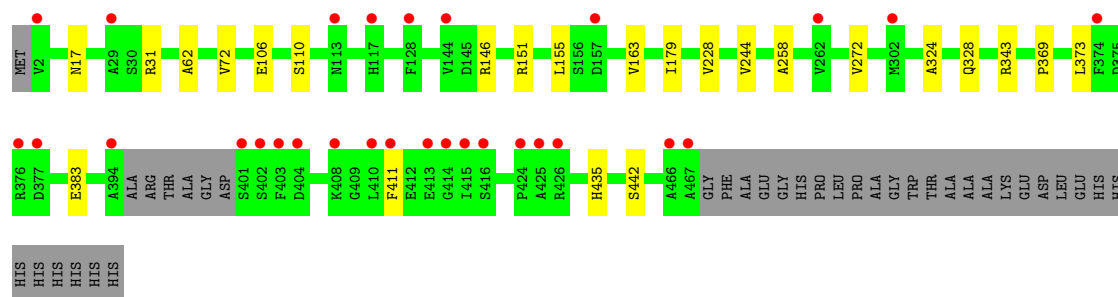
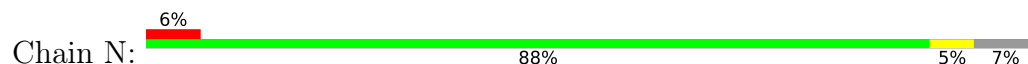
• Molecule 1: GMP reductase



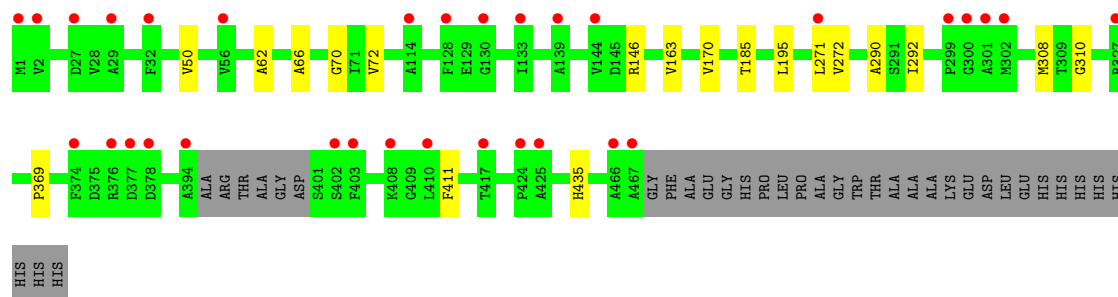
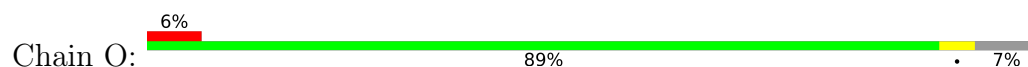




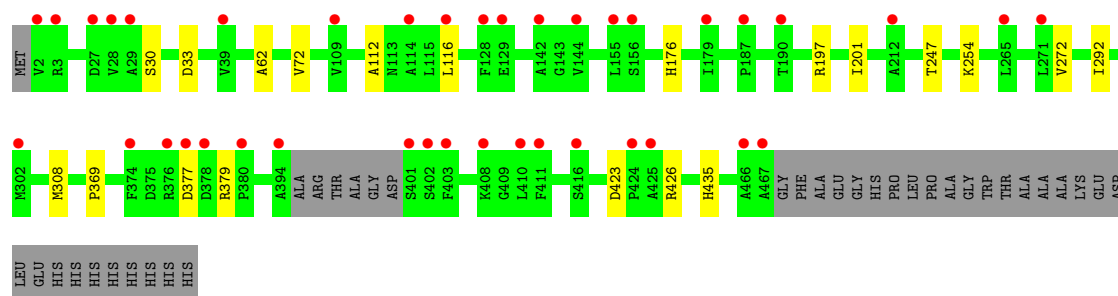
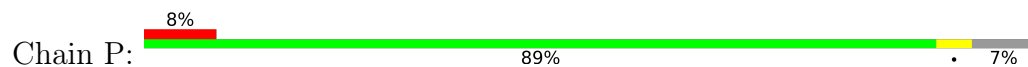
• Molecule 1: GMP reductase



• Molecule 1: GMP reductase



• Molecule 1: GMP reductase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	113.94Å 115.74Å 178.92Å 77.75° 83.32° 65.60°	Depositor
Resolution (Å)	45.50 – 1.60 45.50 – 1.60	Depositor EDS
% Data completeness (in resolution range)	98.6 (45.50-1.60) 99.1 (45.50-1.60)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.04 (at 1.60Å)	Xtriage
Refinement program	PHENIX 1.21_5204	Depositor
R, R_{free}	0.202 , 0.225 0.202 , 0.225	Depositor DCC
R_{free} test set	53152 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	25.5	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 28.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	57872	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.96% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ATP, IMP

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	0.19	0/3420	0.38	0/4655
1	B	0.18	0/3398	0.38	0/4629
1	C	0.20	0/3421	0.40	0/4658
1	D	0.21	0/3407	0.41	0/4640
1	E	0.21	0/3413	0.41	0/4647
1	F	0.22	0/3415	0.42	0/4650
1	G	0.21	0/3400	0.40	0/4632
1	H	0.21	0/3407	0.41	0/4640
1	I	0.15	0/3404	0.34	0/4637
1	J	0.19	0/3407	0.39	0/4639
1	K	0.19	0/3401	0.38	0/4632
1	L	0.17	0/3400	0.35	0/4632
1	M	0.14	0/3398	0.32	0/4630
1	N	0.16	0/3395	0.34	0/4625
1	O	0.16	0/3410	0.33	0/4644
1	P	0.14	0/3393	0.33	0/4623
All	All	0.19	0/54489	0.38	0/74213

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

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	3359	0	3369	13	0
1	B	3339	0	3330	10	0
1	C	3360	0	3357	12	0
1	D	3348	0	3343	13	0
1	E	3352	0	3355	9	0
1	F	3356	0	3355	10	0
1	G	3341	0	3331	7	0
1	H	3348	0	3344	7	0
1	I	3345	0	3338	12	0
1	J	3348	0	3351	14	0
1	K	3342	0	3336	13	0
1	L	3341	0	3331	15	0
1	M	3339	0	3326	13	0
1	N	3336	0	3328	17	0
1	O	3351	0	3346	13	0
1	P	3334	0	3321	12	0
2	A	23	0	11	0	0
2	B	23	0	11	0	0
2	C	23	0	11	0	0
2	D	23	0	11	0	0
2	E	23	0	11	0	0
2	F	23	0	11	0	0
2	G	23	0	11	0	0
2	H	23	0	11	0	0
2	I	23	0	11	0	0
2	J	23	0	11	0	0
2	K	23	0	11	0	0
2	L	23	0	11	0	0
2	M	23	0	11	0	0
2	N	23	0	11	0	0
2	O	23	0	11	0	0
2	P	23	0	11	0	0
3	A	62	0	24	2	0
3	B	62	0	24	2	0
3	C	62	0	24	2	0
3	D	62	0	24	2	0
3	E	62	0	24	2	0
3	F	62	0	24	2	0
3	G	62	0	24	2	0
3	H	62	0	24	1	0
3	I	62	0	24	2	0
3	J	62	0	24	2	0
3	K	62	0	24	4	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	L	62	0	24	2	0
3	M	62	0	24	2	0
3	N	62	0	24	1	0
3	O	62	0	24	1	0
3	P	62	0	24	1	0
4	A	213	0	0	0	0
4	B	189	0	0	0	0
4	C	211	0	0	0	0
4	D	251	0	0	2	0
4	E	241	0	0	0	0
4	F	235	0	0	1	0
4	G	236	0	0	0	0
4	H	244	0	0	1	0
4	I	116	0	0	2	0
4	J	212	0	0	2	0
4	K	200	0	0	1	0
4	L	134	0	0	0	0
4	M	119	0	0	0	0
4	N	149	0	0	0	0
4	O	120	0	0	0	0
4	P	103	0	0	0	0
All	All	57872	0	54021	170	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 2.

All (170) 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:N:31:ARG:NH2	1:O:411:PHE:O	2.29	0.65
1:C:369:PRO:HB3	3:C:502[B]:ATP:H5'2	1.81	0.62
1:M:272:VAL:HG22	1:M:292:ILE:HB	1.83	0.61
1:E:423:ASP:CG	1:E:426:ARG:H	2.07	0.61
1:A:369:PRO:HB3	3:A:502[B]:ATP:H5'2	1.84	0.59
1:B:369:PRO:HB3	3:B:502[B]:ATP:H5'2	1.84	0.59
1:E:369:PRO:HB3	3:E:502[A]:ATP:H5'2	1.85	0.58
1:F:369:PRO:HB3	3:F:502[A]:ATP:H5'2	1.85	0.57
1:C:155:LEU:HD12	1:C:179:ILE:HD11	1.87	0.57
1:J:369:PRO:HB3	3:J:502[A]:ATP:H5'2	1.87	0.57
1:M:423:ASP:OD2	1:M:425:ALA:N	2.29	0.57
1:C:272:VAL:HG22	1:C:292:ILE:HB	1.86	0.57

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:401:SER:HB3	1:D:404:ASP:OD2	2.06	0.56
1:G:369:PRO:HB3	3:G:502[A]:ATP:H5'2	1.87	0.56
1:K:59:ARG:NH1	3:K:502[B]:ATP:O2A	2.38	0.56
1:L:157:ASP:OD2	1:P:176:HIS:NE2	2.34	0.56
1:P:377:ASP:HB2	1:P:379:ARG:HD3	1.88	0.55
1:I:272:VAL:HG22	1:I:292:ILE:HB	1.89	0.55
1:G:401:SER:HB3	1:G:404:ASP:OD2	2.08	0.54
1:N:62:ALA:HA	1:N:72:VAL:HG21	1.90	0.54
1:E:155:LEU:HD12	1:E:179:ILE:HD11	1.90	0.54
1:J:369:PRO:HB3	3:J:502[B]:ATP:H5'2	1.89	0.53
1:C:128:PHE:O	1:C:129:GLU:HG2	2.09	0.53
1:K:369:PRO:HB3	3:K:502[A]:ATP:H5'2	1.91	0.53
1:L:106:GLU:OE2	1:L:151:ARG:NH1	2.39	0.53
1:H:369:PRO:HB3	3:H:502[A]:ATP:H5'2	1.91	0.52
1:A:106:GLU:OE2	1:A:151:ARG:NH2	2.42	0.52
1:D:369:PRO:HB3	3:D:502[A]:ATP:H5'2	1.91	0.52
1:F:369:PRO:HB3	3:F:502[B]:ATP:H5'2	1.90	0.52
1:J:272:VAL:HG22	1:J:292:ILE:HB	1.90	0.52
1:L:369:PRO:HB3	3:L:502[B]:ATP:H5'2	1.91	0.52
1:N:369:PRO:HB3	3:N:502[B]:ATP:H5'2	1.91	0.52
1:P:369:PRO:HB3	3:P:502[A]:ATP:H5'2	1.92	0.52
1:J:106:GLU:OE2	1:J:151:ARG:NH2	2.40	0.51
1:D:105:PRO:HG3	1:D:133:ILE:HD11	1.92	0.51
1:L:369:PRO:HB3	3:L:502[A]:ATP:H5'2	1.91	0.51
1:K:59:ARG:NH1	3:K:502[A]:ATP:O1A	2.44	0.51
1:M:31:ARG:HD3	1:M:442[A]:SER:OG	2.11	0.51
1:D:272:VAL:HG22	1:D:292:ILE:HB	1.93	0.51
1:E:423:ASP:OD2	1:E:426:ARG:HB2	2.12	0.51
1:K:106:GLU:OE2	1:K:151:ARG:NH2	2.38	0.51
1:L:31:ARG:HD3	1:L:442[A]:SER:OG	2.11	0.50
1:J:383:GLU:OE1	1:O:146:ARG:NH2	2.40	0.50
1:N:106:GLU:OE2	1:N:151:ARG:NH2	2.41	0.50
1:F:155:LEU:HD12	1:F:179:ILE:HD11	1.94	0.49
1:L:62:ALA:HA	1:L:72:VAL:HG21	1.95	0.49
1:B:272:VAL:HG22	1:B:292:ILE:HB	1.93	0.49
1:P:112:ALA:O	1:P:116:LEU:HD13	2.12	0.49
1:A:272:VAL:HG22	1:A:292:ILE:HB	1.94	0.49
1:I:62:ALA:HA	1:I:72:VAL:HG21	1.93	0.49
1:I:369:PRO:HB3	3:I:502[B]:ATP:H5'2	1.94	0.49
1:B:31:ARG:HH22	1:C:412:GLU:HA	1.77	0.49
1:P:197:ARG:O	1:P:201:ILE:HG12	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:106:GLU:OE2	1:F:151:ARG:NH2	2.30	0.48
1:H:423:ASP:OD1	1:H:425:ALA:N	2.36	0.48
1:M:369:PRO:HB3	3:M:502[B]:ATP:H5'2	1.94	0.48
1:G:62:ALA:HA	1:G:72:VAL:HG21	1.96	0.48
1:O:435:HIS:NE2	1:P:308:MET:SD	2.87	0.48
1:N:110:SER:HB2	1:N:146:ARG:HG2	1.96	0.48
1:C:91:LYS:HB3	1:C:215:ARG:HE	1.79	0.48
1:C:369:PRO:HB3	3:C:502[A]:ATP:H5'2	1.96	0.48
1:E:272:VAL:HG22	1:E:292:ILE:HB	1.96	0.48
1:K:62:ALA:HA	1:K:72:VAL:HG21	1.96	0.48
1:D:62:ALA:HA	1:D:72:VAL:HG21	1.96	0.47
1:F:131:ARG:NH1	1:F:188:ASP:O	2.47	0.47
1:O:369:PRO:HB3	3:O:502[A]:ATP:H5'2	1.95	0.47
1:A:308:MET:SD	1:D:435:HIS:NE2	2.88	0.47
1:H:62:ALA:HA	1:H:72:VAL:HG21	1.97	0.47
1:G:201:ILE:HG23	1:H:407:ARG:HA	1.96	0.46
1:N:324:ALA:O	1:N:328:GLN:HG3	2.15	0.46
1:H:10:PRO:HB3	4:H:686:HOH:O	2.14	0.46
1:N:31:ARG:HD3	1:N:442[A]:SER:OG	2.15	0.46
1:O:163:VAL:HG12	1:O:185:THR:O	2.16	0.46
1:I:369:PRO:HB3	3:I:502[A]:ATP:H5'2	1.98	0.46
1:A:155:LEU:HD12	1:A:179:ILE:HD11	1.97	0.46
1:D:369:PRO:HB3	3:D:502[B]:ATP:H5'2	1.97	0.46
1:I:308:MET:SD	1:L:435:HIS:NE2	2.89	0.46
1:N:228:VAL:HG11	1:N:258:ALA:HB1	1.98	0.46
1:G:272:VAL:HG22	1:G:292:ILE:HB	1.97	0.46
1:E:426:ARG:HG2	1:E:431:ASP:HB3	1.97	0.45
1:M:261:ALA:O	1:M:265:LEU:HD13	2.15	0.45
1:I:310:GLY:HA2	1:L:343:ARG:HG3	1.98	0.45
1:G:369:PRO:HB3	3:G:502[B]:ATP:H5'2	1.99	0.45
1:M:50:VAL:HB	1:M:72:VAL:HG22	1.99	0.45
1:M:62:ALA:HA	1:M:72:VAL:HG21	1.99	0.45
1:C:62:ALA:HA	1:C:72:VAL:HG21	1.97	0.45
1:H:272:VAL:HG22	1:H:292:ILE:HB	1.99	0.45
4:I:656:HOH:O	1:L:10:PRO:HB3	2.16	0.45
1:K:146:ARG:NH2	1:N:383:GLU:OE1	2.38	0.45
1:I:59:ARG:HG3	1:I:60:ARG:N	2.32	0.45
1:P:30:SER:HB3	1:P:33:ASP:OD2	2.17	0.45
1:E:369:PRO:HB3	3:E:502[B]:ATP:H5'2	1.98	0.45
1:M:283:ARG:HD3	1:M:287:GLU:OE2	2.17	0.45
1:N:106:GLU:OE2	1:N:151:ARG:NH1	2.49	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:267:LEU:HB2	4:J:780:HOH:O	2.17	0.44
1:O:50:VAL:HB	1:O:72:VAL:HG22	1.99	0.44
1:N:17:ASN:OD1	1:N:343:ARG:NH2	2.49	0.44
1:B:62:ALA:HA	1:B:72:VAL:HG21	1.99	0.44
1:B:176:HIS:NE2	1:F:157:ASP:OD2	2.48	0.44
1:A:81:THR:O	1:A:85:GLU:HG3	2.18	0.44
1:I:343:ARG:HG3	1:J:310:GLY:HA2	1.99	0.44
1:N:343:ARG:HG3	1:O:310:GLY:HA2	2.00	0.44
1:P:254:LYS:HE2	1:P:254:LYS:HB3	1.84	0.44
1:A:97:VAL:HG21	1:A:163:VAL:HG23	2.00	0.43
1:D:283:ARG:HD3	4:D:740:HOH:O	2.18	0.43
1:I:411:PHE:O	1:L:31:ARG:NH2	2.51	0.43
1:A:50:VAL:HB	1:A:72:VAL:HG22	2.00	0.43
1:I:201:ILE:HD13	1:I:201:ILE:HA	1.91	0.43
1:L:170:VAL:HG11	1:L:195:LEU:HD23	2.01	0.43
1:O:272:VAL:HG22	1:O:292:ILE:HB	2.01	0.43
1:A:369:PRO:HB3	3:A:502[A]:ATP:H5'2	2.00	0.43
1:D:137:THR:HG21	1:D:179:ILE:HD12	2.00	0.43
1:M:369:PRO:HB3	3:M:502[A]:ATP:H5'2	2.01	0.43
1:K:262:VAL:HG11	1:K:271:LEU:HD21	2.00	0.43
1:L:333:VAL:HG23	1:L:353:ALA:HA	1.99	0.43
1:B:155:LEU:HD12	1:B:179:ILE:HD11	2.01	0.43
1:D:50:VAL:HB	1:D:72:VAL:HG22	2.00	0.43
1:D:283:ARG:O	1:D:287:GLU:HG3	2.19	0.43
1:J:31:ARG:NE	1:J:442[A]:SER:OG	2.51	0.43
1:J:168:ARG:N	1:J:168:ARG:HD3	2.34	0.43
1:C:105:PRO:HG3	1:C:133:ILE:HD11	2.01	0.43
1:F:62:ALA:HA	1:F:72:VAL:HG21	2.00	0.43
1:N:435:HIS:NE2	1:O:308:MET:SD	2.92	0.43
1:K:112:ALA:O	1:K:116:LEU:HD23	2.19	0.42
1:P:62:ALA:HA	1:P:72:VAL:HG21	2.01	0.42
1:P:272:VAL:HG22	1:P:292:ILE:HB	2.01	0.42
1:A:62:ALA:HA	1:A:72:VAL:HG21	2.00	0.42
1:C:50:VAL:HB	1:C:72:VAL:HG22	2.00	0.42
1:M:302:MET:HE3	1:M:302:MET:HB2	1.86	0.42
1:O:62:ALA:HA	1:O:72:VAL:HG21	2.01	0.42
1:B:128:PHE:O	1:B:129:GLU:HG2	2.19	0.42
1:K:78:LEU:HD23	1:K:83:VAL:HG22	2.01	0.42
1:A:106:GLU:OE2	1:A:151:ARG:NH1	2.49	0.42
1:M:381:TYR:CE2	1:M:419:ARG:HG2	2.55	0.42
1:N:155:LEU:HD12	1:N:179:ILE:HD11	2.02	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:244:VAL:HG22	1:N:272:VAL:HB	2.02	0.42
1:O:66:ALA:HA	1:O:70:GLY:O	2.20	0.42
1:B:343:ARG:HG3	1:C:310:GLY:HA2	2.02	0.41
1:L:254:LYS:HE3	1:L:254:LYS:HB3	1.82	0.41
1:E:106:GLU:OE2	1:E:151:ARG:NH1	2.49	0.41
1:K:327:ARG:HD2	4:K:717:HOH:O	2.19	0.41
1:L:31:ARG:HH11	1:L:442[A]:SER:HG	1.65	0.41
1:B:369:PRO:HB3	3:B:502[A]:ATP:H5'2	2.00	0.41
1:J:435:HIS:NE2	1:K:308:MET:SD	2.94	0.41
1:O:170:VAL:HG11	1:O:195:LEU:HD23	2.02	0.41
1:D:327:ARG:HD2	4:D:757:HOH:O	2.20	0.41
1:J:10:PRO:HB3	4:J:727:HOH:O	2.21	0.41
1:K:272:VAL:HG22	1:K:292:ILE:HB	2.01	0.41
1:F:50:VAL:HB	1:F:72:VAL:HG22	2.02	0.41
1:F:152:ASP:OD1	4:F:601:HOH:O	2.22	0.41
1:F:343:ARG:HG3	1:G:310:GLY:HA2	2.03	0.41
1:L:20:PHE:CZ	1:L:462:GLY:HA3	2.56	0.41
1:A:59:ARG:HG3	1:A:60:ARG:N	2.36	0.41
1:D:441:ARG:O	1:D:445:THR:HG23	2.21	0.41
1:J:283:ARG:O	1:J:287:GLU:HG3	2.21	0.41
1:C:106:GLU:OE2	1:C:151:ARG:NH2	2.50	0.41
1:N:62:ALA:HA	1:N:72:VAL:CG2	2.50	0.41
1:B:113:ASN:HD22	1:B:141:CYS:CB	2.34	0.40
1:P:423:ASP:HB3	1:P:426:ARG:HB2	2.04	0.40
1:E:50:VAL:HB	1:E:72:VAL:HG22	2.04	0.40
1:J:166:ASP:OD1	1:J:168:ARG:NE	2.53	0.40
1:M:308:MET:SD	1:P:435:HIS:NE2	2.94	0.40
1:O:271:LEU:HD22	1:O:290:ALA:HA	2.02	0.40
1:A:423:ASP:HB3	1:A:426:ARG:HB2	2.03	0.40
1:I:10:PRO:HG3	1:I:14:LEU:HD21	2.03	0.40
1:K:369:PRO:HB3	3:K:502[B]:ATP:H5'2	2.03	0.40
1:H:423:ASP:HB3	1:H:426:ARG:HB2	2.03	0.40
1:I:254:LYS:HG3	4:I:659:HOH:O	2.21	0.40
1:J:128:PHE:CE1	1:J:129:GLU:HG3	2.57	0.40
1:M:438:SER:HB2	1:N:411:PHE:CD1	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	461/496 (93%)	452 (98%)	9 (2%)	0	100	100
1	B	459/496 (92%)	452 (98%)	7 (2%)	0	100	100
1	C	461/496 (93%)	452 (98%)	9 (2%)	0	100	100
1	D	459/496 (92%)	451 (98%)	8 (2%)	0	100	100
1	E	460/496 (93%)	450 (98%)	10 (2%)	0	100	100
1	F	460/496 (93%)	451 (98%)	9 (2%)	0	100	100
1	G	460/496 (93%)	451 (98%)	9 (2%)	0	100	100
1	H	460/496 (93%)	451 (98%)	9 (2%)	0	100	100
1	I	460/496 (93%)	451 (98%)	9 (2%)	0	100	100
1	J	460/496 (93%)	450 (98%)	10 (2%)	0	100	100
1	K	459/496 (92%)	450 (98%)	9 (2%)	0	100	100
1	L	460/496 (93%)	448 (97%)	12 (3%)	0	100	100
1	M	460/496 (93%)	450 (98%)	10 (2%)	0	100	100
1	N	459/496 (92%)	450 (98%)	9 (2%)	0	100	100
1	O	460/496 (93%)	450 (98%)	10 (2%)	0	100	100
1	P	459/496 (92%)	452 (98%)	7 (2%)	0	100	100
All	All	7357/7936 (93%)	7211 (98%)	146 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	347/372 (93%)	346 (100%)	1 (0%)	86	78
1	B	344/372 (92%)	343 (100%)	1 (0%)	86	78
1	C	348/372 (94%)	347 (100%)	1 (0%)	86	78
1	D	347/372 (93%)	346 (100%)	1 (0%)	86	78
1	E	347/372 (93%)	347 (100%)	0	100	100
1	F	348/372 (94%)	348 (100%)	0	100	100
1	G	344/372 (92%)	343 (100%)	1 (0%)	86	78
1	H	345/372 (93%)	343 (99%)	2 (1%)	78	66
1	I	345/372 (93%)	344 (100%)	1 (0%)	86	78
1	J	345/372 (93%)	345 (100%)	0	100	100
1	K	344/372 (92%)	343 (100%)	1 (0%)	86	78
1	L	344/372 (92%)	344 (100%)	0	100	100
1	M	343/372 (92%)	341 (99%)	2 (1%)	78	66
1	N	343/372 (92%)	341 (99%)	2 (1%)	78	66
1	O	346/372 (93%)	346 (100%)	0	100	100
1	P	343/372 (92%)	342 (100%)	1 (0%)	86	78
All	All	5523/5952 (93%)	5509 (100%)	14 (0%)	86	78

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	327	ARG
1	B	333	VAL
1	C	333	VAL
1	D	247	THR
1	G	247	THR
1	H	378	ASP
1	H	422	LEU
1	I	355	ASN
1	K	116	LEU
1	M	163	VAL
1	M	333	VAL
1	N	163	VAL
1	N	373	LEU
1	P	247	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	233	GLN
1	A	252	GLN
1	B	113	ASN
1	B	328	GLN
1	C	252	GLN
1	D	252	GLN
1	D	328	GLN
1	F	117	HIS
1	F	176	HIS
1	H	176	HIS
1	I	140	ASN
1	I	252	GLN
1	J	176	HIS
1	K	140	ASN
1	L	117	HIS
1	N	113	ASN
1	N	328	GLN
1	O	117	HIS
1	P	113	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

