



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 29, 2020 – 11:56 pm BST

PDB ID : 3WC0  
Title : Crystal structure of *C. albicans* tRNA(His) guanylyltransferase (Thg1) with GTP  
Authors : Nakamura, A.; Nemoto, T.; Sonoda, T.; Yamashita, K.; Tanaka, I.; Yao, M.  
Deposited on : 2013-05-24  
Resolution : 3.03 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

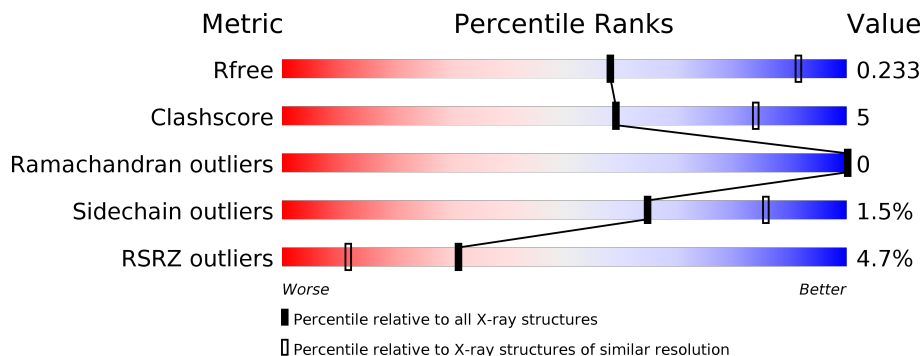
MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins)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

# 1 Overall quality at a glance

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

The reported resolution of this entry is 3.03 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.







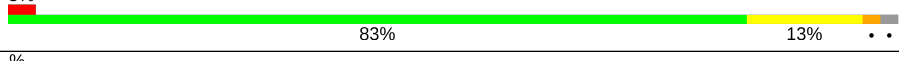
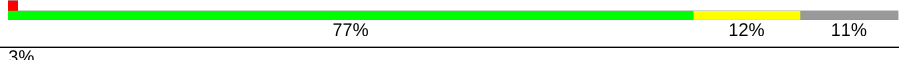

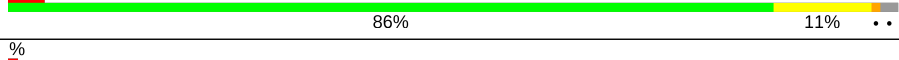


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	130704	2752 (3.08-3.00)
Clashscore	141614	3096 (3.08-3.00)
Ramachandran outliers	138981	2986 (3.08-3.00)
Sidechain outliers	138945	2988 (3.08-3.00)
RSRZ outliers	127900	2636 (3.08-3.00)

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

Mol	Chain	Length	Quality of chain
1	A	271	
1	B	271	
1	C	271	
1	D	271	
1	E	271	
1	F	271	

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Mol	Chain	Length	Quality of chain
1	G	271	
1	H	271	
1	I	271	
1	J	271	
1	K	271	
1	L	271	
1	M	271	
1	N	271	
1	O	271	
1	P	271	

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
3	MG	K	402	-	-	-	X

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 36341 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 Likely histidyl tRNA-specific guanylyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	265	2267	1464	375	416	12	0	0	0
1	B	265	2267	1464	375	416	12	0	0	0
1	C	265	2267	1464	375	416	12	0	0	0
1	D	265	2267	1464	375	416	12	0	0	0
1	E	241	2057	1337	335	373	12	0	0	0
1	F	242	2065	1341	336	376	12	0	0	0
1	G	265	2267	1464	375	416	12	0	0	0
1	H	265	2267	1464	375	416	12	0	0	0
1	I	242	2065	1341	336	376	12	0	0	0
1	J	265	2267	1464	375	416	12	0	0	0
1	K	265	2267	1464	375	416	12	0	0	0
1	L	242	2065	1341	336	376	12	0	0	0
1	M	265	2267	1464	375	416	12	0	0	0
1	N	265	2267	1464	375	416	12	0	0	0
1	O	265	2267	1464	375	416	12	0	0	0
1	P	242	2065	1341	336	376	12	0	0	0