48 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	IMP	P	501	-	25,25,25	0.35	0	37,38,38	0.50	0
3	ATP	C	502[A]	-	32,33,33	0.56	0	48,52,52	0.57	0
3	ATP	C	502[B]	-	32,33,33	0.55	0	48,52,52	0.58	0
3	ATP	M	502[A]	-	32,33,33	0.55	0	48,52,52	0.53	0
3	ATP	M	502[B]	-	32,33,33	0.56	0	48,52,52	0.55	0
2	IMP	H	501	-	25,25,25	0.37	0	37,38,38	0.51	0
2	IMP	E	501	-	25,25,25	0.36	0	37,38,38	0.50	0
2	IMP	D	501	-	25,25,25	0.37	0	37,38,38	0.60	0
2	IMP	M	501	-	25,25,25	0.34	0	37,38,38	0.49	0
2	IMP	N	501	-	25,25,25	0.33	0	37,38,38	0.51	0
3	ATP	N	502[A]	-	32,33,33	0.55	0	48,52,52	0.55	0
3	ATP	A	502[A]	-	32,33,33	0.55	0	48,52,52	0.57	0
3	ATP	A	502[B]	-	32,33,33	0.56	0	48,52,52	0.53	0
3	ATP	F	502[A]	-	32,33,33	0.54	0	48,52,52	0.56	0
3	ATP	F	502[B]	-	32,33,33	0.55	0	48,52,52	0.58	0
3	ATP	N	502[B]	-	32,33,33	0.55	0	48,52,52	0.53	0
2	IMP	C	501	-	25,25,25	0.35	0	37,38,38	0.55	0
3	ATP	I	502[A]	-	32,33,33	0.55	0	48,52,52	0.55	0
2	IMP	L	501	-	25,25,25	0.37	0	37,38,38	0.53	0
3	ATP	I	502[B]	-	32,33,33	0.55	0	48,52,52	0.55	0
3	ATP	L	502[A]	-	32,33,33	0.55	0	48,52,52	0.55	0
3	ATP	D	502[B]	-	32,33,33	0.55	0	48,52,52	0.58	0
3	ATP	L	502[B]	-	32,33,33	0.55	0	48,52,52	0.58	0
2	IMP	A	501	-	25,25,25	0.36	0	37,38,38	0.53	0
3	ATP	J	502[B]	-	32,33,33	0.55	0	48,52,52	0.57	0
3	ATP	D	502[A]	-	32,33,33	0.55	0	48,52,52	0.57	0
2	IMP	G	501	-	25,25,25	0.36	0	37,38,38	0.54	0
3	ATP	G	502[A]	-	32,33,33	0.55	0	48,52,52	0.59	0
2	IMP	F	501	-	25,25,25	0.40	0	37,38,38	0.59	0
3	ATP	G	502[B]	-	32,33,33	0.57	0	48,52,52	0.56	0
3	ATP	H	502[A]	-	32,33,33	0.55	0	48,52,52	0.57	0
3	ATP	E	502[A]	-	32,33,33	0.54	0	48,52,52	0.57	0
3	ATP	E	502[B]	-	32,33,33	0.57	0	48,52,52	0.58	0
2	IMP	J	501	-	25,25,25	0.34	0	37,38,38	0.52	0
3	ATP	H	502[B]	-	32,33,33	0.55	0	48,52,52	0.55	0
2	IMP	I	501	-	25,25,25	0.36	0	37,38,38	0.50	0
3	ATP	K	502[A]	-	32,33,33	0.54	0	48,52,52	0.52	0
3	ATP	K	502[B]	-	32,33,33	0.55	0	48,52,52	0.57	0
3	ATP	O	502[A]	-	32,33,33	0.56	0	48,52,52	0.54	0
3	ATP	O	502[B]	-	32,33,33	0.56	0	48,52,52	0.55	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	ATP	B	502[A]	-	32,33,33	0.56	0	48,52,52	0.55	0
3	ATP	B	502[B]	-	32,33,33	0.55	0	48,52,52	0.55	0
2	IMP	O	501	-	25,25,25	0.34	0	37,38,38	0.54	0
3	ATP	P	502[A]	-	32,33,33	0.55	0	48,52,52	0.53	0
3	ATP	P	502[B]	-	32,33,33	0.56	0	48,52,52	0.54	0
2	IMP	B	501	-	25,25,25	0.36	0	37,38,38	0.49	0
3	ATP	J	502[A]	-	32,33,33	0.56	0	48,52,52	0.57	0
2	IMP	K	501	-	25,25,25	0.38	0	37,38,38	0.50	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	IMP	P	501	-	-	0/10/26/26	0/3/3/3
3	ATP	C	502[A]	-	-	2/22/38/38	0/3/3/3
3	ATP	C	502[B]	-	-	4/22/38/38	0/3/3/3
3	ATP	M	502[A]	-	-	0/22/38/38	0/3/3/3
3	ATP	M	502[B]	-	-	4/22/38/38	0/3/3/3
2	IMP	H	501	-	-	0/10/26/26	0/3/3/3
2	IMP	E	501	-	-	0/10/26/26	0/3/3/3
2	IMP	D	501	-	-	0/10/26/26	0/3/3/3
2	IMP	M	501	-	-	0/10/26/26	0/3/3/3
2	IMP	N	501	-	-	0/10/26/26	0/3/3/3
3	ATP	N	502[A]	-	-	0/22/38/38	0/3/3/3
3	ATP	A	502[A]	-	-	3/22/38/38	0/3/3/3
3	ATP	A	502[B]	-	-	4/22/38/38	0/3/3/3
3	ATP	F	502[A]	-	-	2/22/38/38	0/3/3/3
3	ATP	F	502[B]	-	-	2/22/38/38	0/3/3/3
3	ATP	N	502[B]	-	-	4/22/38/38	0/3/3/3
2	IMP	C	501	-	-	0/10/26/26	0/3/3/3
3	ATP	I	502[A]	-	-	7/22/38/38	0/3/3/3
2	IMP	L	501	-	-	0/10/26/26	0/3/3/3
3	ATP	I	502[B]	-	-	0/22/38/38	0/3/3/3
3	ATP	L	502[A]	-	-	3/22/38/38	0/3/3/3
3	ATP	D	502[B]	-	-	2/22/38/38	0/3/3/3
3	ATP	L	502[B]	-	-	1/22/38/38	0/3/3/3
2	IMP	A	501	-	-	0/10/26/26	0/3/3/3
3	ATP	J	502[B]	-	-	3/22/38/38	0/3/3/3

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	ATP	D	502[A]	-	-	3/22/38/38	0/3/3/3
2	IMP	G	501	-	-	0/10/26/26	0/3/3/3
3	ATP	G	502[A]	-	-	7/22/38/38	0/3/3/3
2	IMP	F	501	-	-	0/10/26/26	0/3/3/3
3	ATP	G	502[B]	-	-	2/22/38/38	0/3/3/3
3	ATP	H	502[A]	-	-	2/22/38/38	0/3/3/3
3	ATP	E	502[A]	-	-	2/22/38/38	0/3/3/3
3	ATP	E	502[B]	-	-	3/22/38/38	0/3/3/3
2	IMP	J	501	-	-	0/10/26/26	0/3/3/3
3	ATP	H	502[B]	-	-	0/22/38/38	0/3/3/3
2	IMP	I	501	-	-	0/10/26/26	0/3/3/3
3	ATP	K	502[A]	-	-	2/22/38/38	0/3/3/3
3	ATP	K	502[B]	-	-	1/22/38/38	0/3/3/3
3	ATP	O	502[A]	-	-	7/22/38/38	0/3/3/3
3	ATP	O	502[B]	-	-	3/22/38/38	0/3/3/3
3	ATP	B	502[A]	-	-	3/22/38/38	0/3/3/3
3	ATP	B	502[B]	-	-	3/22/38/38	0/3/3/3
2	IMP	O	501	-	-	0/10/26/26	0/3/3/3
3	ATP	P	502[A]	-	-	4/22/38/38	0/3/3/3
3	ATP	P	502[B]	-	-	2/22/38/38	0/3/3/3
2	IMP	B	501	-	-	0/10/26/26	0/3/3/3
3	ATP	J	502[A]	-	-	3/22/38/38	0/3/3/3
2	IMP	K	501	-	-	0/10/26/26	0/3/3/3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (88) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502[A]	ATP	PB-O3B-PG-O3G
3	B	502[A]	ATP	PB-O3B-PG-O3G
3	C	502[A]	ATP	PB-O3B-PG-O3G
3	C	502[B]	ATP	PB-O3B-PG-O2G
3	D	502[B]	ATP	PB-O3B-PG-O3G
3	E	502[B]	ATP	PB-O3B-PG-O3G
3	F	502[B]	ATP	PB-O3B-PG-O3G
3	G	502[A]	ATP	PB-O3B-PG-O2G

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	I	502[A]	ATP	PB-O3B-PG-O2G
3	O	502[A]	ATP	PB-O3B-PG-O2G
3	P	502[A]	ATP	PB-O3B-PG-O2G
3	P	502[A]	ATP	PB-O3B-PG-O3G
3	P	502[B]	ATP	PB-O3B-PG-O3G
3	K	502[B]	ATP	PB-O3B-PG-O1G
3	N	502[B]	ATP	PB-O3B-PG-O1G
3	B	502[B]	ATP	PB-O3A-PA-O1A
3	D	502[A]	ATP	PA-O3A-PB-O1B
3	G	502[B]	ATP	PA-O3A-PB-O1B
3	J	502[B]	ATP	PA-O3A-PB-O1B
3	C	502[B]	ATP	PB-O3B-PG-O1G
3	A	502[B]	ATP	PB-O3B-PG-O2G
3	A	502[B]	ATP	PB-O3B-PG-O3G
3	M	502[B]	ATP	PB-O3B-PG-O2G
3	N	502[B]	ATP	PB-O3B-PG-O2G
3	F	502[A]	ATP	PB-O3A-PA-O2A
3	G	502[A]	ATP	PB-O3A-PA-O2A
3	J	502[A]	ATP	PG-O3B-PB-O2B
3	K	502[A]	ATP	PB-O3A-PA-O2A
3	L	502[B]	ATP	PA-O3A-PB-O1B
3	M	502[B]	ATP	PB-O3A-PA-O2A
3	O	502[A]	ATP	PB-O3A-PA-O2A
3	O	502[B]	ATP	PG-O3B-PB-O2B
3	F	502[B]	ATP	PB-O3B-PG-O1G
3	A	502[B]	ATP	PB-O3A-PA-O2A
3	B	502[B]	ATP	PB-O3A-PA-O2A
3	C	502[B]	ATP	PG-O3B-PB-O2B
3	C	502[B]	ATP	PB-O3A-PA-O2A
3	E	502[A]	ATP	PB-O3A-PA-O2A
3	H	502[A]	ATP	PA-O3A-PB-O1B
3	H	502[A]	ATP	PB-O3A-PA-O2A
3	I	502[A]	ATP	PB-O3A-PA-O2A
3	L	502[A]	ATP	PA-O3A-PB-O1B
3	M	502[B]	ATP	PB-O3A-PA-O1A
3	N	502[B]	ATP	PB-O3A-PA-O2A
3	P	502[A]	ATP	PB-O3A-PA-O1A
3	P	502[A]	ATP	PB-O3A-PA-O2A
3	B	502[A]	ATP	PB-O3B-PG-O1G
3	C	502[A]	ATP	PB-O3B-PG-O1G
3	D	502[B]	ATP	PB-O3B-PG-O1G
3	O	502[B]	ATP	PB-O3B-PG-O1G

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
3	A	502[B]	ATP	PA-O3A-PB-O1B
3	G	502[B]	ATP	PA-O3A-PB-O2B
3	I	502[A]	ATP	PA-O3A-PB-O1B
3	A	502[A]	ATP	PB-O3B-PG-O1G
3	I	502[A]	ATP	PB-O3B-PG-O1G
3	P	502[B]	ATP	PB-O3B-PG-O1G
3	A	502[A]	ATP	PB-O3B-PG-O2G
3	B	502[A]	ATP	PB-O3B-PG-O2G
3	B	502[B]	ATP	PB-O3B-PG-O3G
3	E	502[A]	ATP	PB-O3B-PG-O3G
3	E	502[B]	ATP	PB-O3B-PG-O2G
3	G	502[A]	ATP	PB-O3B-PG-O3G
3	I	502[A]	ATP	PB-O3B-PG-O3G
3	M	502[B]	ATP	PB-O3B-PG-O3G
3	O	502[A]	ATP	PB-O3B-PG-O3G
3	D	502[A]	ATP	PB-O3A-PA-O1A
3	G	502[A]	ATP	PG-O3B-PB-O2B
3	G	502[A]	ATP	PB-O3A-PA-O1A
3	I	502[A]	ATP	PG-O3B-PB-O1B
3	I	502[A]	ATP	PB-O3A-PA-O1A
3	J	502[A]	ATP	PG-O3B-PB-O1B
3	J	502[B]	ATP	PB-O3A-PA-O2A
3	K	502[A]	ATP	PB-O3A-PA-O1A
3	L	502[A]	ATP	PB-O3A-PA-O2A
3	N	502[B]	ATP	PB-O3A-PA-O1A
3	O	502[A]	ATP	PG-O3B-PB-O2B
3	O	502[A]	ATP	PB-O3A-PA-O1A
3	E	502[B]	ATP	PB-O3B-PG-O1G
3	G	502[A]	ATP	PB-O3B-PG-O1G
3	O	502[A]	ATP	PB-O3B-PG-O1G
3	D	502[A]	ATP	PB-O3A-PA-O2A
3	F	502[A]	ATP	PB-O3A-PA-O1A
3	G	502[A]	ATP	PG-O3B-PB-O1B
3	J	502[A]	ATP	PA-O3A-PB-O2B
3	J	502[B]	ATP	PB-O3A-PA-O1A
3	L	502[A]	ATP	PB-O3A-PA-O1A
3	O	502[A]	ATP	PG-O3B-PB-O1B
3	O	502[B]	ATP	PG-O3B-PB-O1B

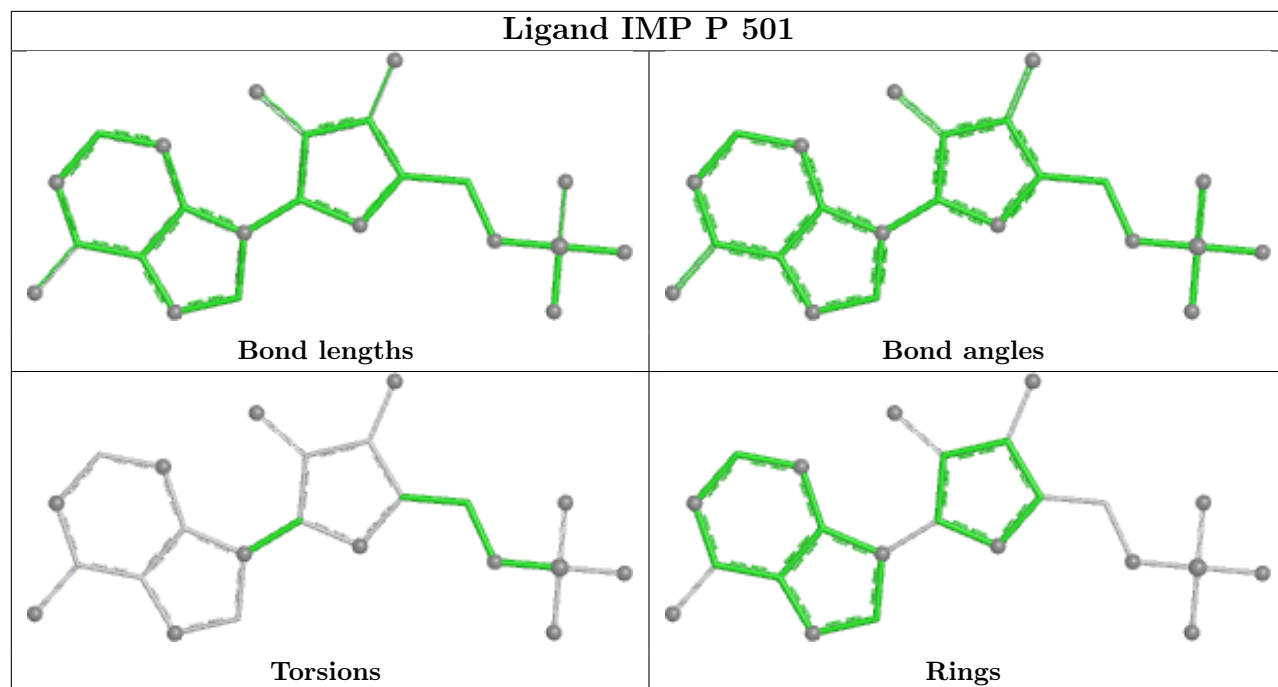
There are no ring outliers.

28 monomers are involved in 30 short contacts:

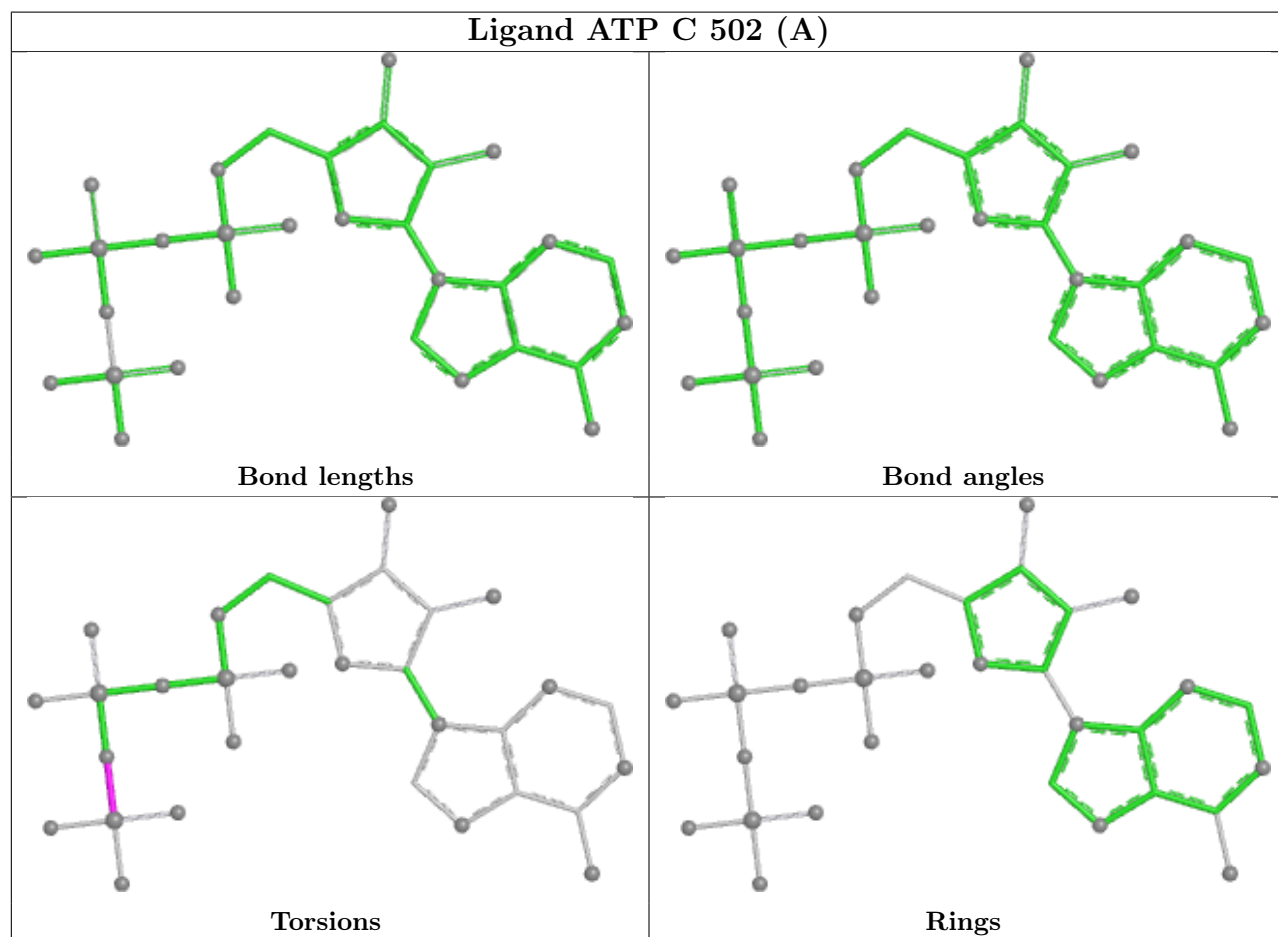
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	502[A]	ATP	1	0
3	C	502[B]	ATP	1	0
3	M	502[A]	ATP	1	0
3	M	502[B]	ATP	1	0
3	A	502[A]	ATP	1	0
3	A	502[B]	ATP	1	0
3	F	502[A]	ATP	1	0
3	F	502[B]	ATP	1	0
3	N	502[B]	ATP	1	0
3	I	502[A]	ATP	1	0
3	I	502[B]	ATP	1	0
3	L	502[A]	ATP	1	0
3	D	502[B]	ATP	1	0
3	L	502[B]	ATP	1	0
3	J	502[B]	ATP	1	0
3	D	502[A]	ATP	1	0
3	G	502[A]	ATP	1	0
3	G	502[B]	ATP	1	0
3	H	502[A]	ATP	1	0
3	E	502[A]	ATP	1	0
3	E	502[B]	ATP	1	0
3	K	502[A]	ATP	2	0
3	K	502[B]	ATP	2	0
3	O	502[A]	ATP	1	0
3	B	502[A]	ATP	1	0
3	B	502[B]	ATP	1	0
3	P	502[A]	ATP	1	0
3	J	502[A]	ATP	1	0

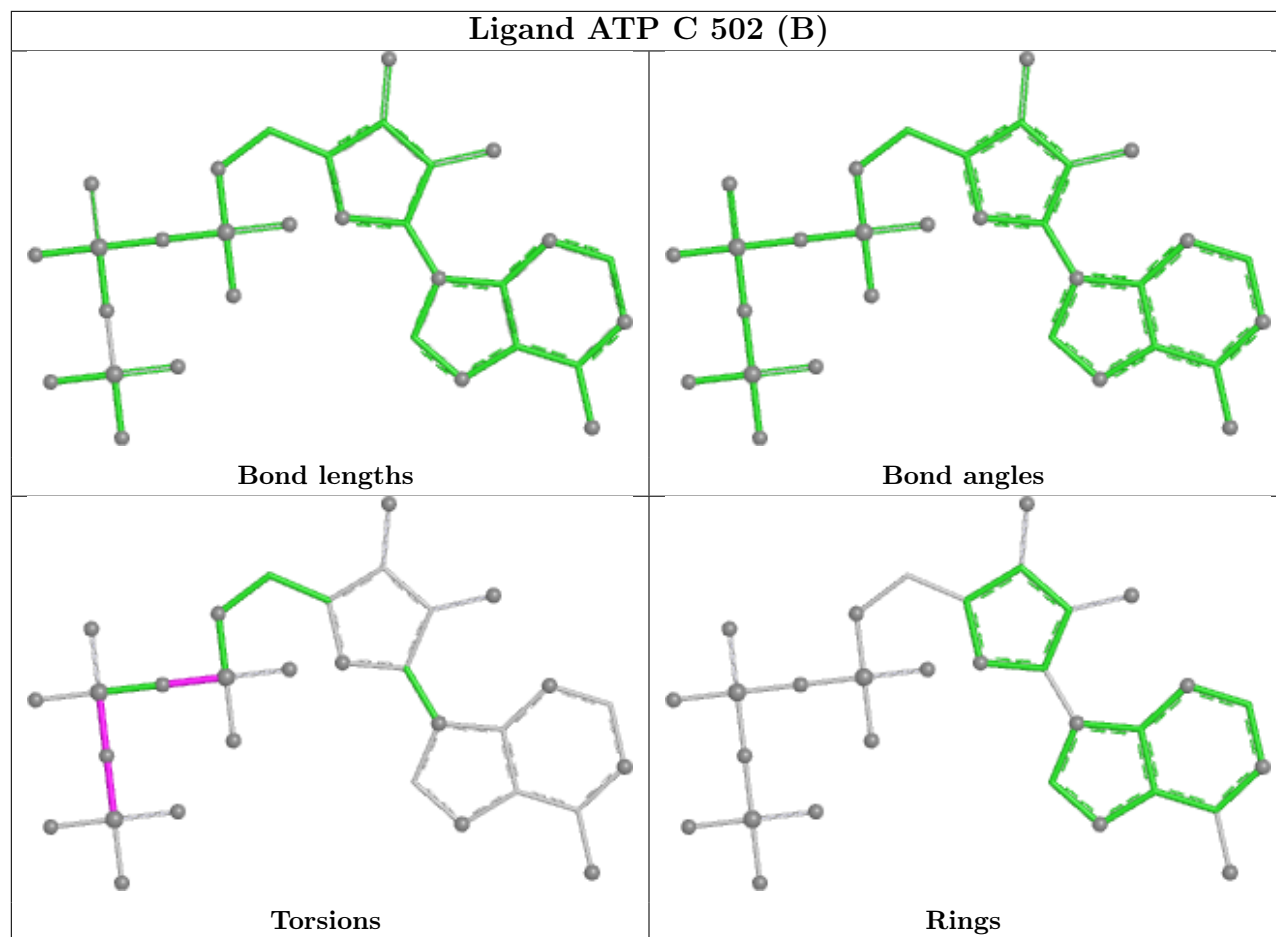
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

Ligand IMP P 501

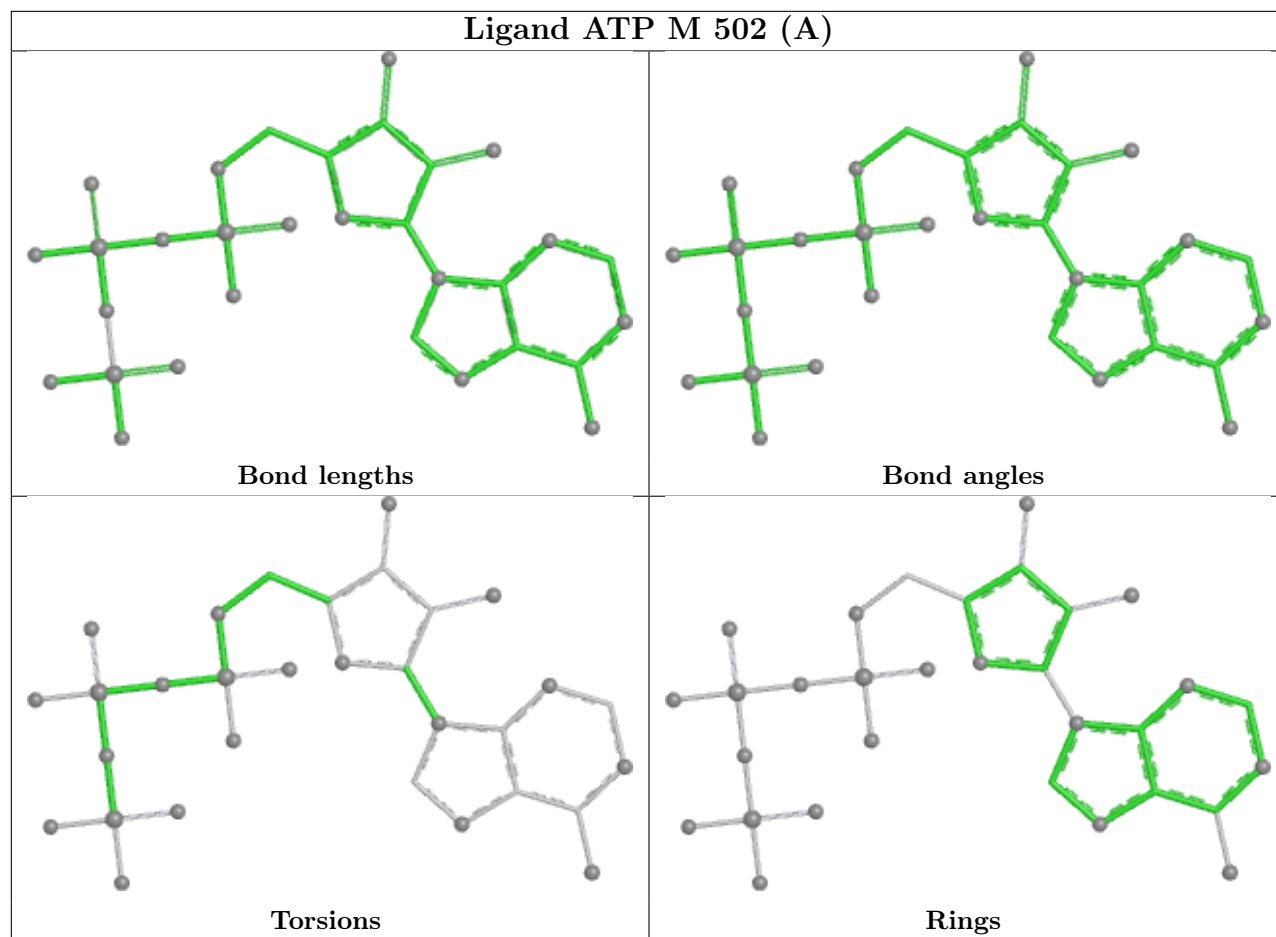


Ligand ATP C 502 (A)

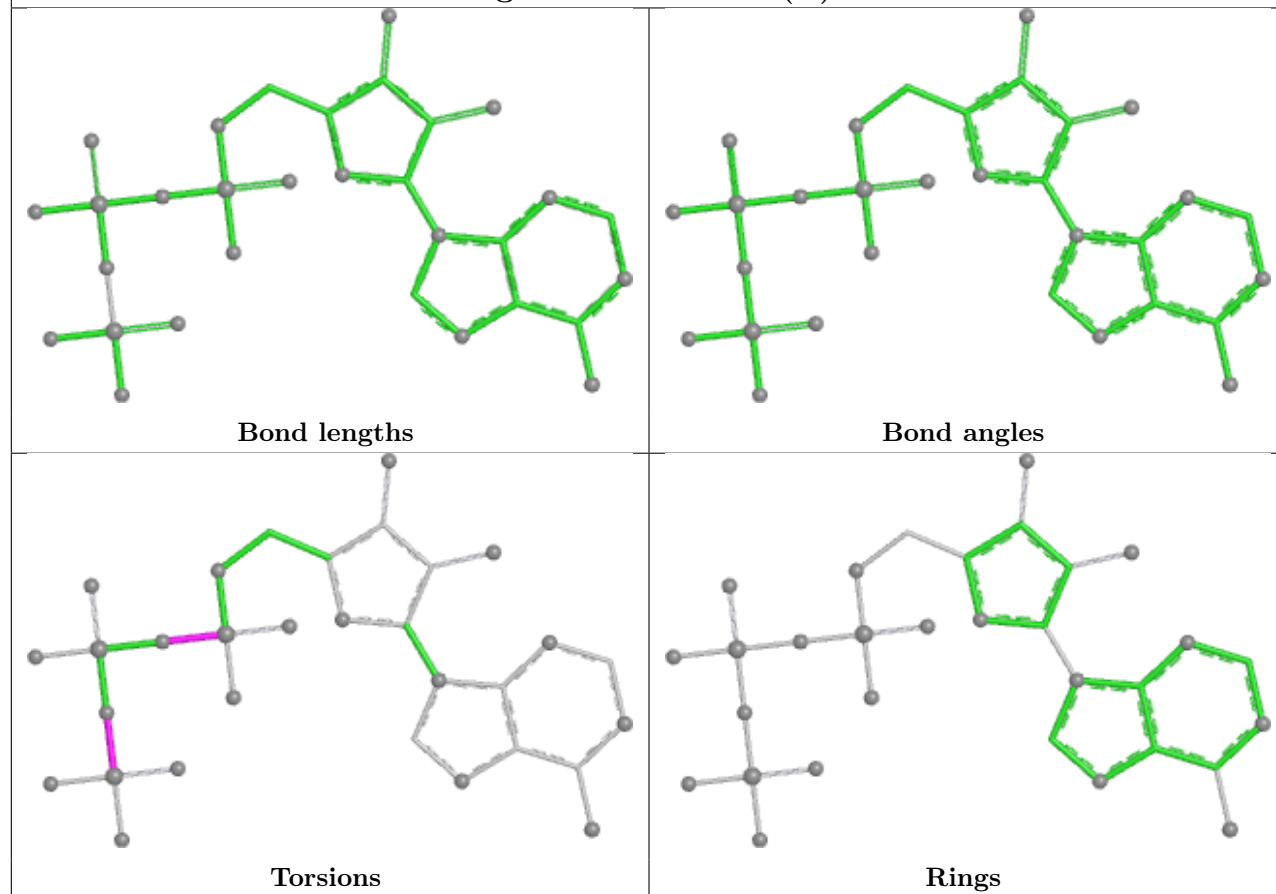




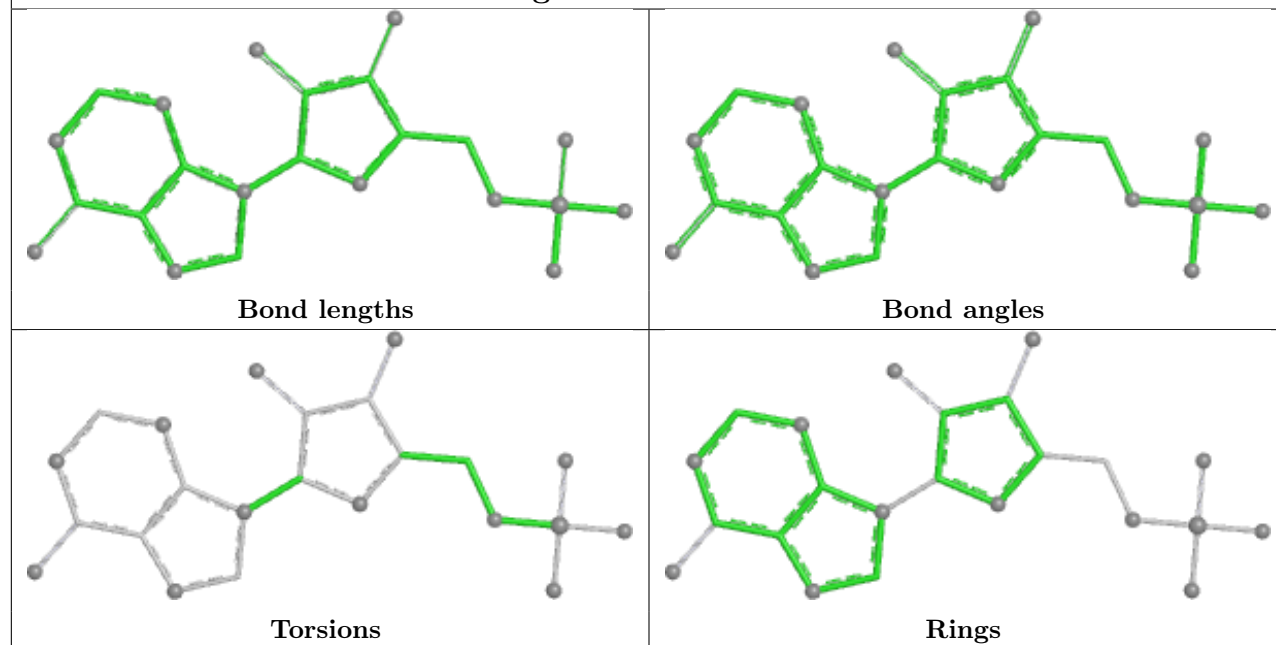
Ligand ATP M 502 (A)



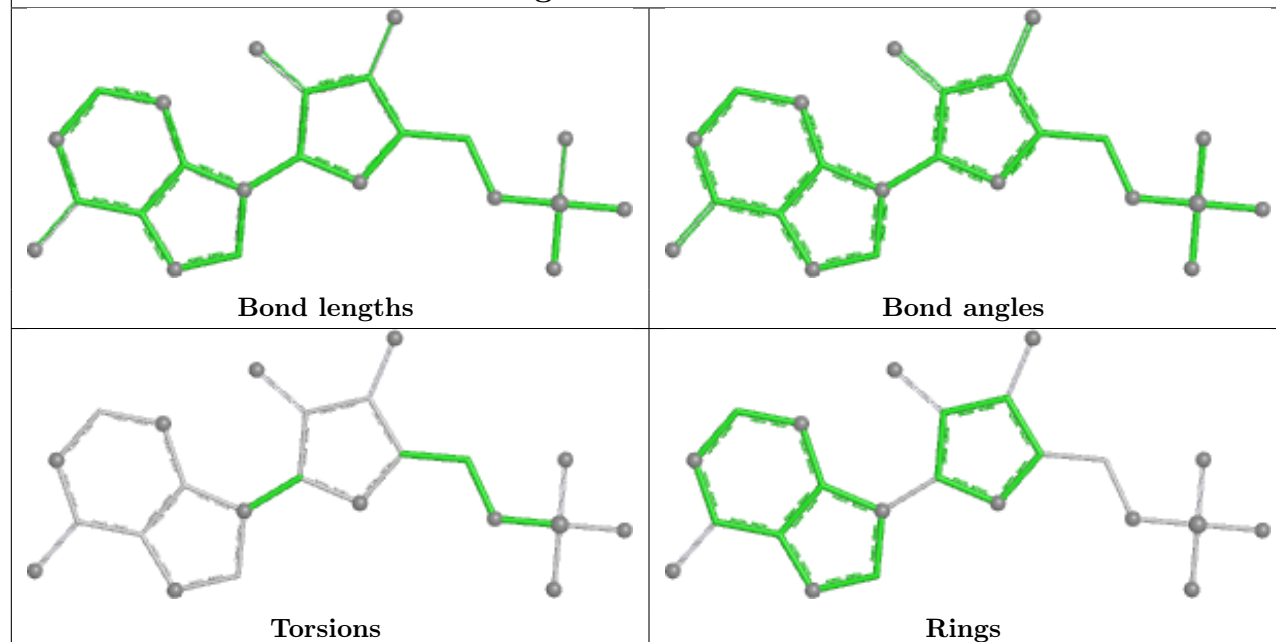
Ligand ATP M 502 (B)



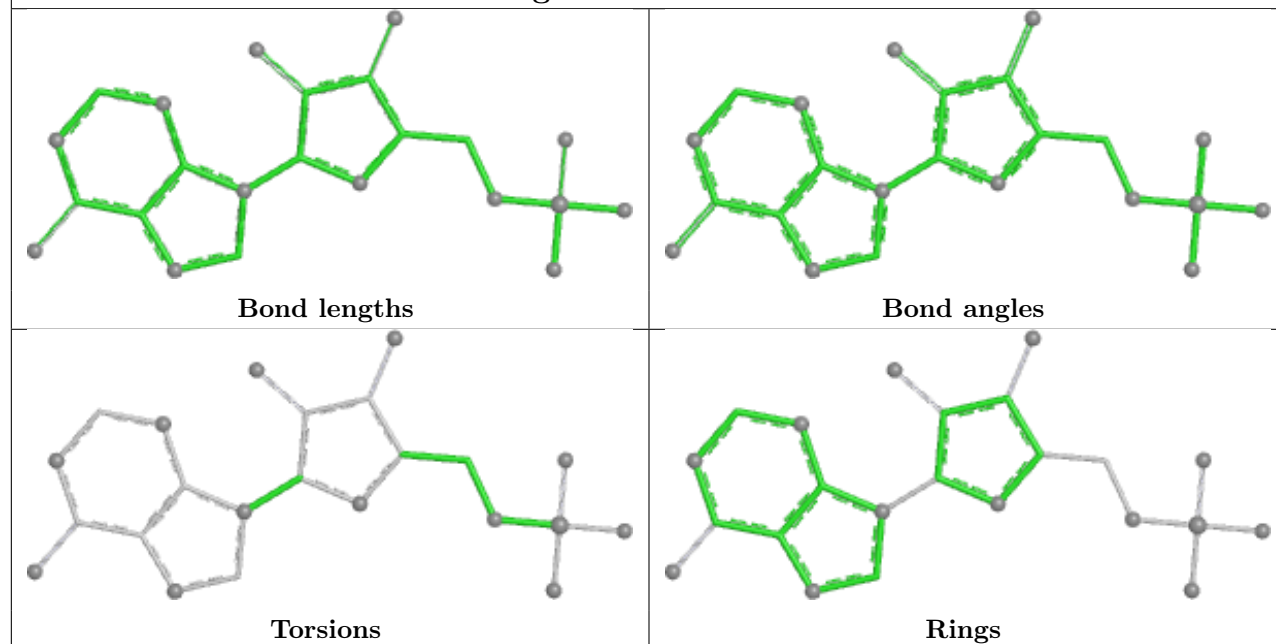
Ligand IMP H 501

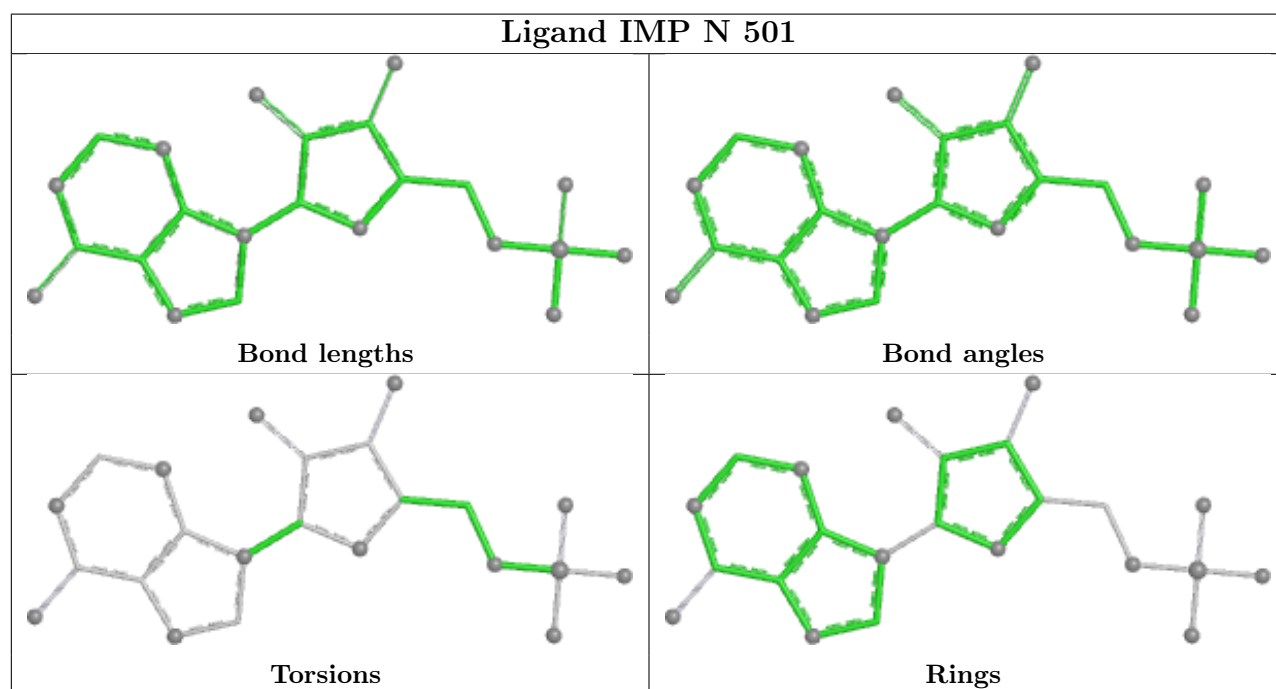
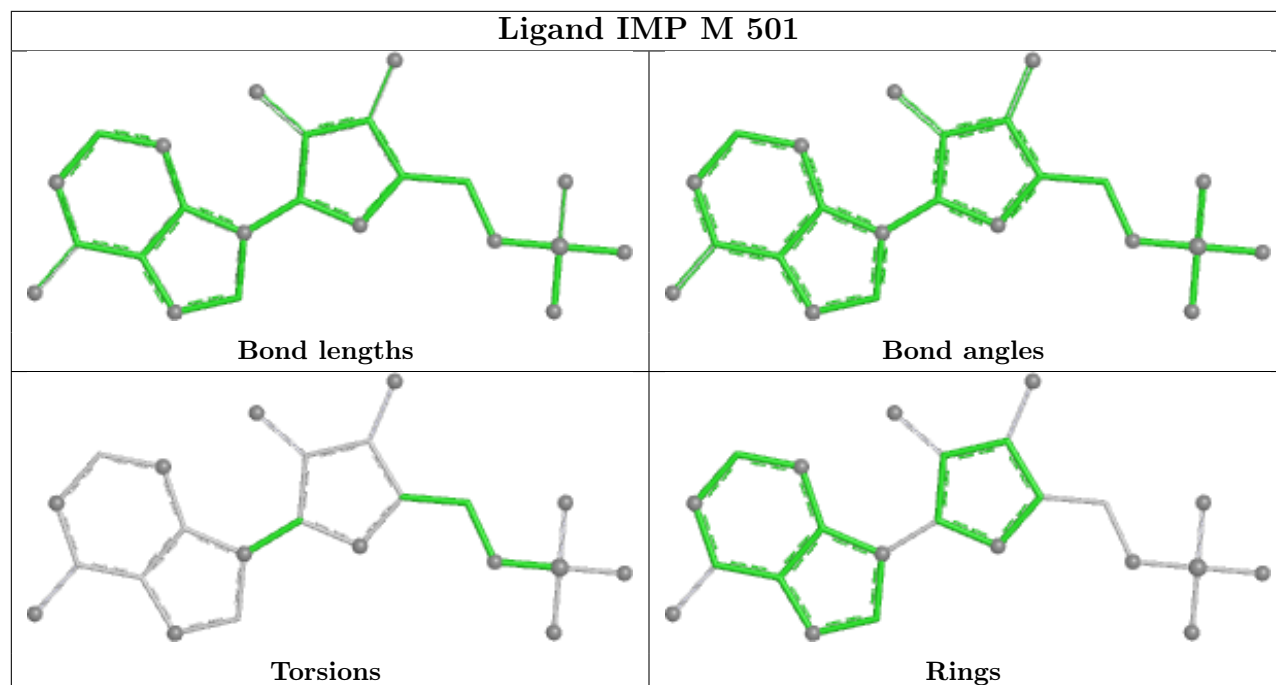