There are 48 discrepancies between the modelled and reference sequences:

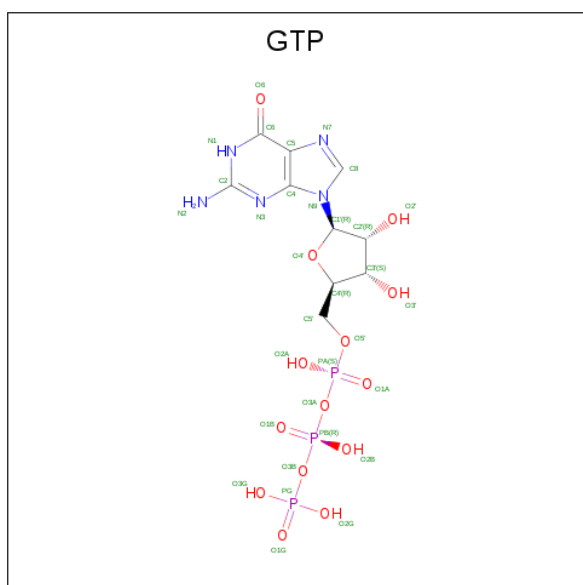
Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
A	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
A	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
B	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
B	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
B	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
C	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
C	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
C	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
D	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
D	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
D	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
E	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
E	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
E	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
F	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
F	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
F	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
G	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
G	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
G	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
H	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
H	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
H	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
I	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
I	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
I	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
J	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
J	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
J	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
K	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
K	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
K	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
L	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
L	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
L	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
M	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
M	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
M	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
N	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
N	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
N	0	SER	-	EXPRESSION TAG	UNP Q5AFK5

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Chain	Residue	Modelled	Actual	Comment	Reference
O	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
O	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
O	0	SER	-	EXPRESSION TAG	UNP Q5AFK5
P	-2	GLY	-	EXPRESSION TAG	UNP Q5AFK5
P	-1	GLY	-	EXPRESSION TAG	UNP Q5AFK5
P	0	SER	-	EXPRESSION TAG	UNP Q5AFK5

- Molecule 2 is GUANOSINE-5'-TRIPHOSPHATE (three-letter code: GTP) (formula:  $C_{10}H_{16}N_5O_{14}P_3$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	N	O			P
2	A	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	A	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	B	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	B	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	C	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	C	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	D	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	D	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	E	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	E	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	F	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	F	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	G	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	G	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	H	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	H	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	I	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	I	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	J	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	J	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	K	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	L	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	L	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	L	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	M	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	M	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	N	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	N	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	O	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	O	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	O	1	Total	C	N	O	P	0	0
			32	10	5	14	3		
2	P	1	Total	C	N	O	P	0	0
			32	10	5	14	3		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	P	3	Total	Mg	0	0
			3	3		
3	G	3	Total	Mg	0	0
			3	3		
3	J	3	Total	Mg	0	0
			3	3		
3	D	3	Total	Mg	0	0
			3	3		
3	K	2	Total	Mg	0	0
			2	2		
3	E	2	Total	Mg	0	0
			2	2		
3	H	3	Total	Mg	0	0
			3	3		
3	B	3	Total	Mg	0	0
			3	3		
3	I	3	Total	Mg	0	0
			3	3		
3	C	3	Total	Mg	0	0
			3	3		
3	A	3	Total	Mg	0	0
			3	3		
3	N	3	Total	Mg	0	0
			3	3		
3	O	3	Total	Mg	0	0
			3	3		
3	L	4	Total	Mg	0	0
			4	4		
3	F	4	Total	Mg	0	0
			4	4		
3	M	3	Total	Mg	0	0
			3	3		