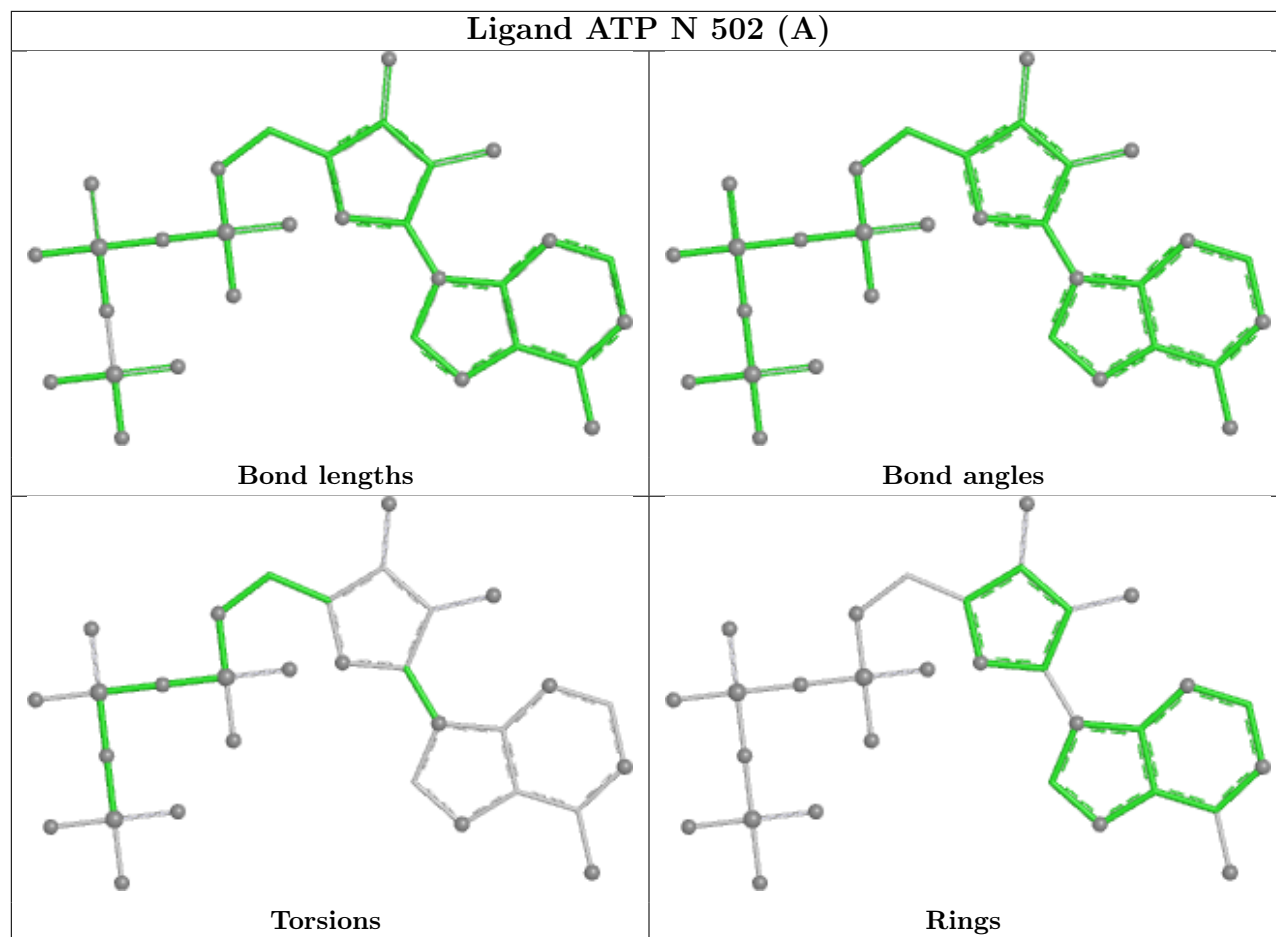
Ligand IMP E 501

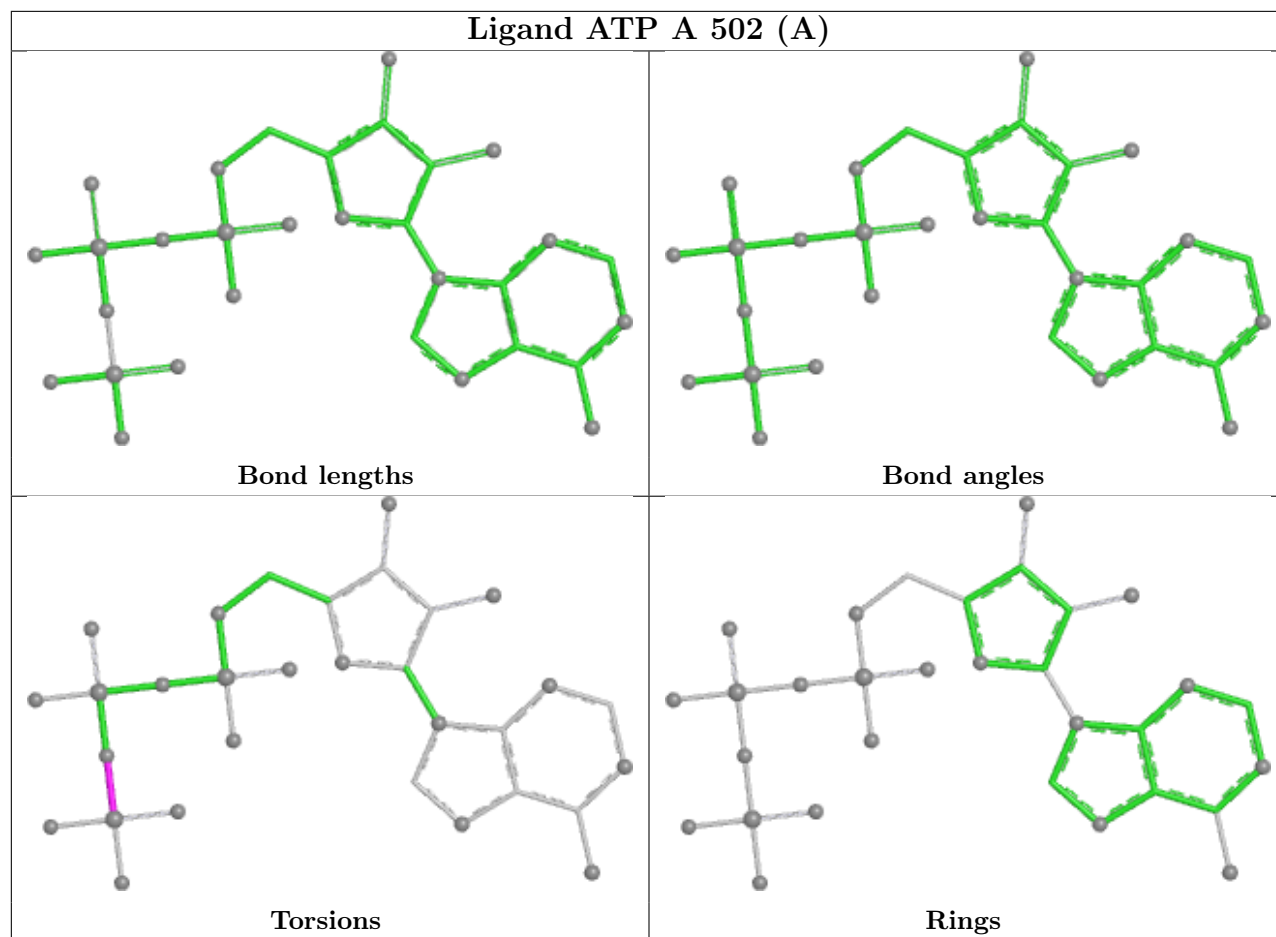


Ligand IMP D 501

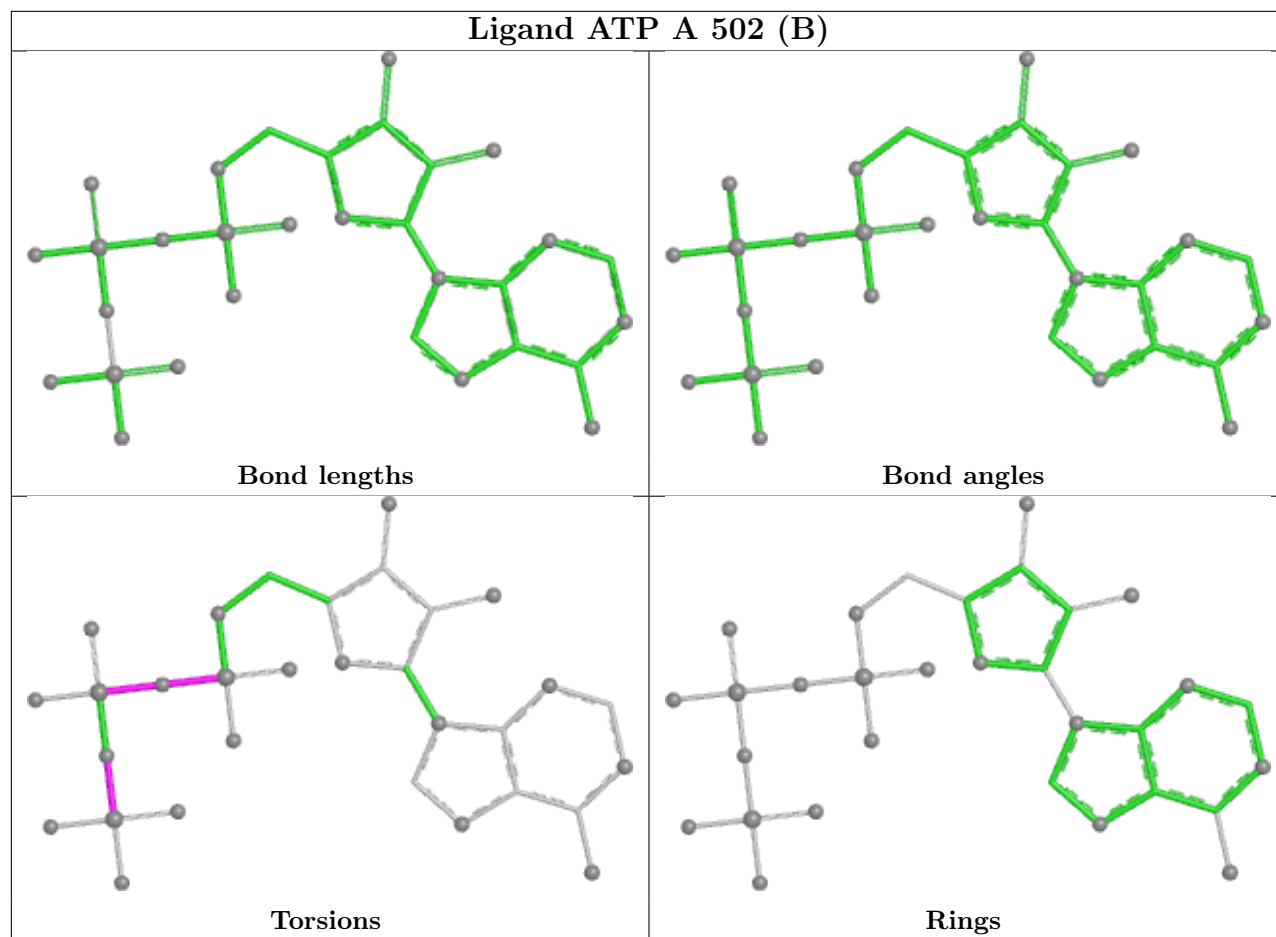


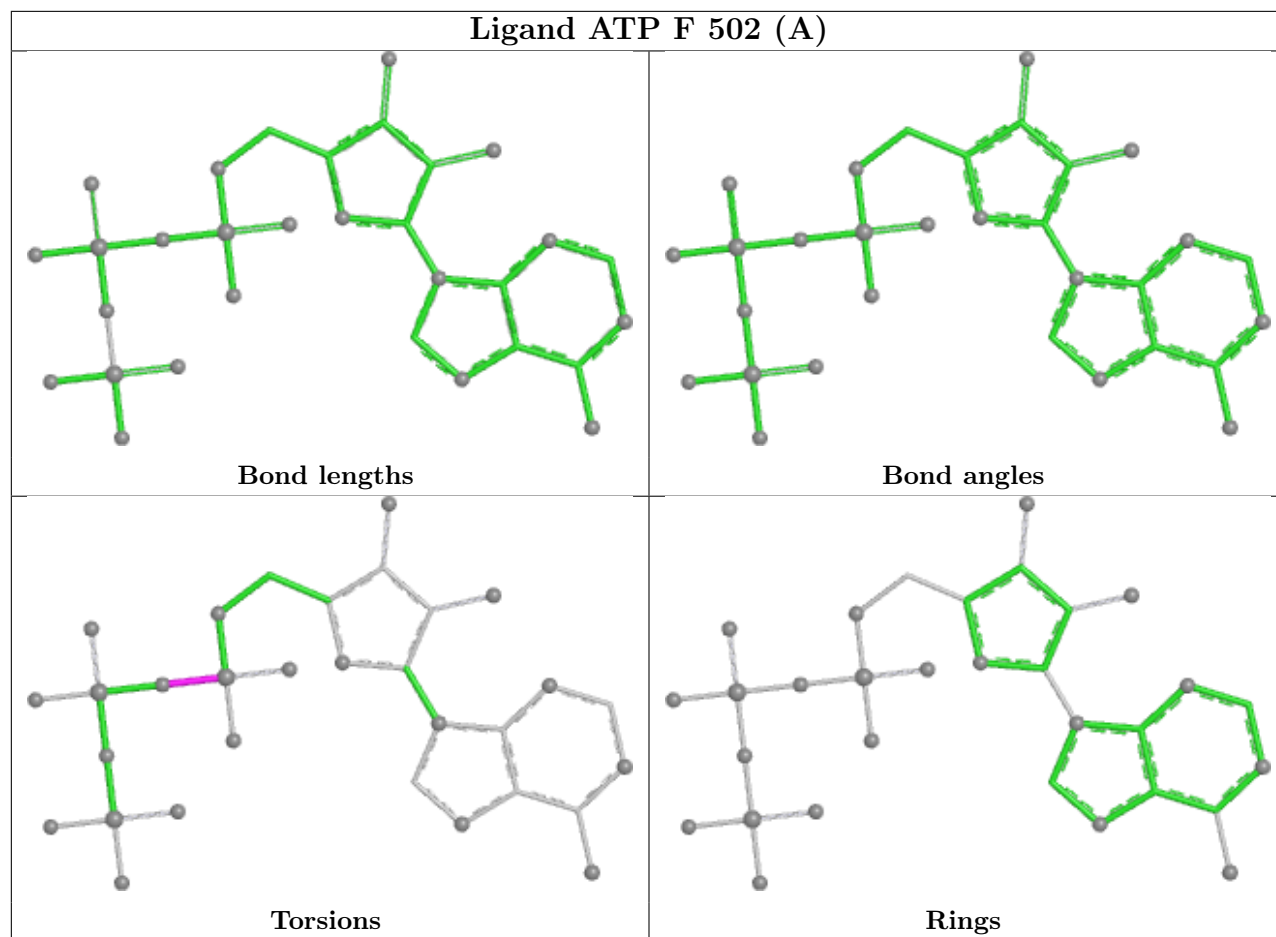


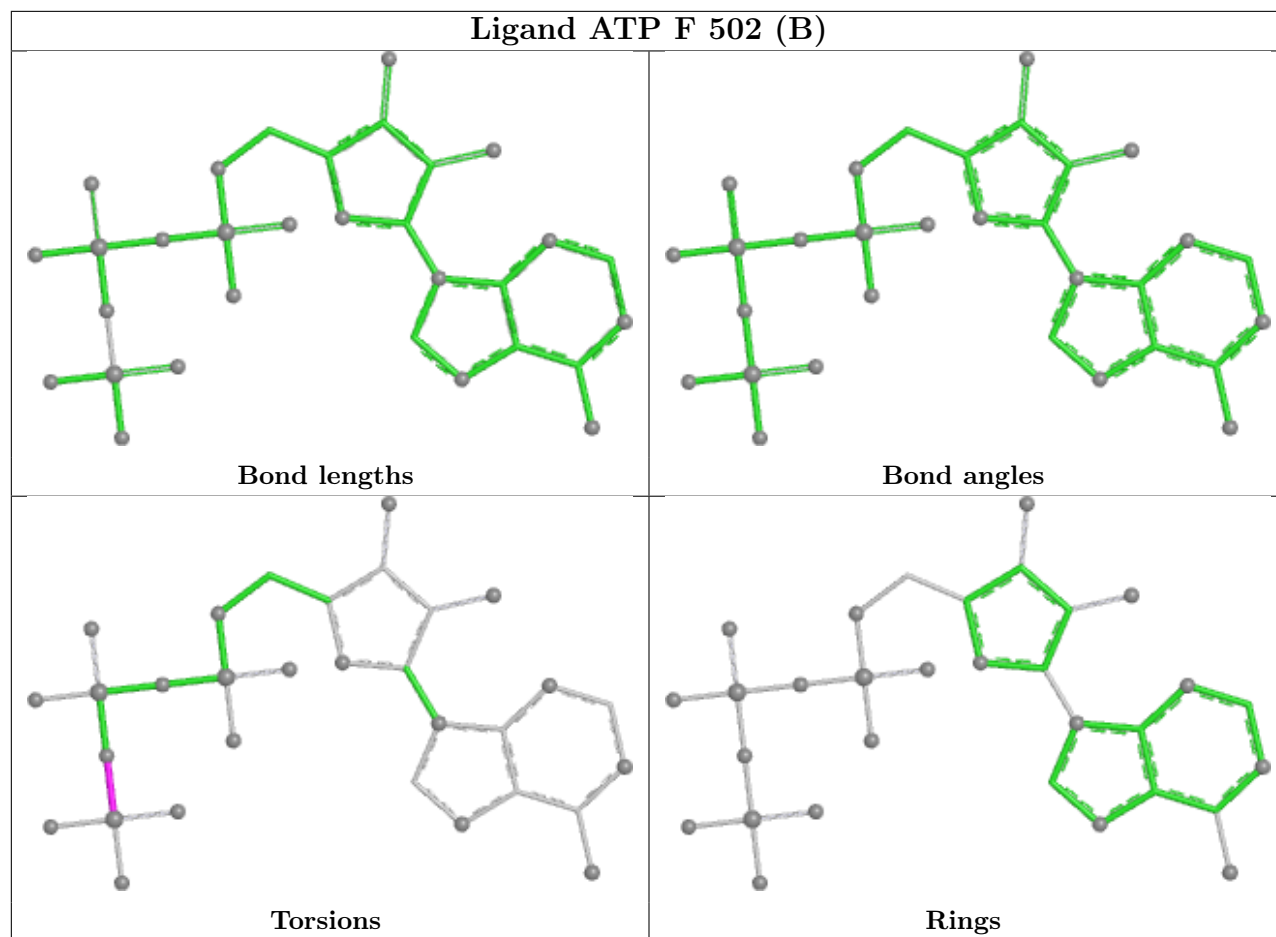




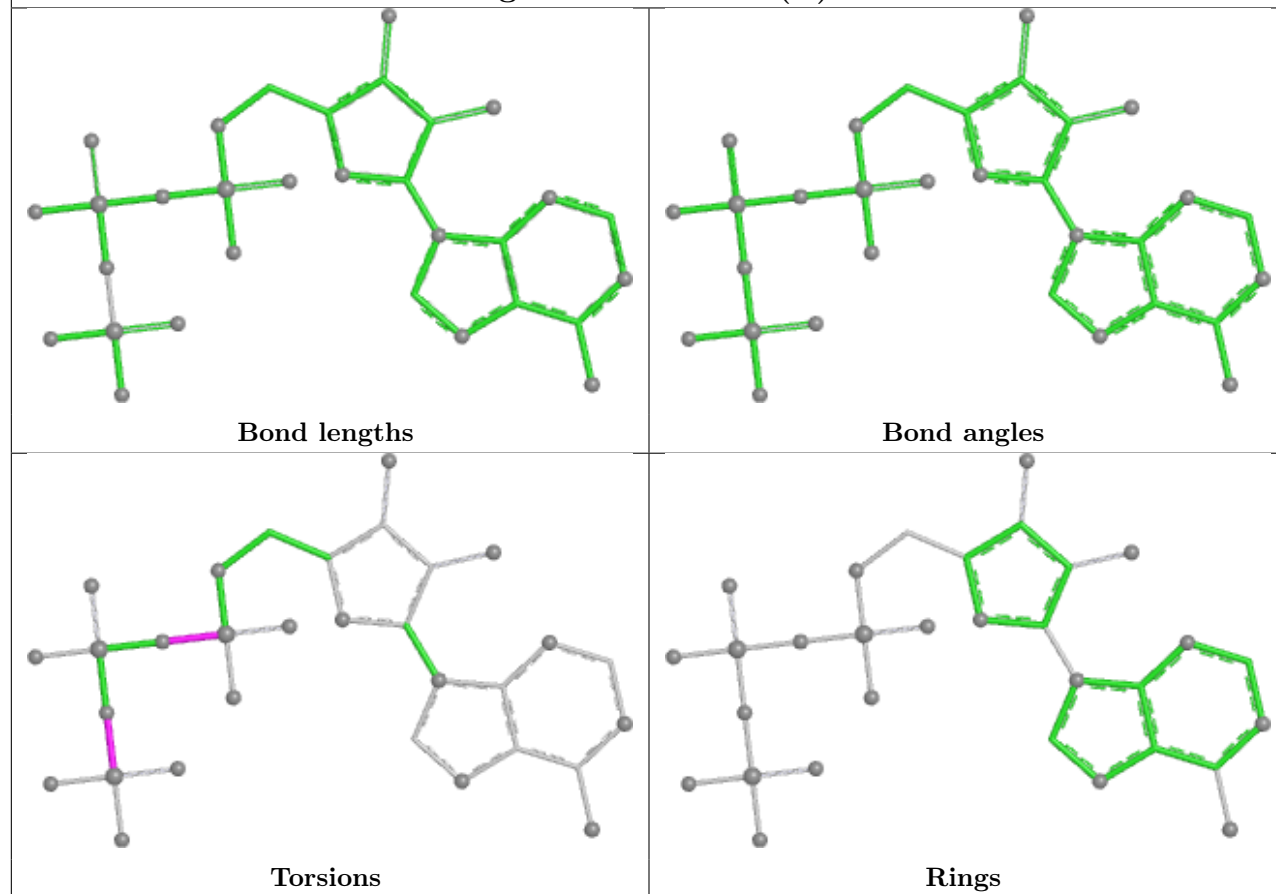
Ligand ATP A 502 (B)



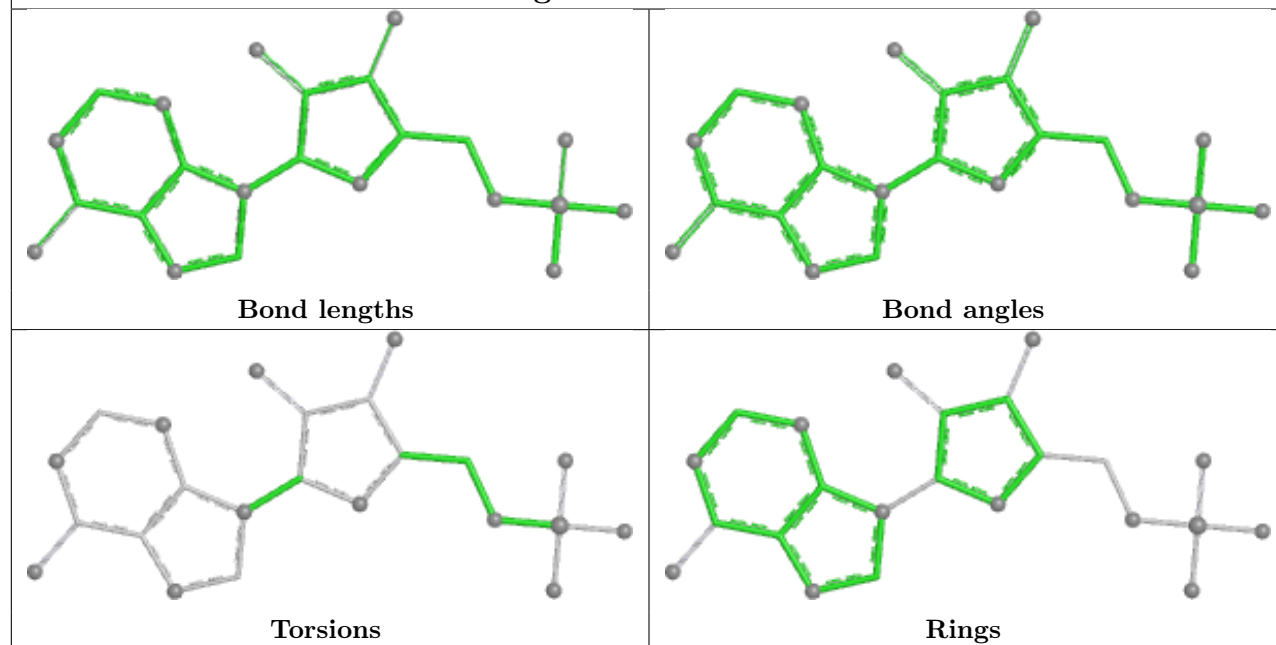




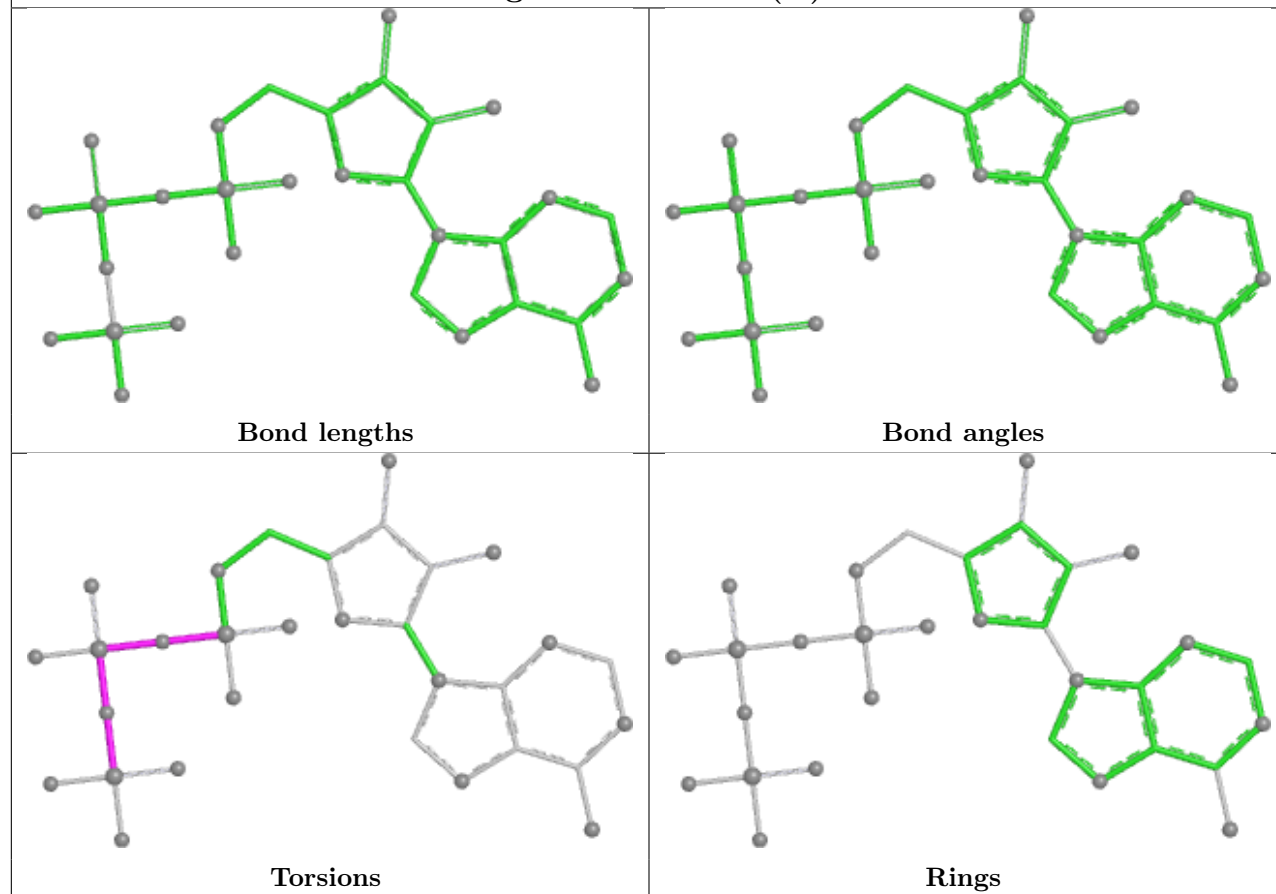
Ligand ATP N 502 (B)



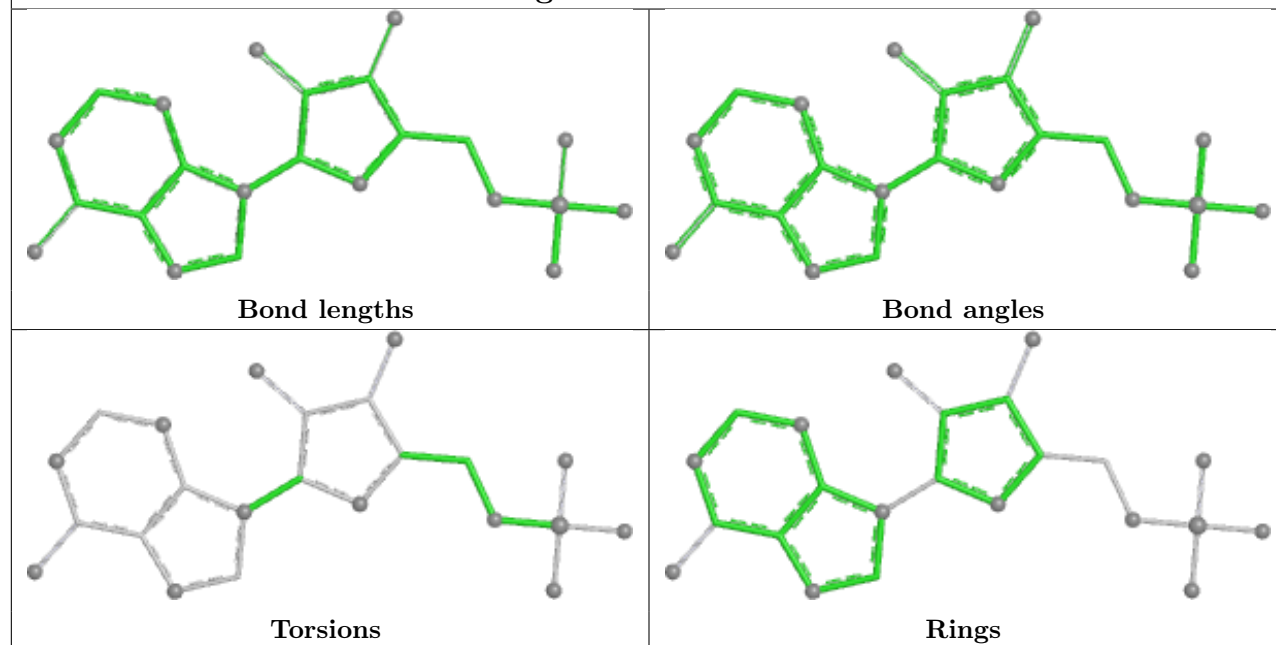
Ligand IMP C 501



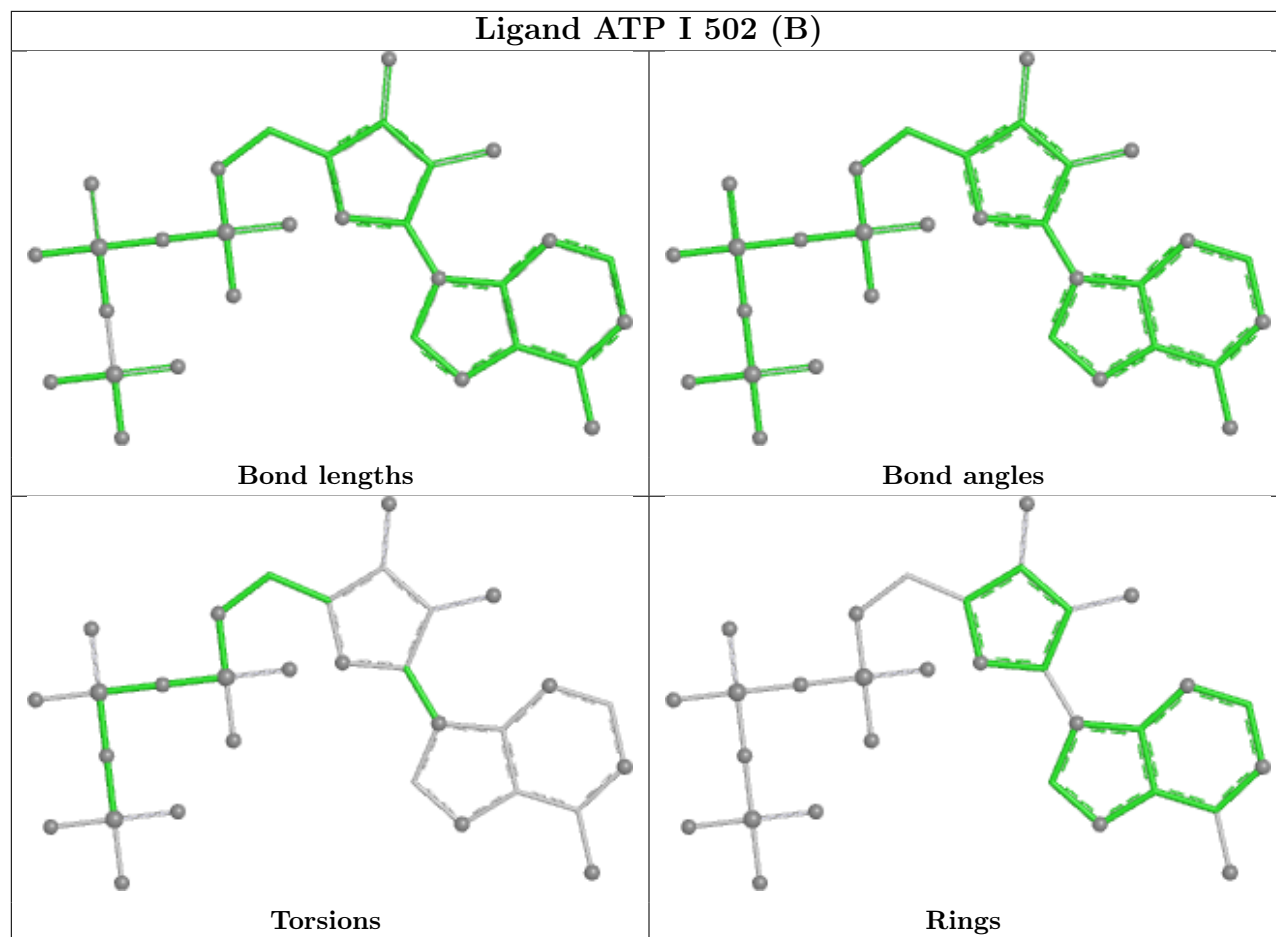
Ligand ATP I 502 (A)

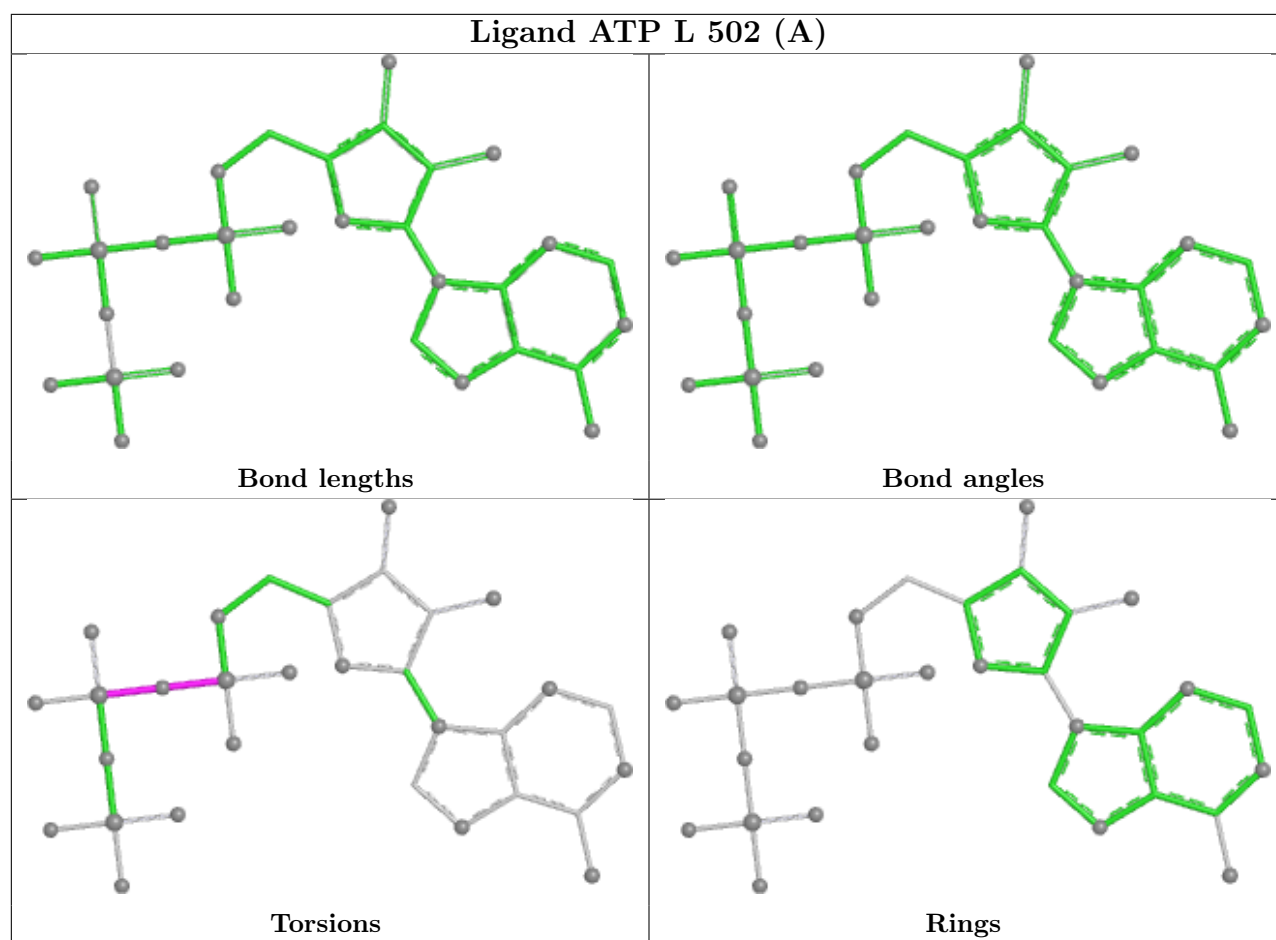


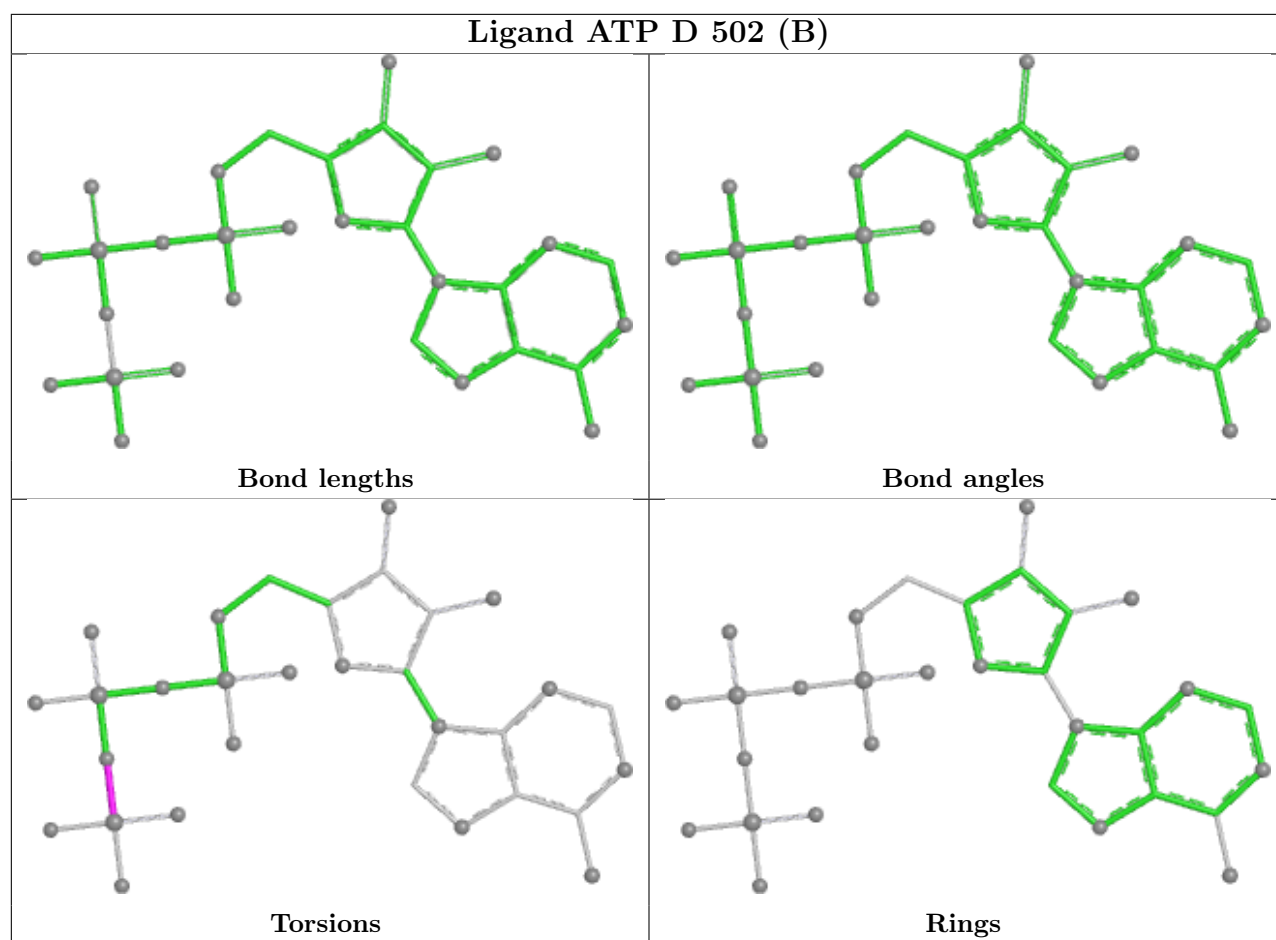
Ligand IMP L 501



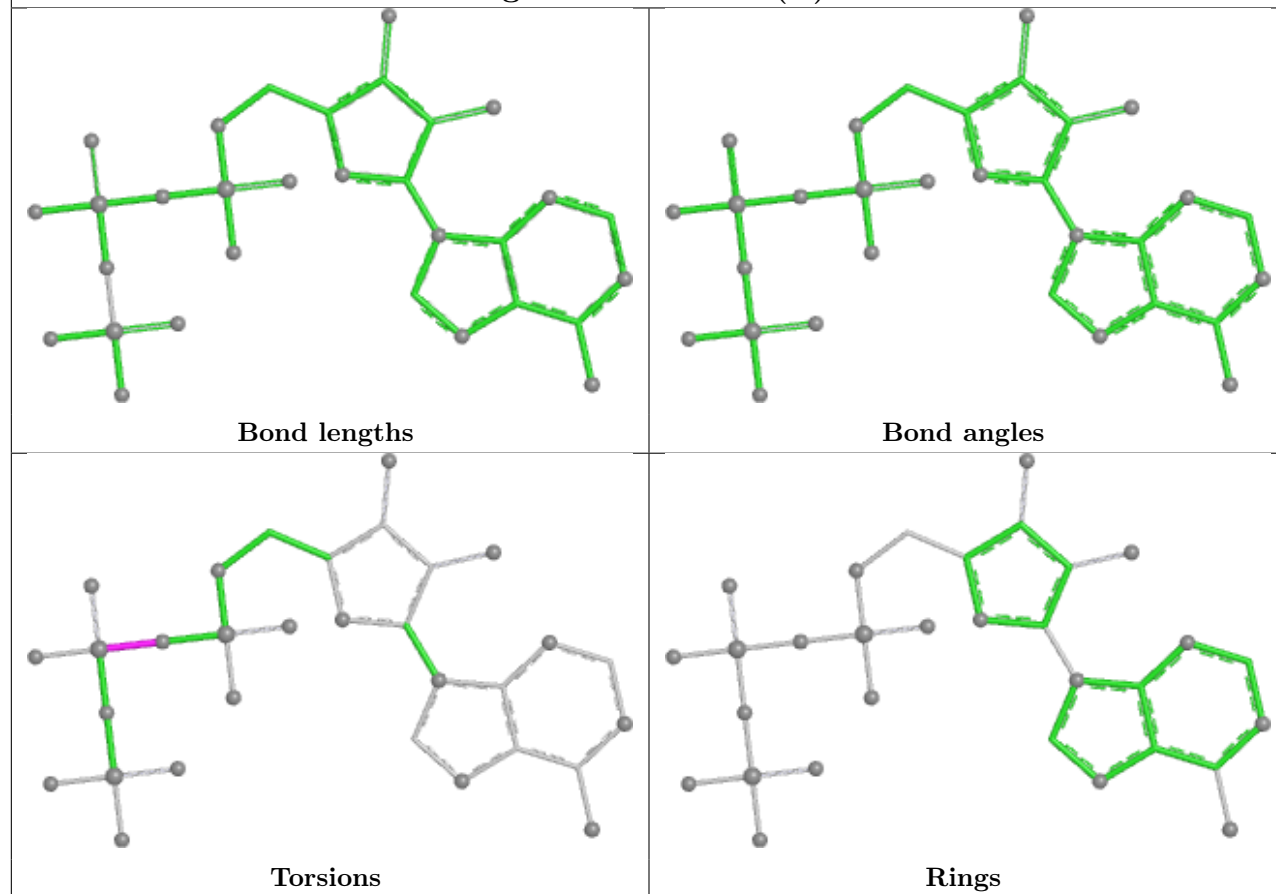
Ligand ATP I 502 (B)



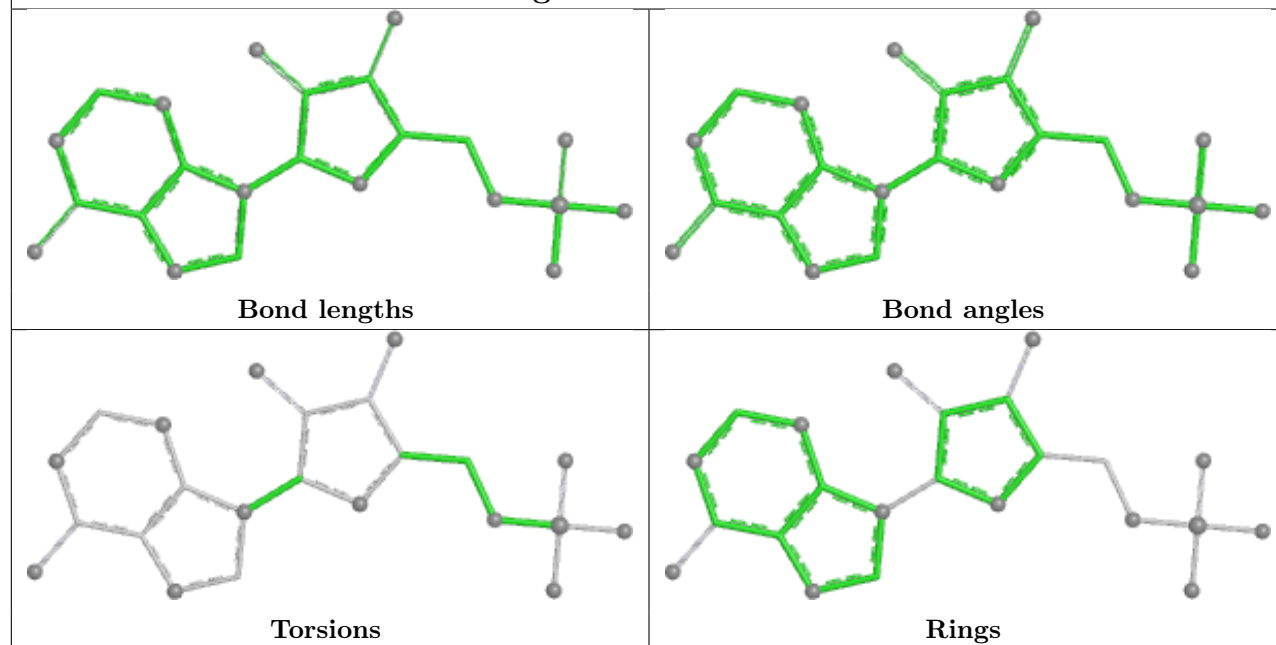




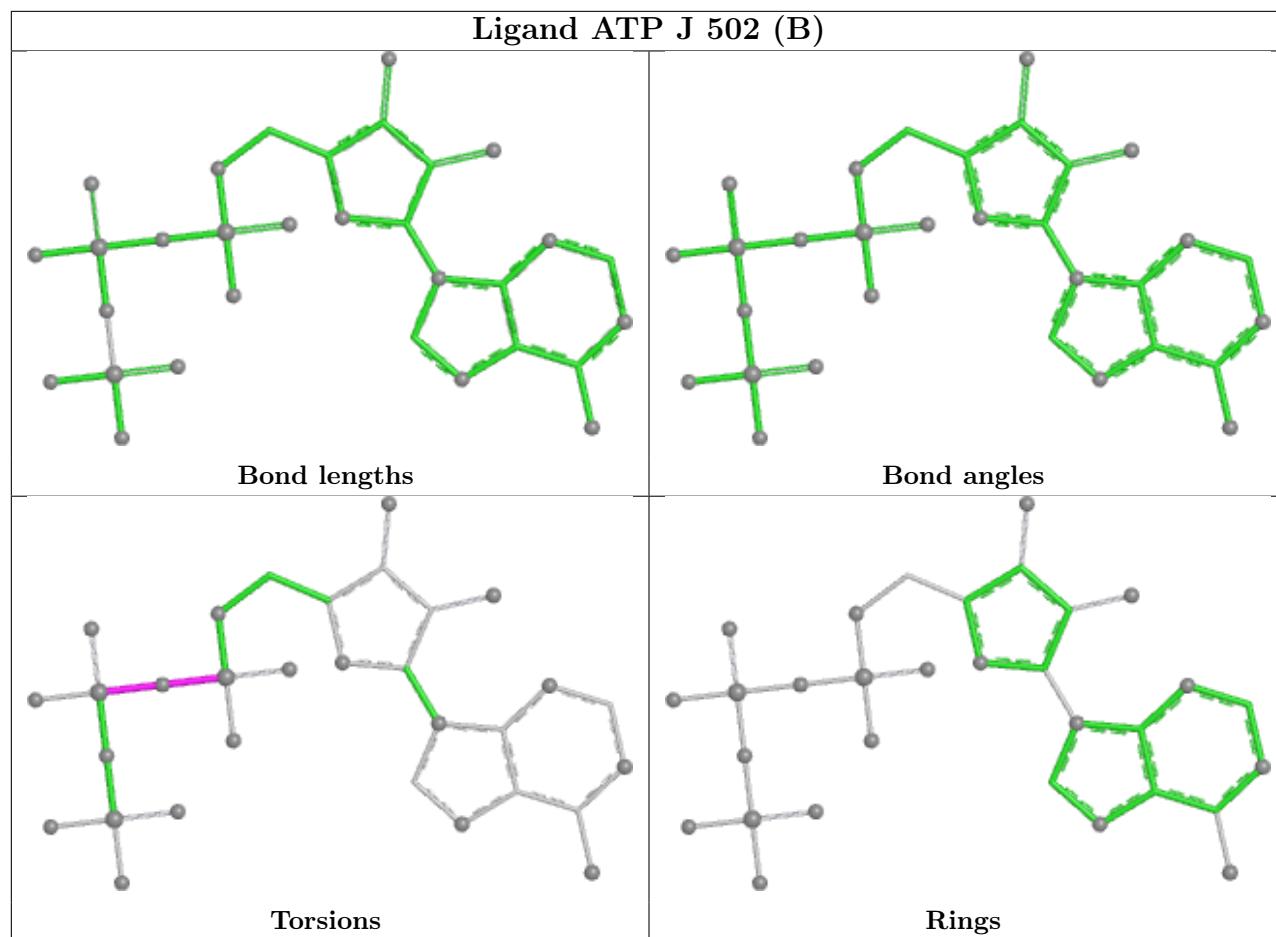
Ligand ATP L 502 (B)



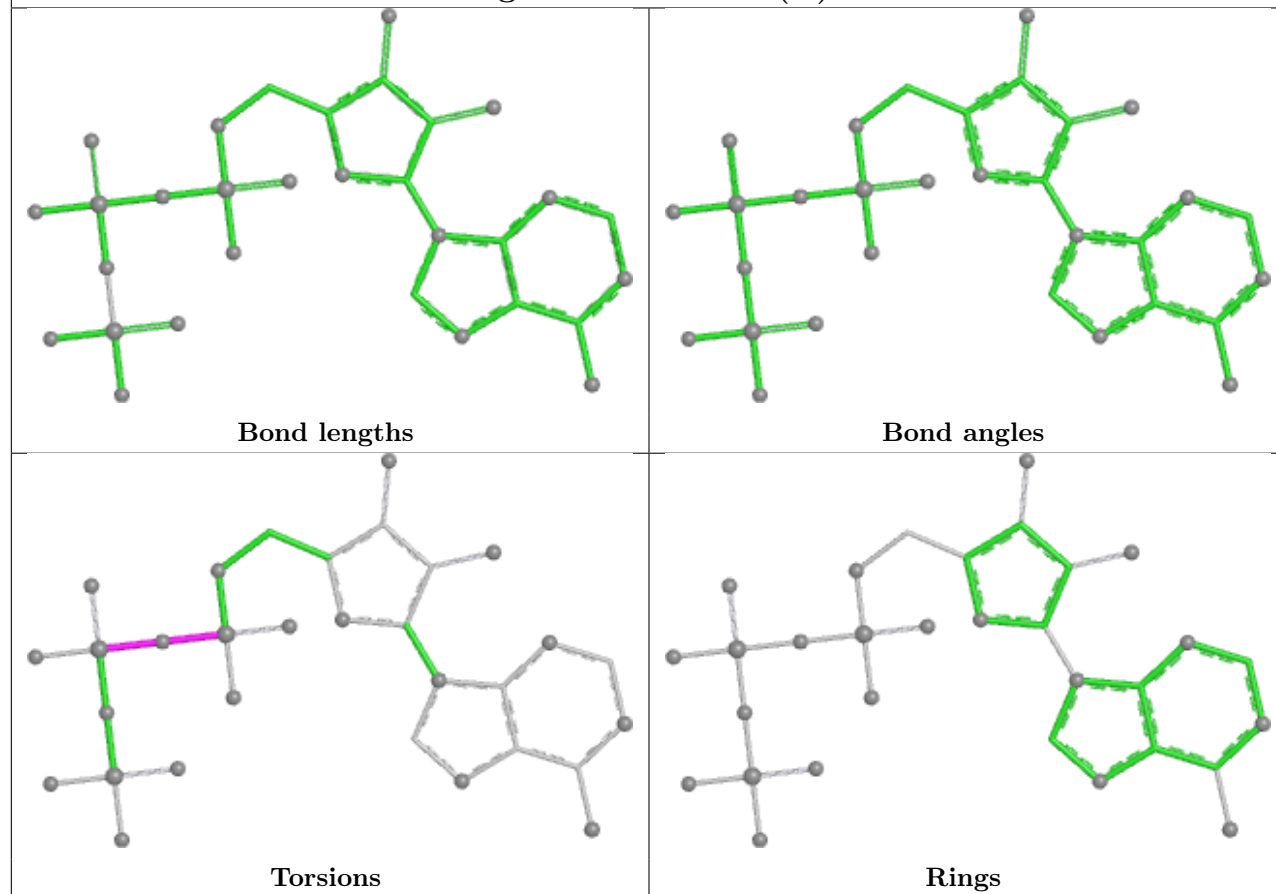
Ligand IMP A 501



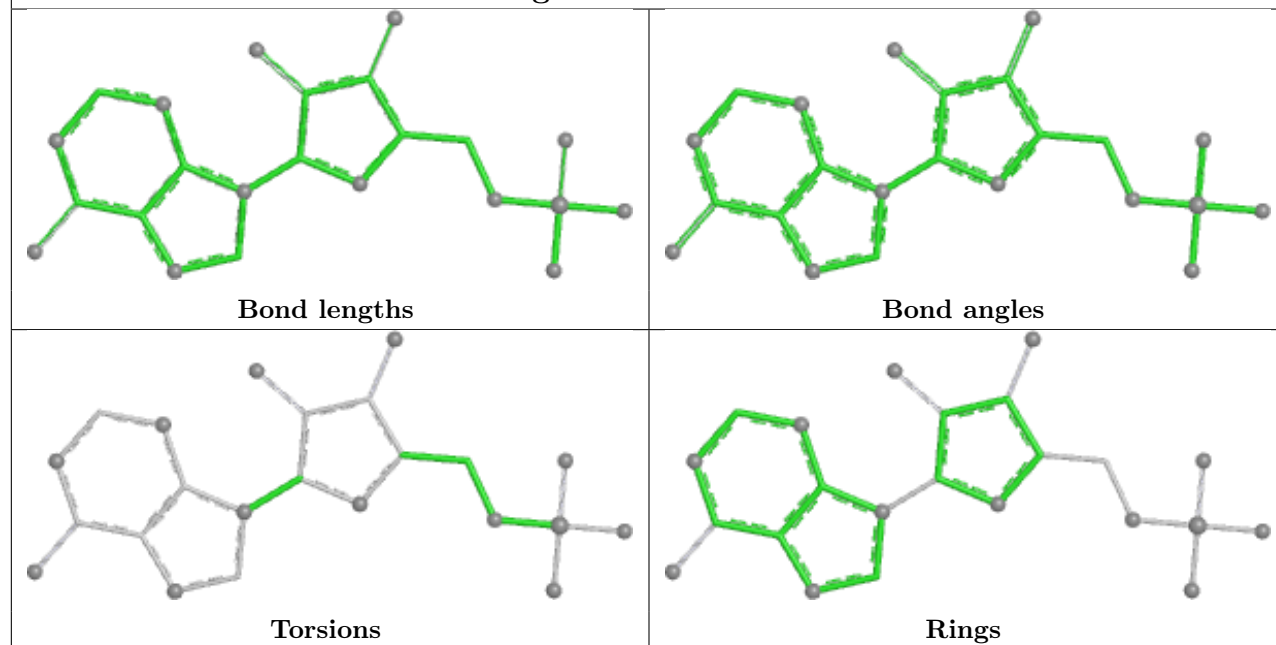
Ligand ATP J 502 (B)



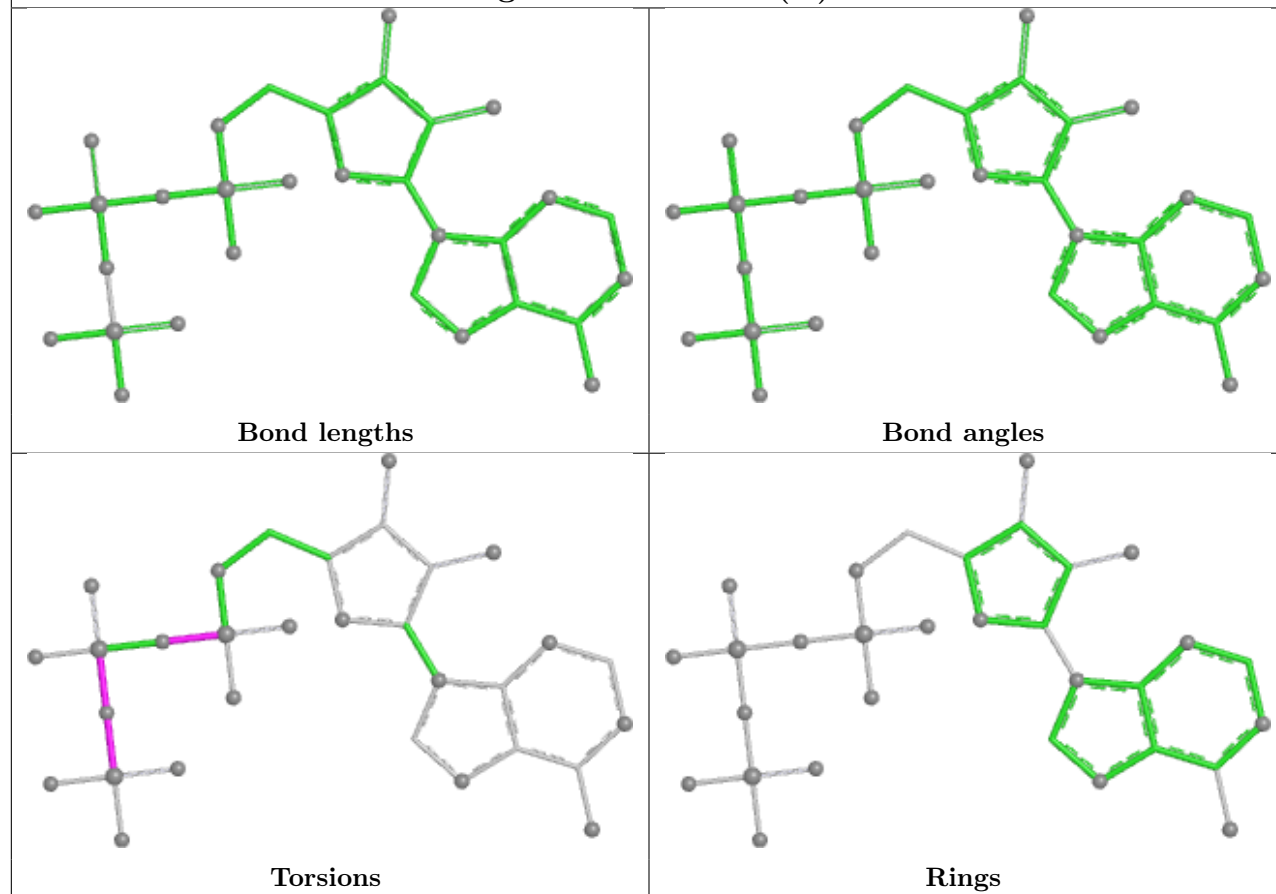
Ligand ATP D 502 (A)



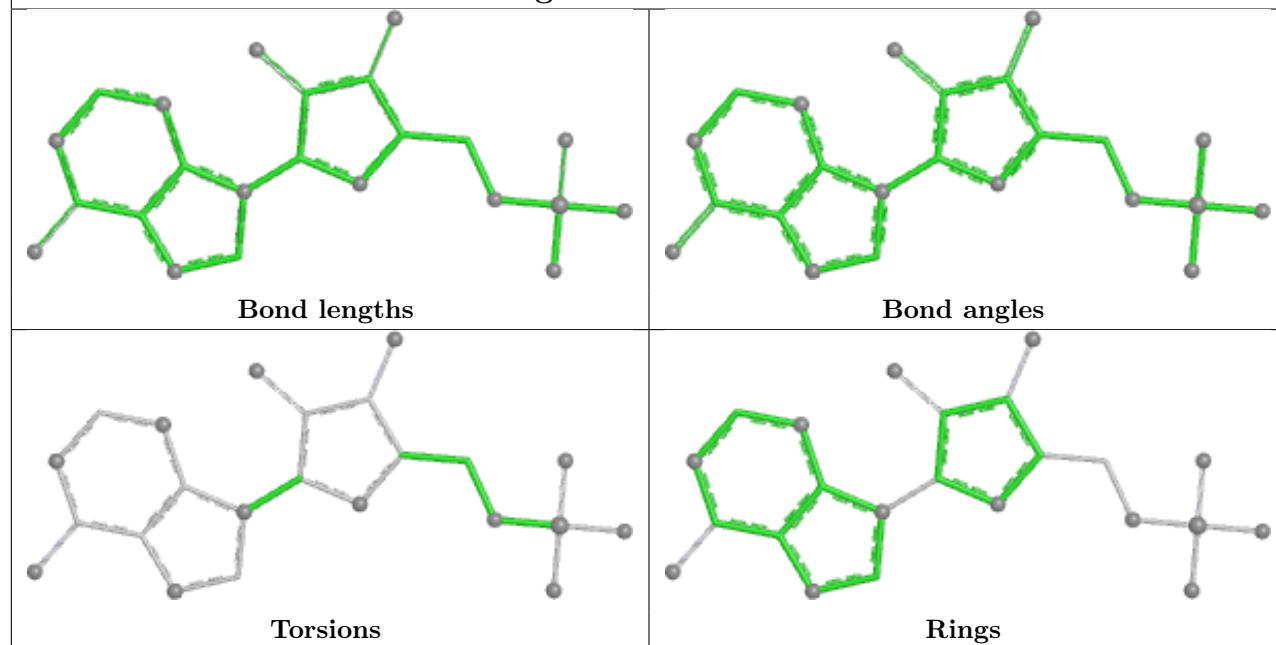
Ligand IMP G 501

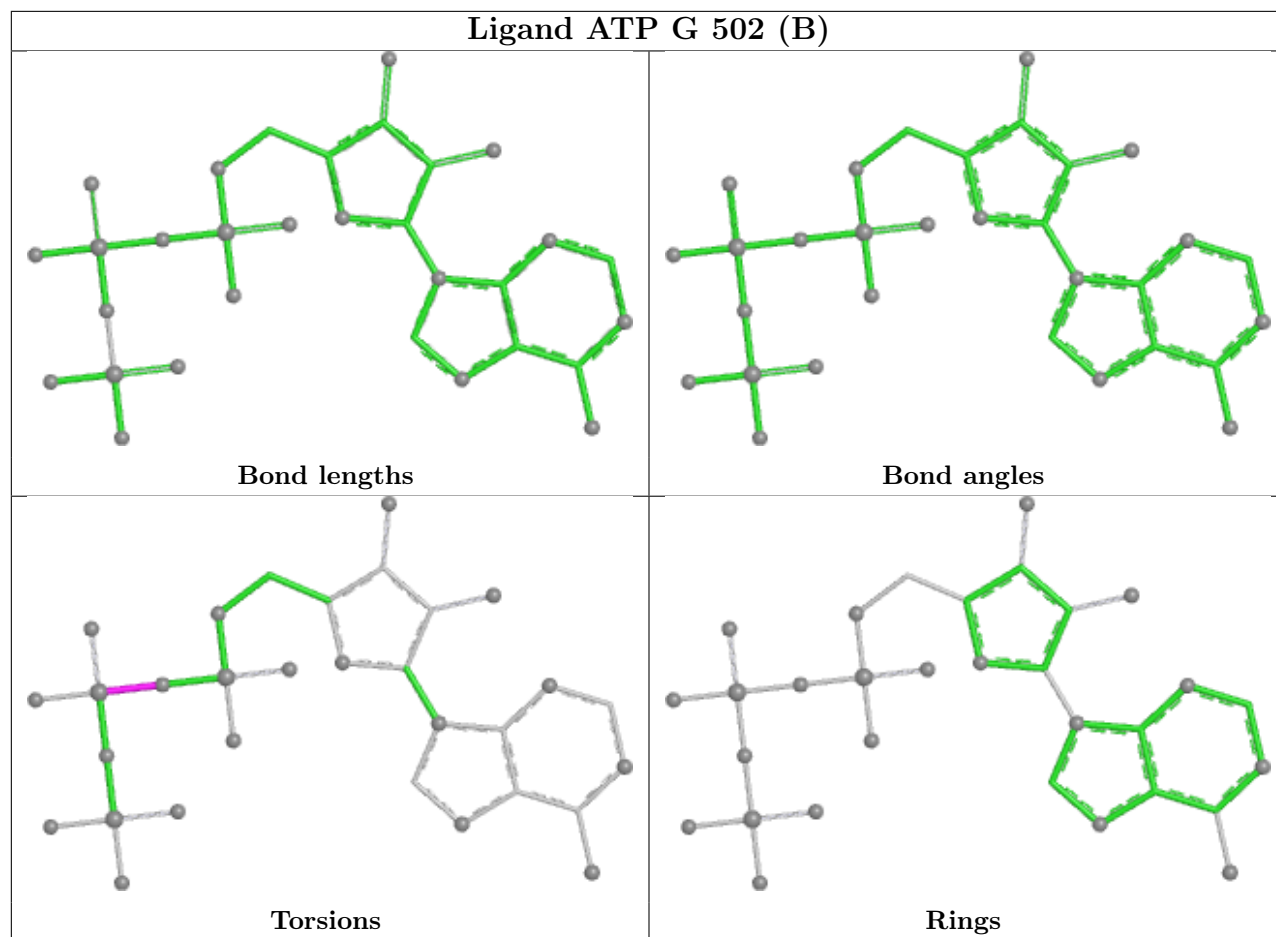


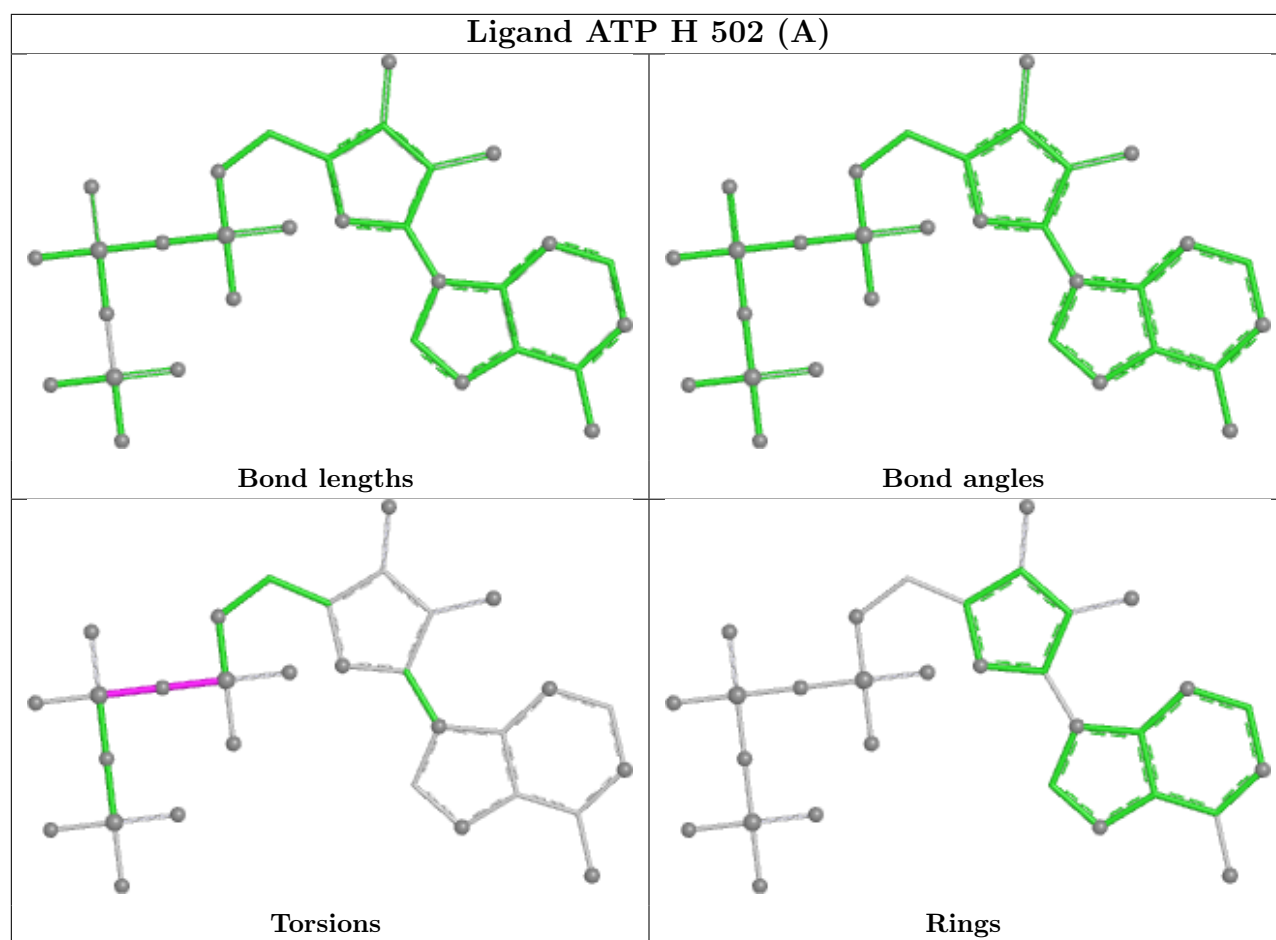
Ligand ATP G 502 (A)

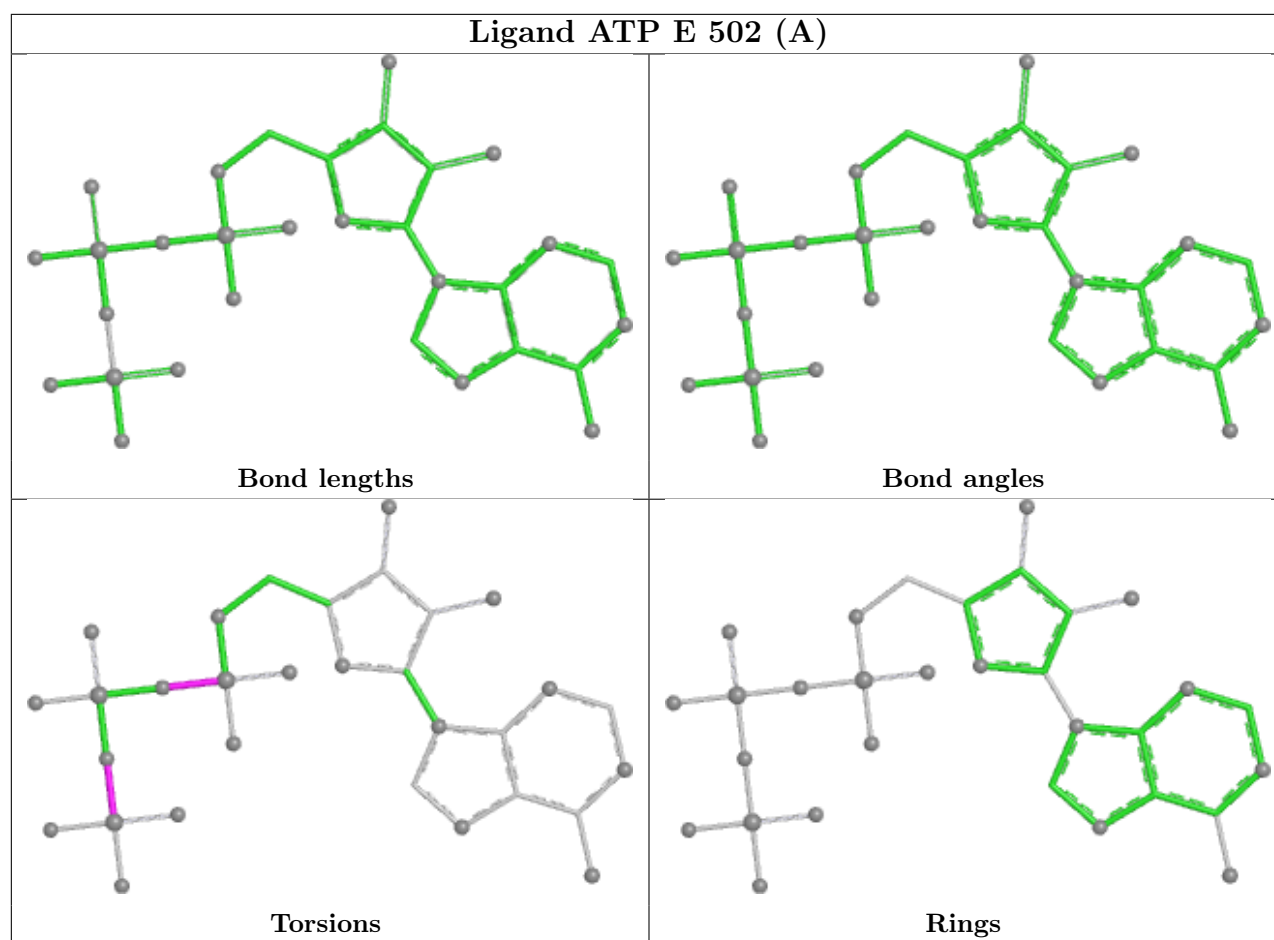


Ligand IMP F 501

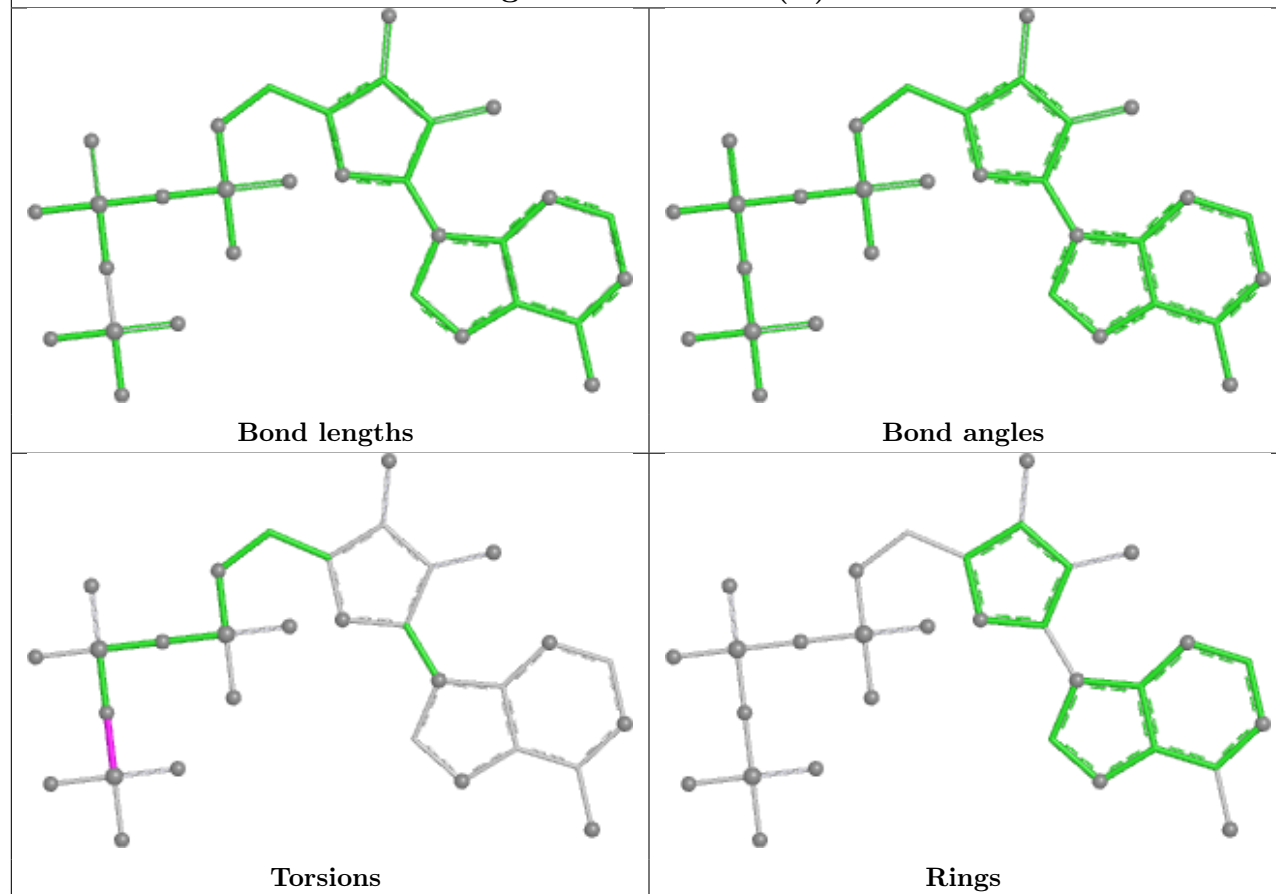




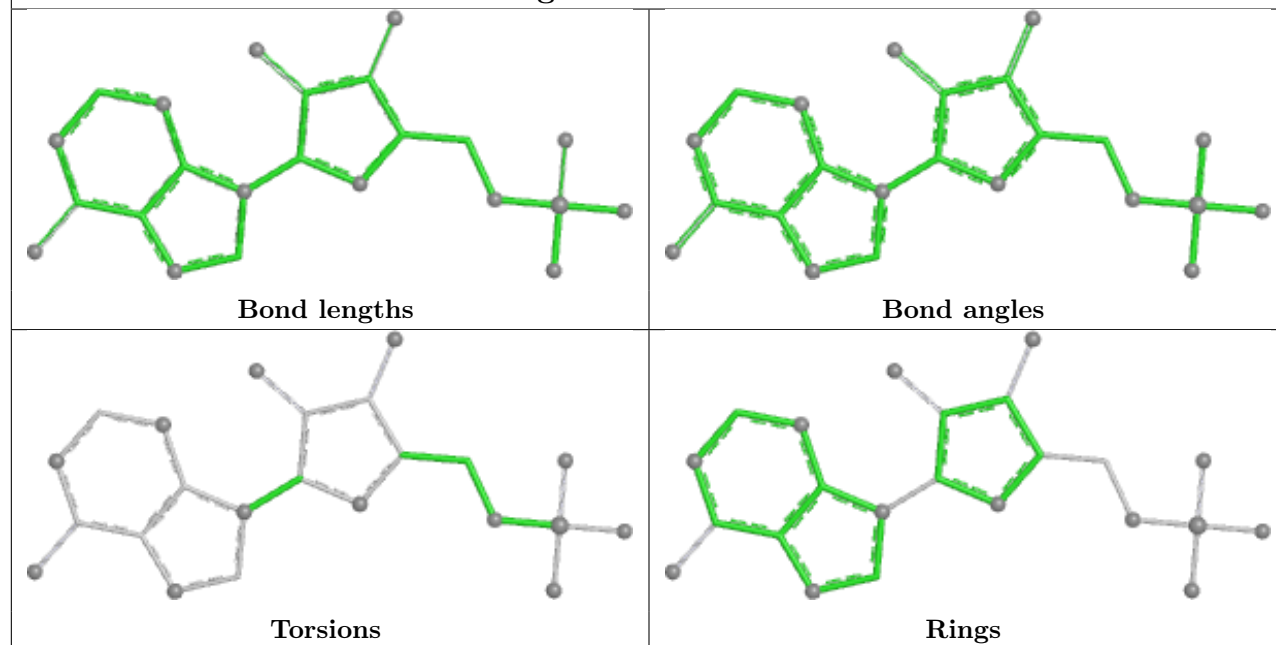




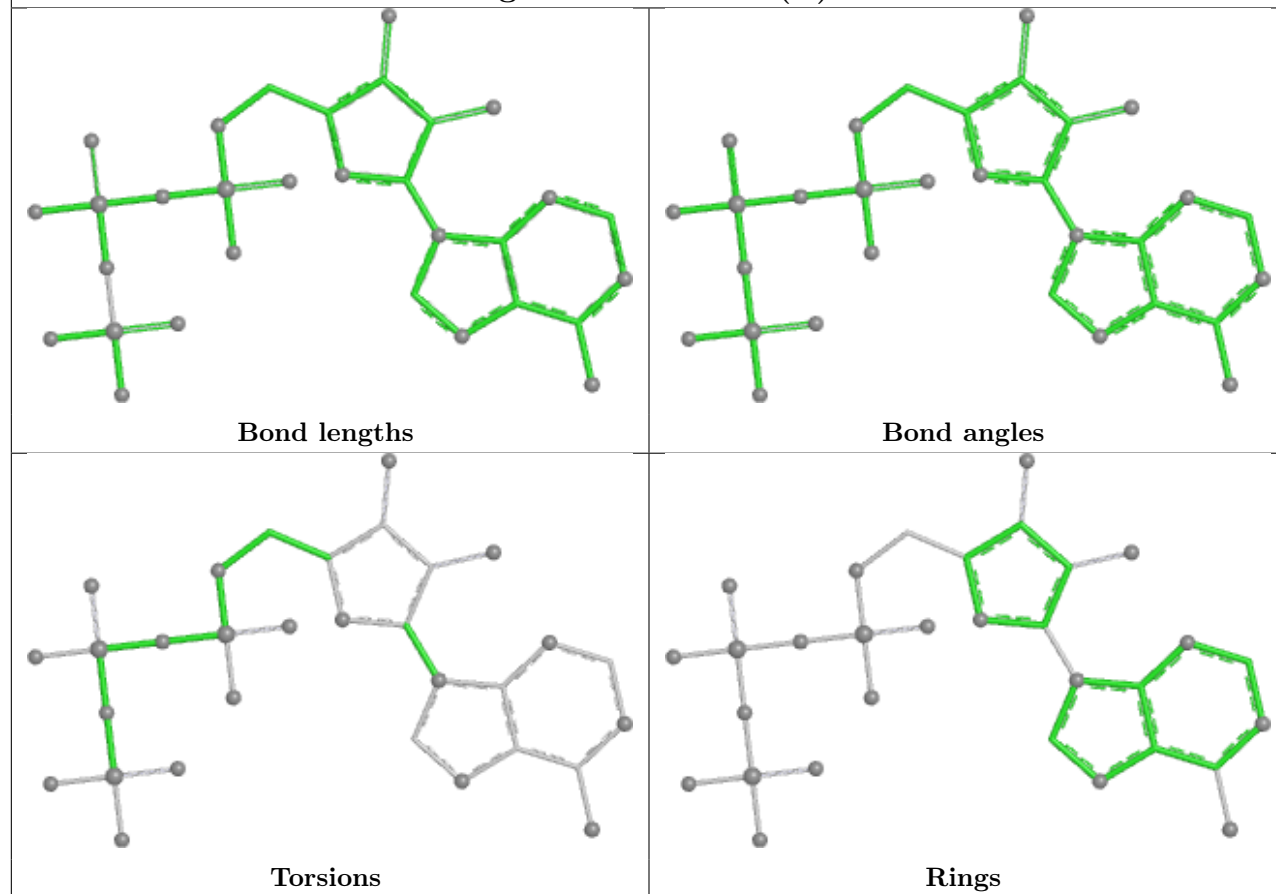
Ligand ATP E 502 (B)



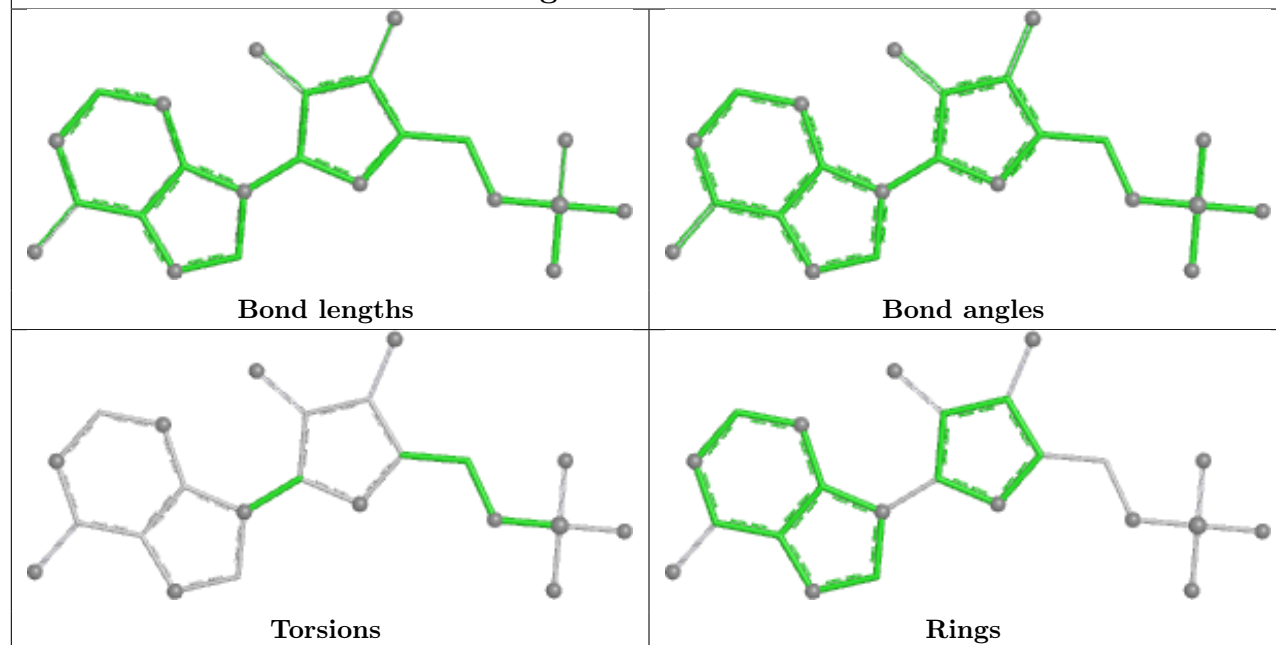
Ligand IMP J 501

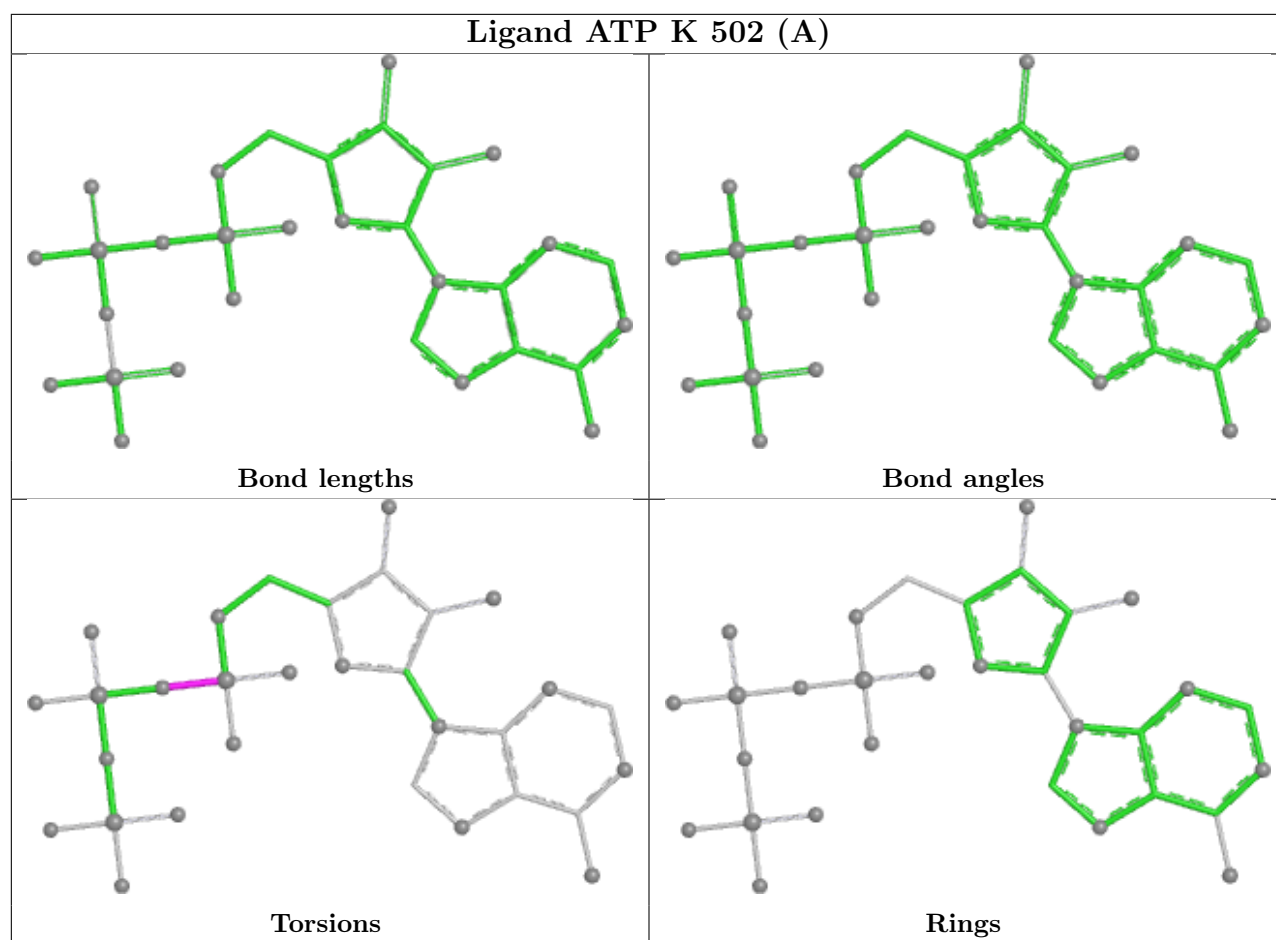


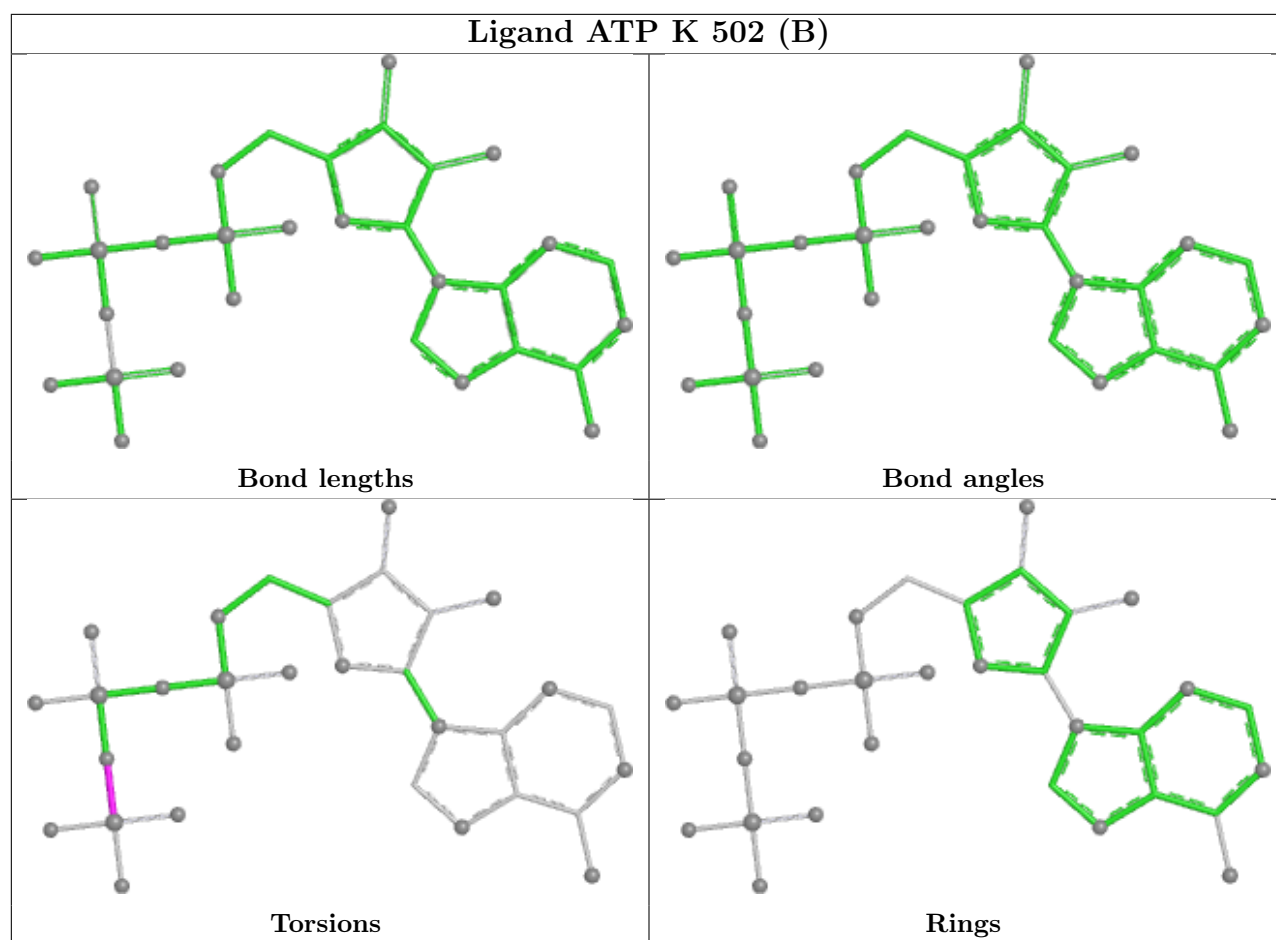
Ligand ATP H 502 (B)

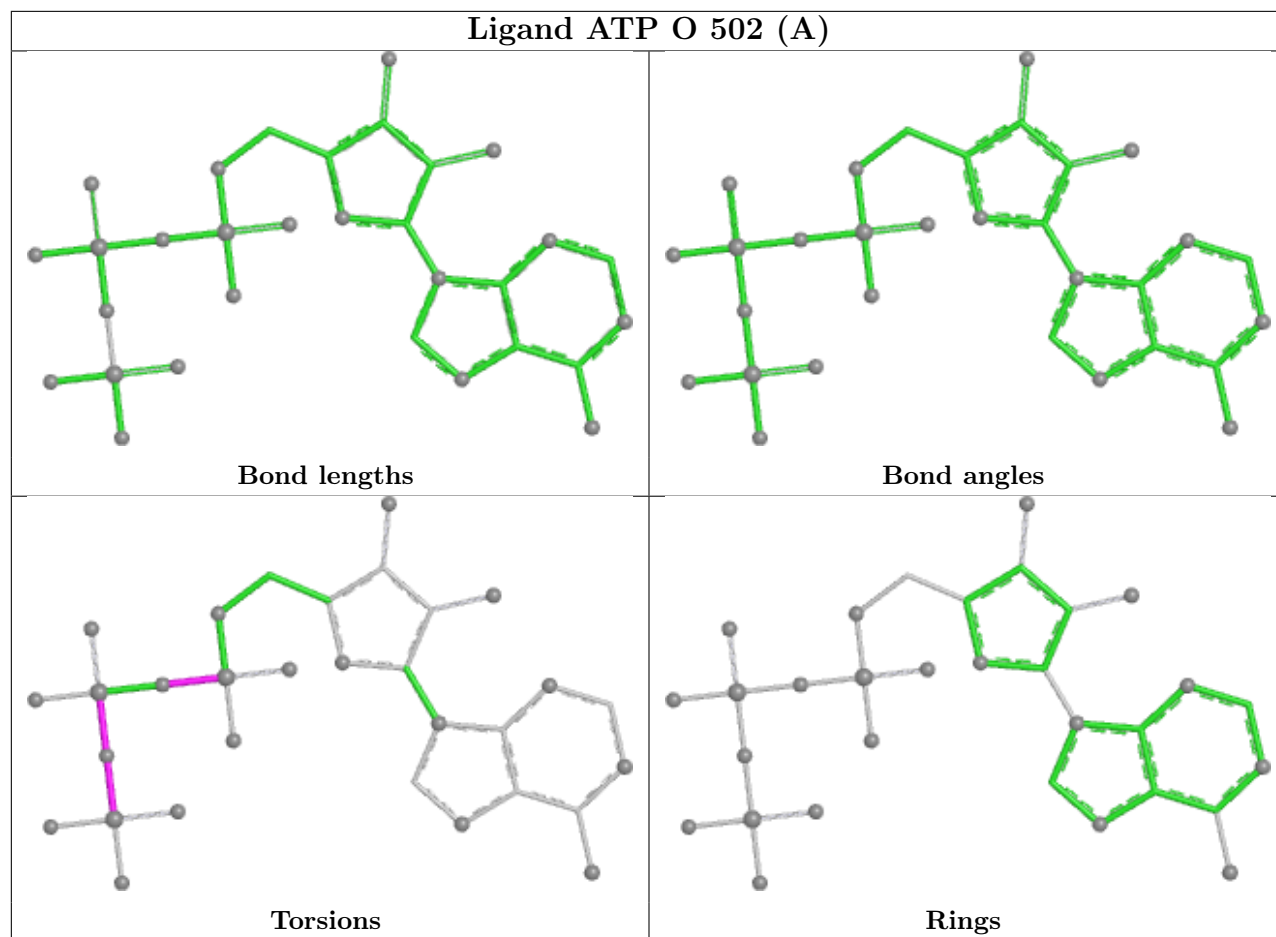


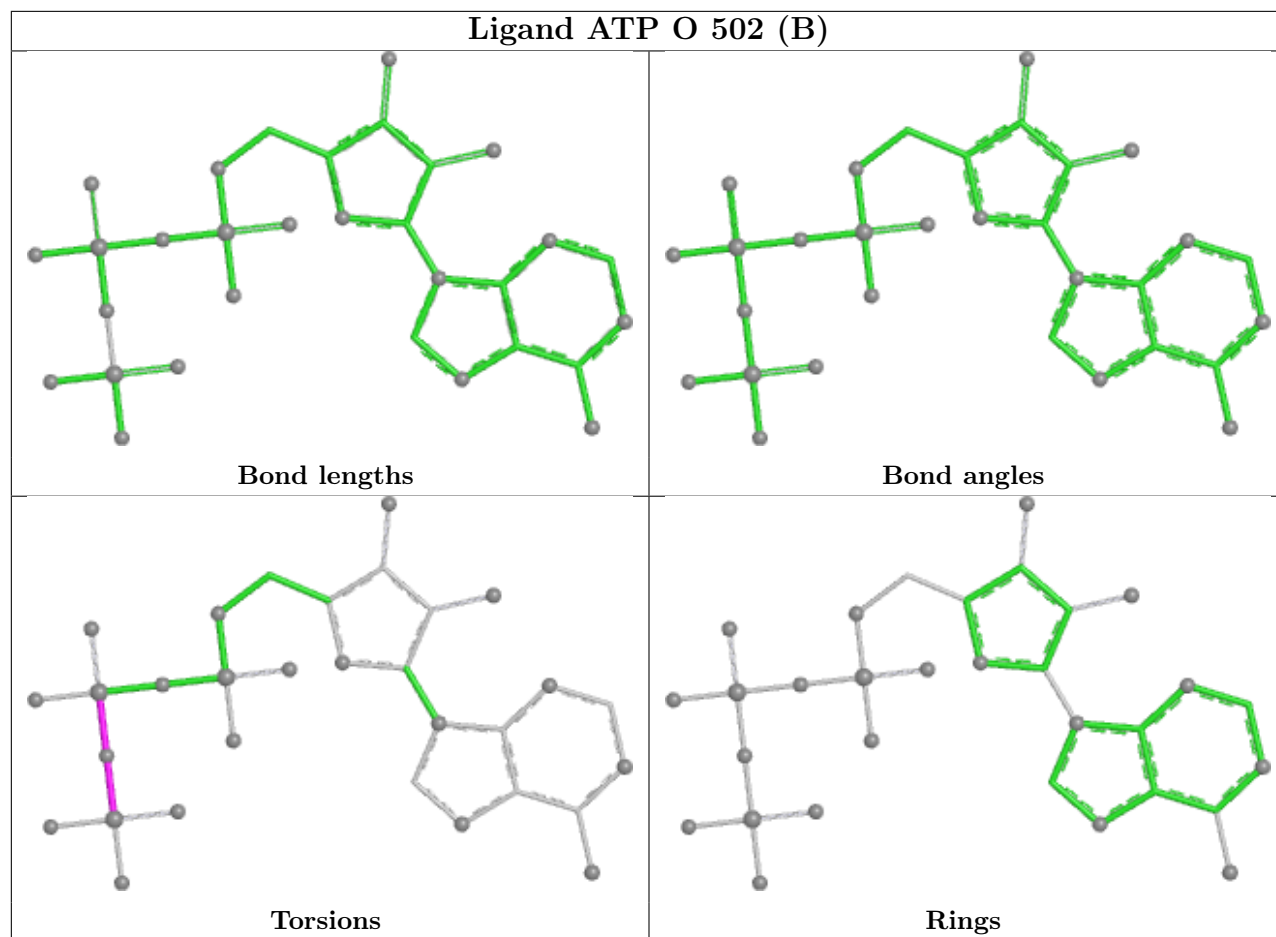
Ligand IMP I 501

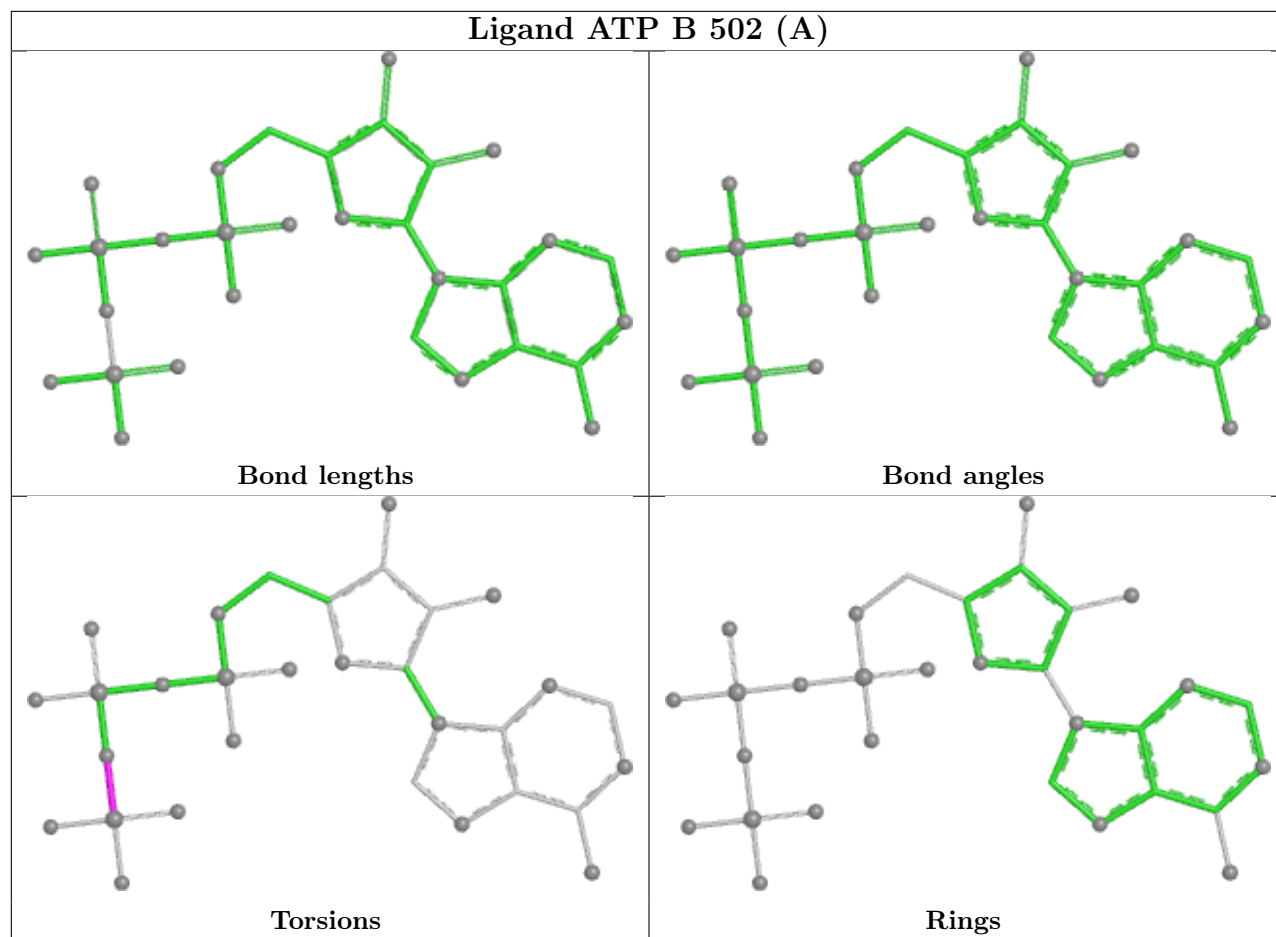




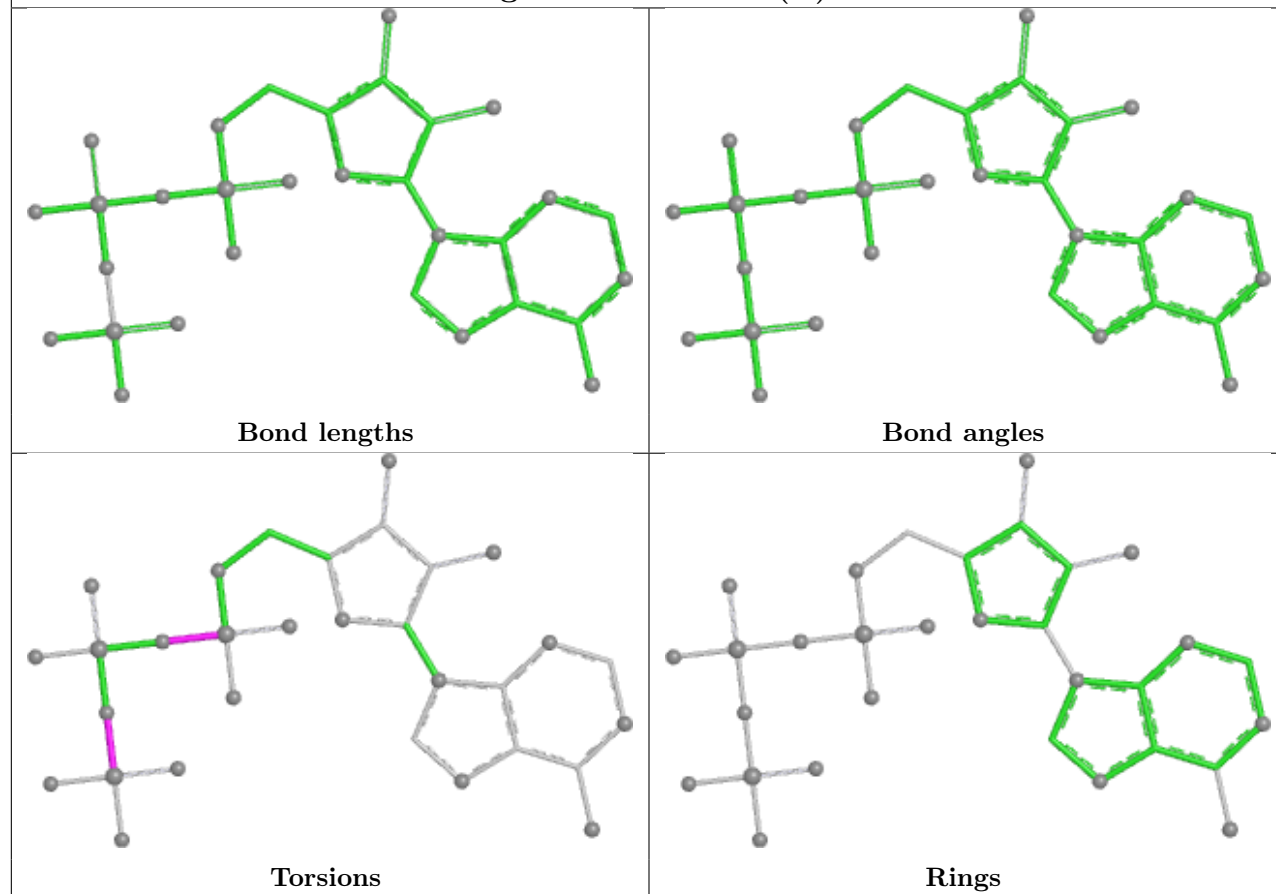




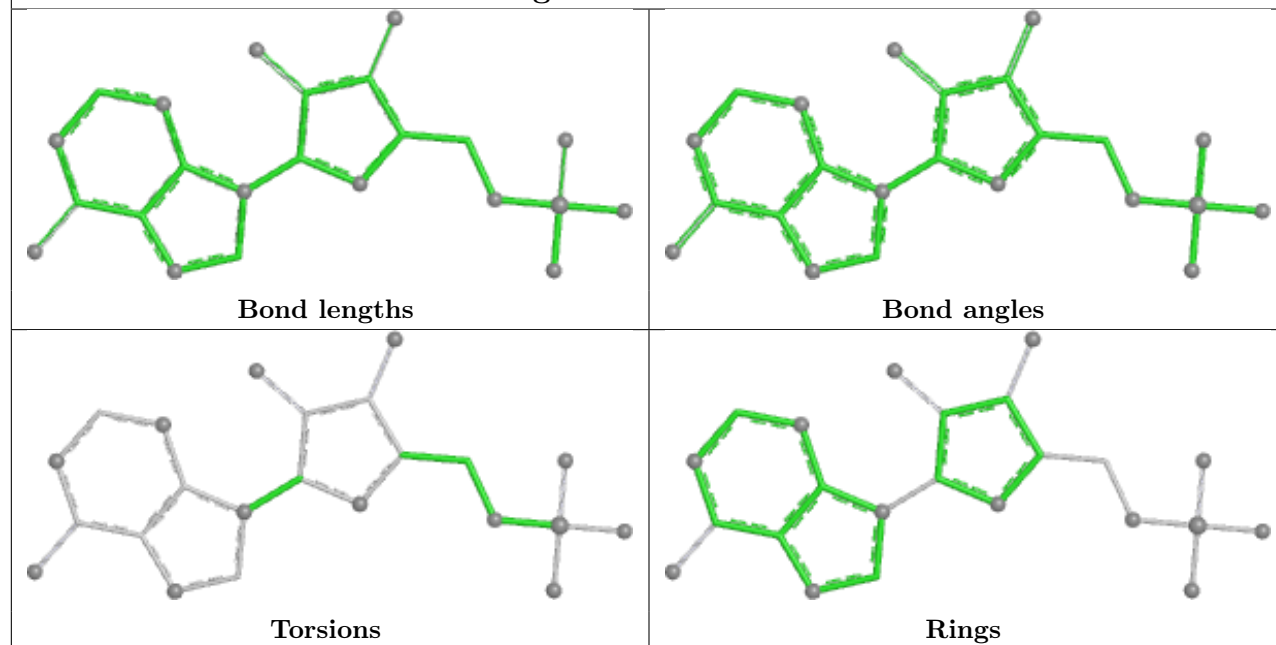


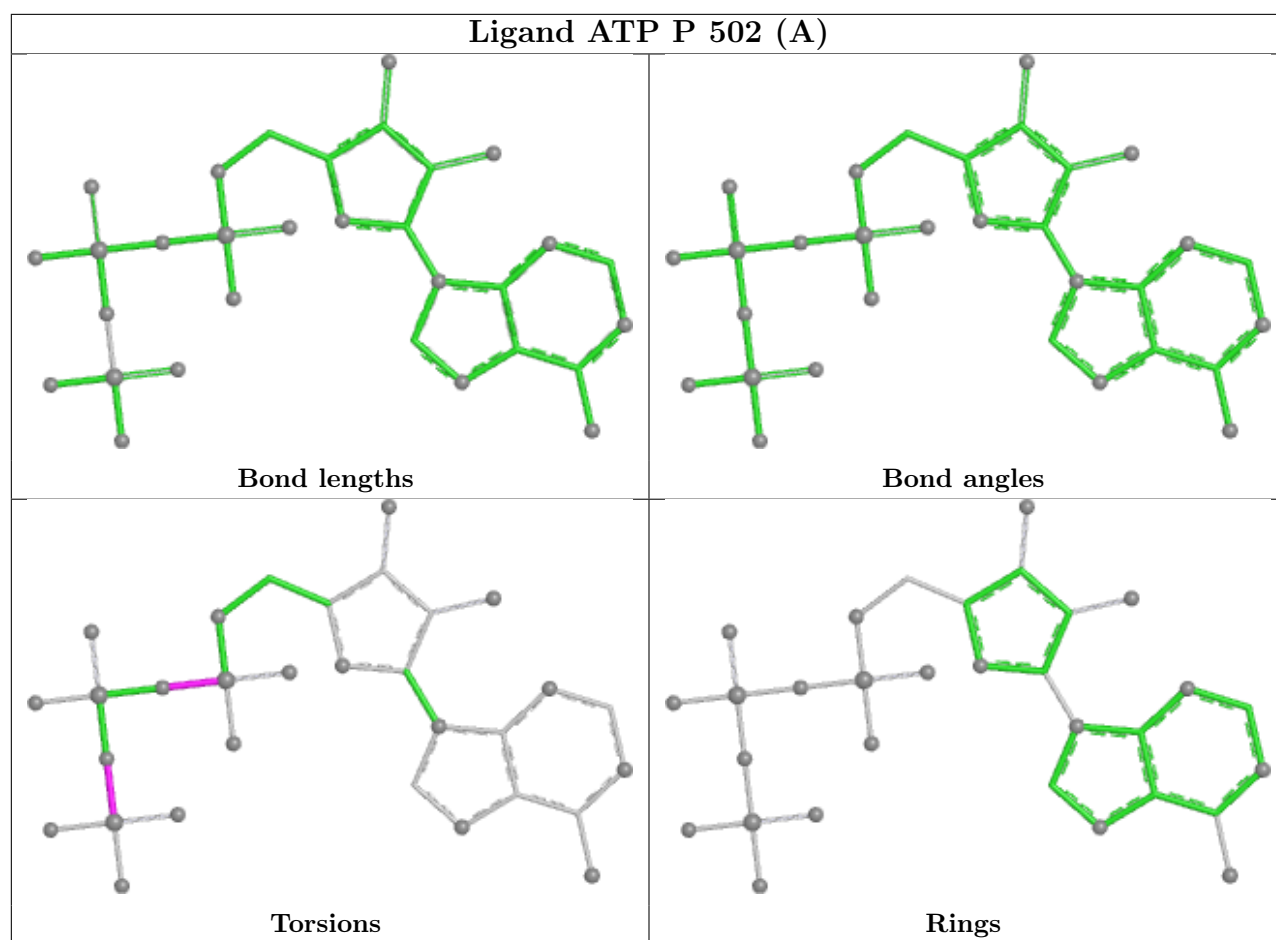


Ligand ATP B 502 (B)

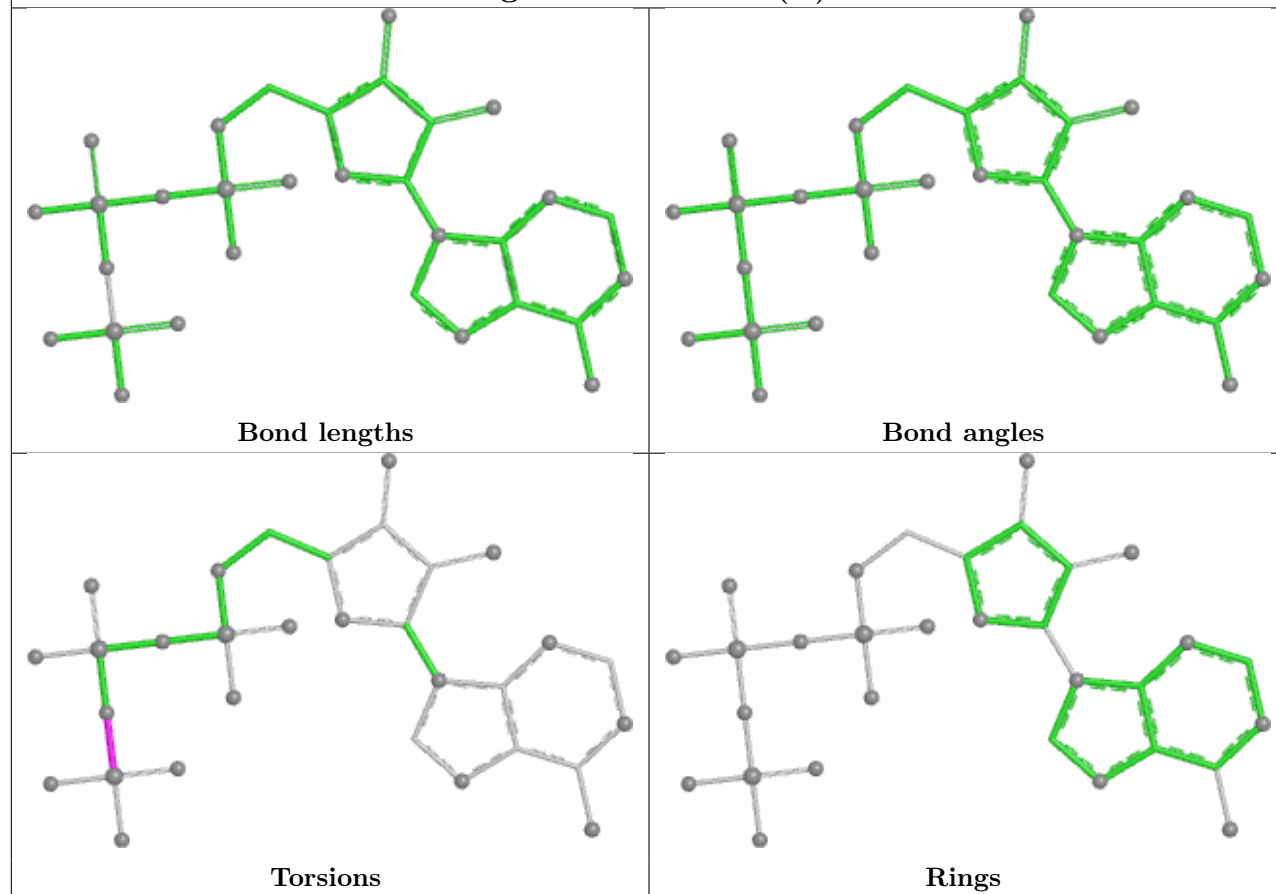


Ligand IMP O 501

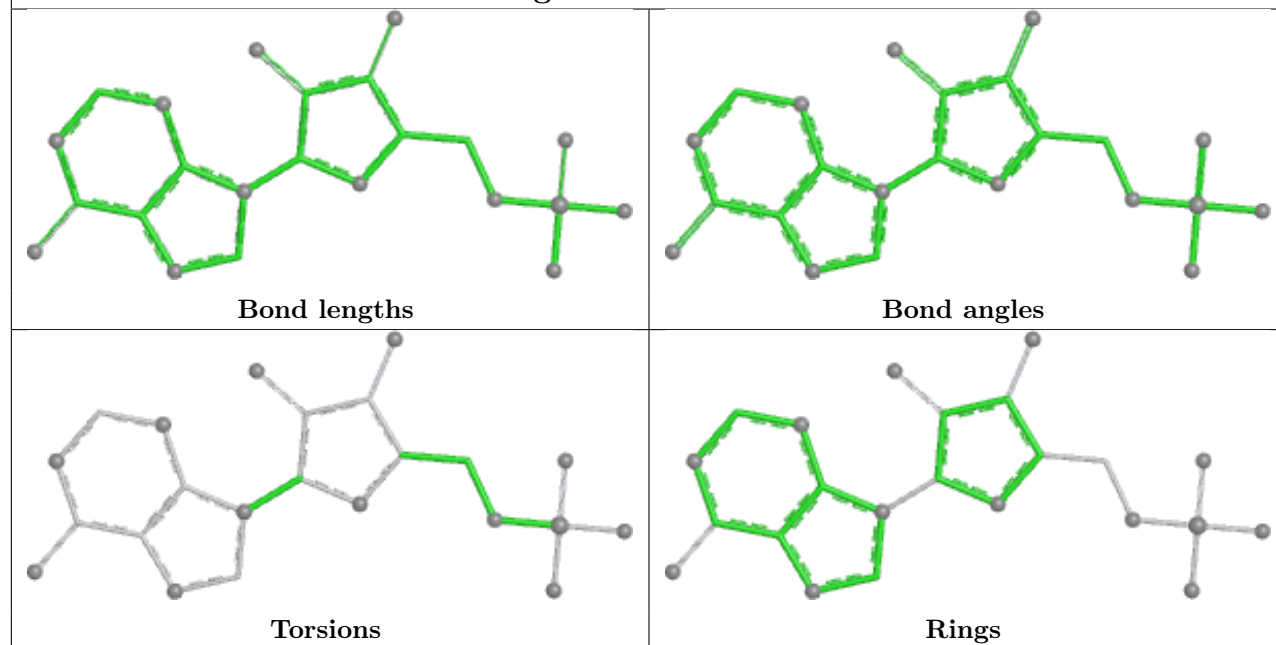




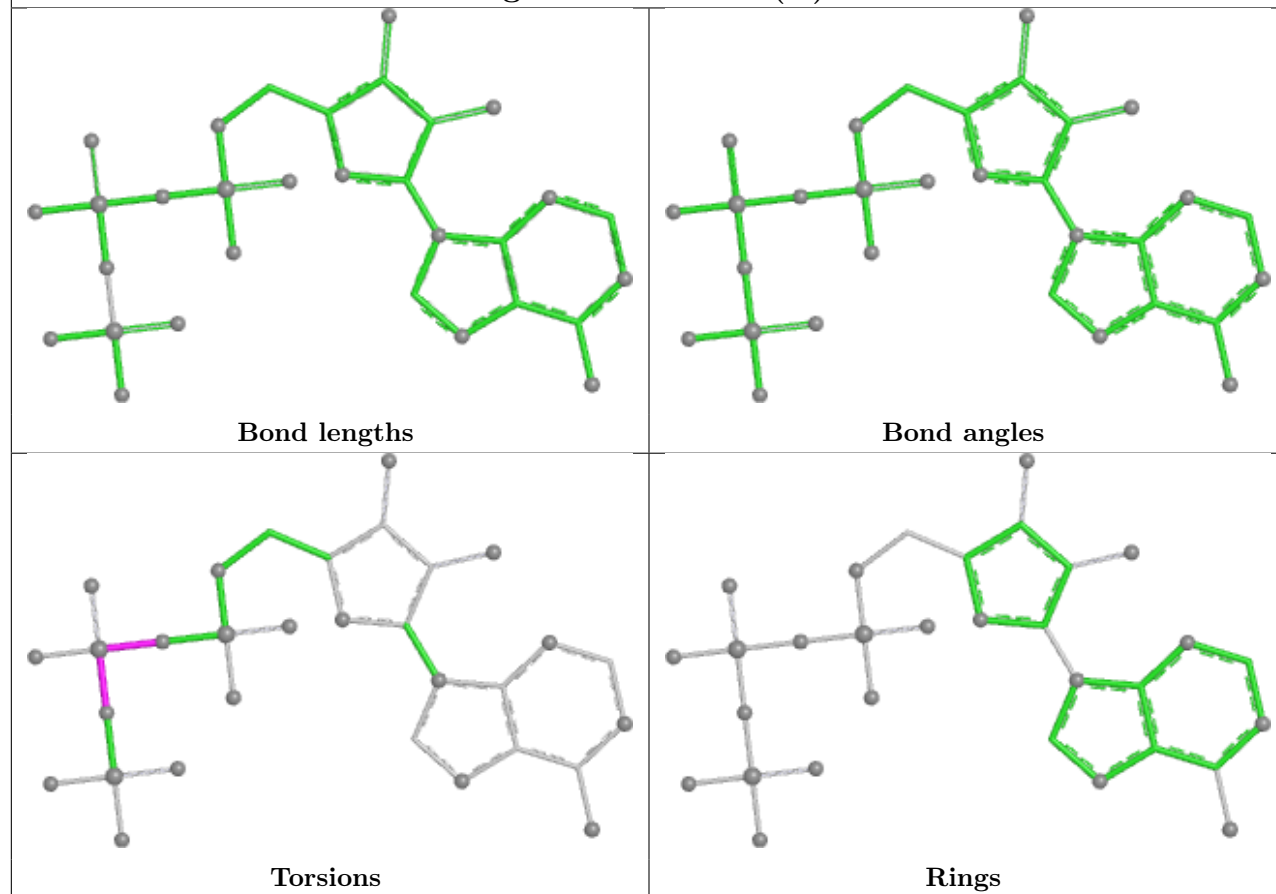
Ligand ATP P 502 (B)



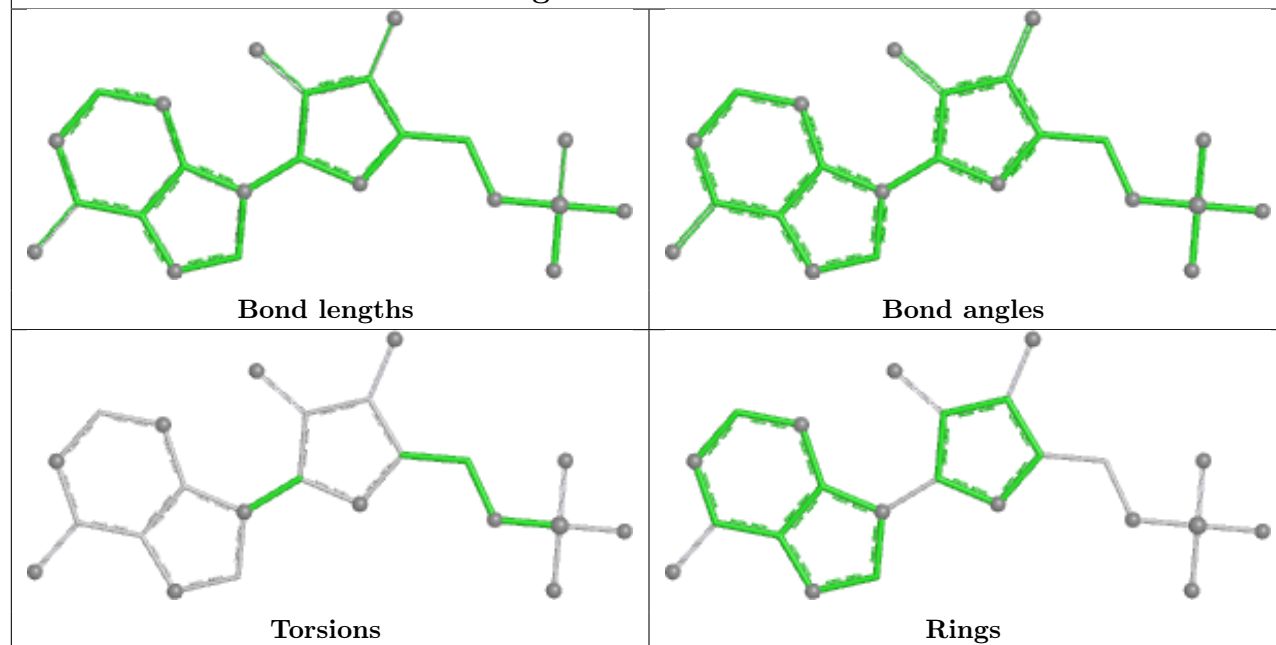
Ligand IMP B 501



Ligand ATP J 502 (A)



Ligand IMP K 501



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	461/496 (92%)	0.33	21 (4%)	37	39	13, 28, 45, 63	4 (0%)
1	B	460/496 (92%)	0.33	24 (5%)	33	33	14, 29, 50, 67	3 (0%)
1	C	461/496 (92%)	0.29	27 (5%)	28	29	13, 26, 46, 66	4 (0%)
1	D	460/496 (92%)	0.12	21 (4%)	37	39	12, 25, 43, 59	3 (0%)
1	E	460/496 (92%)	0.16	19 (4%)	41	43	13, 25, 45, 62	4 (0%)
1	F	461/496 (92%)	0.21	25 (5%)	31	32	12, 25, 45, 64	3 (0%)
1	G	461/496 (92%)	0.16	23 (4%)	34	35	12, 25, 44, 63	3 (0%)
1	H	461/496 (92%)	0.14	18 (3%)	43	45	13, 25, 43, 64	3 (0%)
1	I	461/496 (92%)	0.81	45 (9%)	13	13	17, 36, 57, 74	3 (0%)
1	J	461/496 (92%)	0.28	29 (6%)	26	26	14, 27, 49, 66	3 (0%)
1	K	460/496 (92%)	0.39	28 (6%)	27	27	14, 28, 49, 66	3 (0%)
1	L	461/496 (92%)	0.70	39 (8%)	16	16	16, 33, 53, 67	3 (0%)
1	M	461/496 (92%)	0.88	37 (8%)	18	18	18, 38, 56, 73	3 (0%)
1	N	460/496 (92%)	0.67	29 (6%)	26	26	17, 34, 52, 71	3 (0%)
1	O	461/496 (92%)	0.83	32 (6%)	23	23	17, 36, 54, 72	3 (0%)
1	P	460/496 (92%)	0.87	39 (8%)	16	16	20, 38, 56, 73	3 (0%)
All	All	7370/7936 (92%)	0.45	456 (6%)	26	27	12, 30, 51, 74	51 (0%)

All (456) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	425	ALA	6.4
1	H	29	ALA	5.6
1	N	425	ALA	5.4
1	L	424	PRO	5.2
1	J	467	ALA	5.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	425	ALA	4.8
1	J	128	PHE	4.8
1	E	29	ALA	4.7
1	E	425	ALA	4.7
1	J	425	ALA	4.7
1	D	29	ALA	4.7
1	G	425	ALA	4.7
1	M	467	ALA	4.7
1	H	467	ALA	4.6
1	P	29	ALA	4.6
1	K	424	PRO	4.5
1	P	425	ALA	4.5
1	B	424	PRO	4.5
1	A	467	ALA	4.5
1	C	29	ALA	4.5
1	N	466	ALA	4.5
1	B	425	ALA	4.4
1	B	467	ALA	4.4
1	C	467	ALA	4.4
1	L	394	ALA	4.4
1	I	128	PHE	4.4
1	G	376	ARG	4.4
1	L	374	PHE	4.3
1	I	467	ALA	4.3
1	O	467	ALA	4.3
1	M	466	ALA	4.3
1	O	425	ALA	4.3
1	C	424	PRO	4.3
1	K	466	ALA	4.2
1	K	467	ALA	4.2
1	N	29	ALA	4.2
1	G	29	ALA	4.2
1	I	425	ALA	4.1
1	D	424	PRO	4.1
1	F	424	PRO	4.1
1	L	467	ALA	4.1
1	M	424	PRO	4.1
1	F	29	ALA	4.0
1	F	466	ALA	4.0
1	K	425	ALA	4.0
1	G	424	PRO	4.0
1	G	467	ALA	4.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	424	PRO	3.9
1	F	425	ALA	3.9
1	D	467	ALA	3.9
1	O	300	GLY	3.8
1	P	128	PHE	3.8
1	E	467	ALA	3.8
1	C	425	ALA	3.8
1	C	410	LEU	3.7
1	C	403	PHE	3.7
1	L	2	VAL	3.7
1	L	425	ALA	3.7
1	C	401	SER	3.7
1	J	376	ARG	3.7
1	O	424	PRO	3.7
1	N	467	ALA	3.7
1	O	394	ALA	3.7
1	M	29	ALA	3.6
1	M	403	PHE	3.6
1	G	378	ASP	3.6
1	C	466	ALA	3.6
1	F	467	ALA	3.6
1	M	425	ALA	3.6
1	H	424	PRO	3.6
1	I	424	PRO	3.6
1	O	128	PHE	3.6
1	O	410	LEU	3.6
1	N	401	SER	3.5
1	N	424	PRO	3.5
1	I	394	ALA	3.5
1	E	426	ARG	3.5
1	G	377	ASP	3.5
1	I	186	ALA	3.4
1	P	142	ALA	3.4
1	G	401	SER	3.4
1	H	466	ALA	3.4
1	M	394	ALA	3.4
1	C	1	MET	3.4
1	B	394	ALA	3.4
1	I	29	ALA	3.4
1	P	394	ALA	3.4
1	P	467	ALA	3.4
1	M	376	ARG	3.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	I	403	PHE	3.4
1	D	466	ALA	3.3
1	I	466	ALA	3.3
1	J	394	ALA	3.3
1	D	410	LEU	3.3
1	F	128	PHE	3.3
1	M	411	PHE	3.3
1	J	2	VAL	3.3
1	P	144	VAL	3.3
1	A	1	MET	3.3
1	I	302	MET	3.3
1	B	29	ALA	3.3
1	O	29	ALA	3.3
1	D	426	ARG	3.3
1	G	302	MET	3.3
1	J	424	PRO	3.3
1	P	402	SER	3.3
1	F	374	PHE	3.3
1	A	417	THR	3.2
1	G	416	SER	3.2
1	K	377	ASP	3.2
1	P	378	ASP	3.2
1	I	410	LEU	3.2
1	N	410	LEU	3.2
1	N	411	PHE	3.2
1	P	403	PHE	3.2
1	B	144	VAL	3.2
1	N	404	ASP	3.2
1	E	466	ALA	3.2
1	A	376	ARG	3.2
1	P	376	ARG	3.2
1	J	378	ASP	3.2
1	M	402	SER	3.2
1	I	426	ARG	3.1
1	M	426	ARG	3.1
1	M	329	LEU	3.1
1	J	29	ALA	3.1
1	B	403	PHE	3.1
1	A	424	PRO	3.1
1	P	401	SER	3.1
1	H	1	MET	3.1
1	O	374	PHE	3.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	L	376	ARG	3.1
1	P	424	PRO	3.1
1	G	394	ALA	3.1
1	K	394	ALA	3.1
1	P	374	PHE	3.0
1	B	376	ARG	3.0
1	C	376	ARG	3.0
1	K	2	VAL	3.0
1	B	466	ALA	3.0
1	I	201	ILE	3.0
1	P	155	LEU	3.0
1	L	411	PHE	3.0
1	C	378	ASP	2.9
1	J	466	ALA	2.9
1	O	466	ALA	2.9
1	I	155	LEU	2.9
1	I	416	SER	2.9
1	P	416	SER	2.9
1	C	408	LYS	2.9
1	D	2	VAL	2.9
1	D	394	ALA	2.9
1	I	187	PRO	2.9
1	D	302	MET	2.9
1	P	302	MET	2.9
1	E	401	SER	2.9
1	C	128	PHE	2.9
1	J	411	PHE	2.9
1	L	29	ALA	2.9
1	K	410	LEU	2.9
1	O	271	LEU	2.9
1	I	153	ILE	2.9
1	D	401	SER	2.9
1	N	426	ARG	2.8
1	A	378	ASP	2.8
1	M	378	ASP	2.8
1	I	271	LEU	2.8
1	E	302	MET	2.8
1	O	378	ASP	2.8
1	B	143	GLY	2.8
1	K	413	GLU	2.8
1	I	417	THR	2.8
1	O	1	MET	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	128	PHE	2.8
1	D	402	SER	2.8
1	K	376	ARG	2.8
1	O	376	ARG	2.8
1	K	302	MET	2.8
1	I	378	ASP	2.8
1	P	466	ALA	2.7
1	L	1	MET	2.7
1	E	404	ASP	2.7
1	G	404	ASP	2.7
1	N	377	ASP	2.7
1	I	185	THR	2.7
1	C	405	ARG	2.7
1	I	142	ALA	2.7
1	E	2	VAL	2.7
1	J	416	SER	2.7
1	N	416	SER	2.7
1	L	409	GLY	2.7
1	C	302	MET	2.7
1	M	408	LYS	2.7
1	L	466	ALA	2.7
1	O	377	ASP	2.7
1	J	379	ARG	2.7
1	M	374	PHE	2.7
1	N	403	PHE	2.7
1	E	402	SER	2.7
1	M	409	GLY	2.7
1	G	1	MET	2.7
1	P	27	ASP	2.7
1	A	29	ALA	2.7
1	F	114	ALA	2.7
1	K	416	SER	2.6
1	G	403	PHE	2.6
1	M	373	LEU	2.6
1	B	378	ASP	2.6
1	A	466	ALA	2.6
1	N	394	ALA	2.6
1	P	114	ALA	2.6
1	D	409	GLY	2.6
1	H	374	PHE	2.6
1	I	189	GLY	2.6
1	I	300	GLY	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	I	411	PHE	2.6
1	L	403	PHE	2.6
1	N	128	PHE	2.6
1	P	411	PHE	2.6
1	N	376	ARG	2.6
1	B	404	ASP	2.6
1	P	271	LEU	2.6
1	A	402	SER	2.6
1	H	425	ALA	2.6
1	N	402	SER	2.6
1	E	376	ARG	2.6
1	F	143	GLY	2.6
1	E	378	ASP	2.6
1	E	128	PHE	2.6
1	K	374	PHE	2.6
1	L	116	LEU	2.6
1	O	56	VAL	2.6
1	H	401	SER	2.6
1	I	133	ILE	2.6
1	K	153	ILE	2.6
1	E	394	ALA	2.6
1	F	394	ALA	2.6
1	J	142	ALA	2.6
1	H	376	ARG	2.6
1	J	426	ARG	2.6
1	C	414	GLY	2.5
1	I	81	THR	2.5
1	H	128	PHE	2.5
1	I	374	PHE	2.5
1	J	374	PHE	2.5
1	K	403	PHE	2.5
1	F	402	SER	2.5
1	L	410	LEU	2.5
1	N	2	VAL	2.5
1	C	394	ALA	2.5
1	O	114	ALA	2.5
1	O	301	ALA	2.5
1	P	408	LYS	2.5
1	B	156	SER	2.5
1	J	1	MET	2.5
1	K	401	SER	2.5
1	I	376	ARG	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	E	374	PHE	2.5
1	G	374	PHE	2.5
1	K	128	PHE	2.5
1	M	128	PHE	2.5
1	B	114	ALA	2.5
1	L	133	ILE	2.5
1	B	401	SER	2.5
1	F	155	LEU	2.5
1	H	377	ASP	2.5
1	L	128	PHE	2.5
1	M	410	LEU	2.5
1	I	301	ALA	2.5
1	K	114	ALA	2.5
1	L	187	PRO	2.5
1	O	299	PRO	2.5
1	P	187	PRO	2.5
1	I	144	VAL	2.5
1	O	408	LYS	2.5
1	I	413	GLU	2.5
1	J	377	ASP	2.4
1	L	377	ASP	2.4
1	K	408	LYS	2.4
1	F	302	MET	2.4
1	I	2	VAL	2.4
1	L	163	VAL	2.4
1	L	417	THR	2.4
1	C	404	ASP	2.4
1	I	116	LEU	2.4
1	L	414	GLY	2.4
1	N	414	GLY	2.4
1	H	394	ALA	2.4
1	H	402	SER	2.4
1	M	401	SER	2.4
1	P	212	ALA	2.4
1	F	413	GLU	2.4
1	I	1	MET	2.4
1	O	302	MET	2.4
1	B	402	SER	2.4
1	L	115	LEU	2.4
1	M	115	LEU	2.4
1	F	403	PHE	2.4
1	N	374	PHE	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	K	404	ASP	2.4
1	A	426	ARG	2.4
1	O	144	VAL	2.4
1	J	413	GLU	2.3
1	A	139	ALA	2.3
1	A	377	ASP	2.3
1	B	405	ARG	2.3
1	G	27	ASP	2.3
1	G	426	ARG	2.3
1	G	466	ALA	2.3
1	L	3	ARG	2.3
1	M	422	LEU	2.3
1	N	157	ASP	2.3
1	P	116	LEU	2.3
1	N	408	LYS	2.3
1	C	411	PHE	2.3
1	I	147	PHE	2.3
1	O	32	PHE	2.3
1	L	117	HIS	2.3
1	K	133	ILE	2.3
1	I	150	VAL	2.3
1	O	2	VAL	2.3
1	E	413	GLU	2.3
1	H	417	THR	2.3
1	O	417	THR	2.3
1	M	155	LEU	2.3
1	M	173	LEU	2.3
1	M	117	HIS	2.3
1	A	328	GLN	2.3
1	L	32	PHE	2.3
1	P	179	ILE	2.3
1	O	402	SER	2.3
1	F	2	VAL	2.3
1	H	426	ARG	2.3
1	P	2	VAL	2.3
1	J	143	GLY	2.3
1	O	130	GLY	2.3
1	I	192	ALA	2.3
1	M	142	ALA	2.3
1	G	410	LEU	2.3
1	L	173	LEU	2.3
1	A	403	PHE	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	403	PHE	2.3
1	G	411	PHE	2.3
1	L	416	SER	2.3
1	B	423	ASP	2.3
1	J	409	GLY	2.3
1	M	163	VAL	2.3
1	P	380	PRO	2.2
1	C	412	GLU	2.2
1	K	29	ALA	2.2
1	L	406	ALA	2.2
1	D	113	ASN	2.2
1	F	426	ARG	2.2
1	J	405	ARG	2.2
1	P	410	LEU	2.2
1	M	418	SER	2.2
1	M	377	ASP	2.2
1	G	128	PHE	2.2
1	N	415	ILE	2.2
1	D	412	GLU	2.2
1	K	28	VAL	2.2
1	F	113	ASN	2.2
1	J	401	SER	2.2
1	I	173	LEU	2.2
1	J	155	LEU	2.2
1	L	302	MET	2.2
1	N	302	MET	2.2
1	D	413	GLU	2.2
1	D	411	PHE	2.2
1	J	147	PHE	2.2
1	J	403	PHE	2.2
1	O	403	PHE	2.2
1	L	153	ILE	2.2
1	O	133	ILE	2.2
1	P	3	ARG	2.2
1	J	417	THR	2.2
1	L	408	LYS	2.2
1	P	109	VAL	2.2
1	A	394	ALA	2.2
1	D	378	ASP	2.2
1	F	377	ASP	2.2
1	F	1	MET	2.2
1	K	139	ALA	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	P	129	GLU	2.2
1	C	409	GLY	2.2
1	I	414	GLY	2.2
1	L	155	LEU	2.2
1	C	374	PHE	2.2
1	G	32	PHE	2.2
1	B	375	ASP	2.1
1	I	377	ASP	2.1
1	M	2	VAL	2.1
1	N	144	VAL	2.1
1	A	413	GLU	2.1
1	C	413	GLU	2.1
1	I	212	ALA	2.1
1	C	215	ARG	2.1
1	C	426	ARG	2.1
1	L	379	ARG	2.1
1	C	143	GLY	2.1
1	B	302	MET	2.1
1	E	377	ASP	2.1
1	J	302	MET	2.1
1	M	302	MET	2.1
1	F	411	PHE	2.1
1	H	378	ASP	2.1
1	P	39	VAL	2.1
1	D	300	GLY	2.1
1	N	113	ASN	2.1
1	B	377	ASP	2.1
1	I	401	SER	2.1
1	L	412	GLU	2.1
1	P	377	ASP	2.1
1	K	417	THR	2.1
1	O	327	ARG	2.1
1	B	374	PHE	2.1
1	J	140	ASN	2.1
1	P	156	SER	2.1
1	P	265	LEU	2.1
1	I	32	PHE	2.1
1	M	201	ILE	2.1
1	F	301	ALA	2.0
1	M	114	ALA	2.0
1	O	139	ALA	2.0
1	E	423	ASP	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	K	378	ASP	2.0
1	A	405	ARG	2.0
1	F	144	VAL	2.0
1	F	376	ARG	2.0
1	G	402	SER	2.0
1	H	418	SER	2.0
1	L	144	VAL	2.0
1	L	150	VAL	2.0
1	M	156	SER	2.0
1	N	262	VAL	2.0
1	O	27	ASP	2.0
1	P	28	VAL	2.0
1	A	141	CYS	2.0
1	K	303	CYS	2.0
1	K	116	LEU	2.0
1	M	372	LEU	2.0
1	N	117	HIS	2.0
1	P	190	THR	2.0
1	N	413	GLU	2.0
1	H	130	GLY	2.0
1	D	374	PHE	2.0
1	I	131	ARG	2.0
1	B	142	ALA	2.0
1	F	378	ASP	2.0
1	L	114	ALA	2.0
1	L	378	ASP	2.0
1	M	301	ALA	2.0
1	C	402	SER	2.0
1	A	144	VAL	2.0
1	B	2	VAL	2.0
1	M	39	VAL	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 oligosaccharides in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ATP	M	502[A]	31/31	0.85	0.12	39,41,49,50	31
3	ATP	M	502[B]	31/31	0.85	0.12	38,41,49,51	31
3	ATP	O	502[A]	31/31	0.87	0.11	31,33,46,48	31
3	ATP	O	502[B]	31/31	0.87	0.11	31,34,46,47	31
3	ATP	N	502[A]	31/31	0.88	0.11	32,34,43,47	31
3	ATP	N	502[B]	31/31	0.88	0.11	32,34,42,44	31
3	ATP	L	502[A]	31/31	0.88	0.11	34,37,46,49	31
3	ATP	L	502[B]	31/31	0.88	0.11	35,37,46,49	31
3	ATP	I	502[A]	31/31	0.89	0.11	35,40,49,50	31
3	ATP	I	502[B]	31/31	0.89	0.11	35,40,49,51	31
3	ATP	J	502[A]	31/31	0.89	0.10	28,32,39,39	31
3	ATP	J	502[B]	31/31	0.89	0.10	28,32,38,38	31
3	ATP	P	502[A]	31/31	0.89	0.12	36,39,48,50	31
3	ATP	P	502[B]	31/31	0.89	0.12	38,40,48,51	31
3	ATP	C	502[A]	31/31	0.92	0.09	26,28,35,36	31
3	ATP	C	502[B]	31/31	0.92	0.09	24,28,35,36	31
3	ATP	G	502[A]	31/31	0.92	0.09	21,27,38,39	31
3	ATP	G	502[B]	31/31	0.92	0.09	23,27,38,39	31
3	ATP	A	502[A]	31/31	0.92	0.09	26,29,39,41	31
3	ATP	A	502[B]	31/31	0.92	0.09	24,29,41,41	31
3	ATP	B	502[A]	31/31	0.92	0.10	26,29,38,40	31
3	ATP	B	502[B]	31/31	0.92	0.10	23,29,38,40	31
3	ATP	K	502[A]	31/31	0.92	0.09	24,29,39,39	31
3	ATP	K	502[B]	31/31	0.92	0.09	27,30,39,39	31
3	ATP	D	502[A]	31/31	0.93	0.08	20,25,36,38	31
3	ATP	D	502[B]	31/31	0.93	0.08	22,25,37,38	31
3	ATP	H	502[A]	31/31	0.93	0.09	19,28,35,37	31
3	ATP	H	502[B]	31/31	0.93	0.09	24,28,35,37	31
3	ATP	E	502[A]	31/31	0.94	0.08	17,25,32,36	31
3	ATP	E	502[B]	31/31	0.94	0.08	23,25,32,35	31
3	ATP	F	502[A]	31/31	0.94	0.09	17,26,38,39	31
3	ATP	F	502[B]	31/31	0.94	0.09	24,27,37,38	31
2	IMP	N	501	23/23	0.97	0.07	26,29,32,32	0
2	IMP	O	501	23/23	0.97	0.07	26,31,34,38	0
2	IMP	P	501	23/23	0.97	0.06	27,32,37,38	0
2	IMP	I	501	23/23	0.97	0.07	27,31,34,35	0
2	IMP	L	501	23/23	0.97	0.06	23,27,30,33	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	IMP	M	501	23/23	0.98	0.06	29,33,38,41	0
2	IMP	C	501	23/23	0.98	0.05	19,22,27,29	0
2	IMP	F	501	23/23	0.98	0.05	17,21,23,27	0
2	IMP	G	501	23/23	0.98	0.05	19,21,24,27	0
2	IMP	H	501	23/23	0.98	0.05	19,21,24,26	0
2	IMP	A	501	23/23	0.98	0.05	18,23,27,28	0
2	IMP	J	501	23/23	0.98	0.05	21,25,30,32	0
2	IMP	K	501	23/23	0.98	0.05	21,23,26,28	0
2	IMP	B	501	23/23	0.98	0.06	20,23,28,30	0
2	IMP	E	501	23/23	0.99	0.04	18,20,22,26	0
2	IMP	D	501	23/23	0.99	0.04	20,22,26,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.