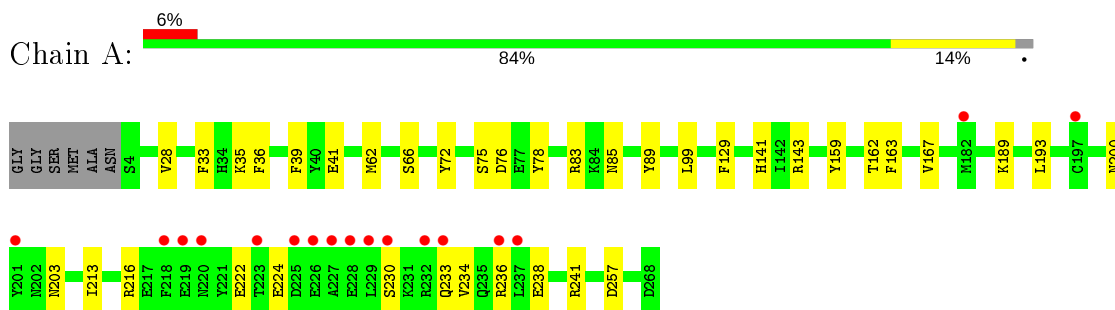
- Molecule 4 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>	<b>ZeroOcc</b>	<b>AltConf</b>
4	C	2	Total O 2 2	0	0
4	G	1	Total O 1 1	0	0
4	H	2	Total O 2 2	0	0
4	I	1	Total O 1 1	0	0
4	K	1	Total O 1 1	0	0
4	M	2	Total O 2 2	0	0
4	N	3	Total O 3 3	0	0
4	O	3	Total O 3 3	0	0

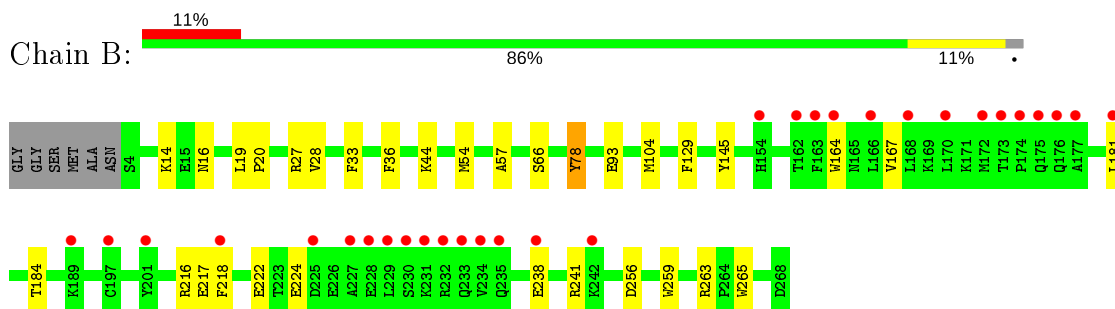
### 3 Residue-property plots [i](#)

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

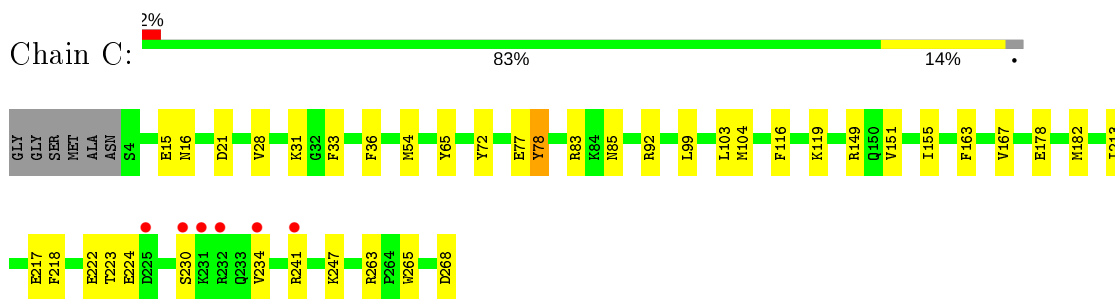
- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase



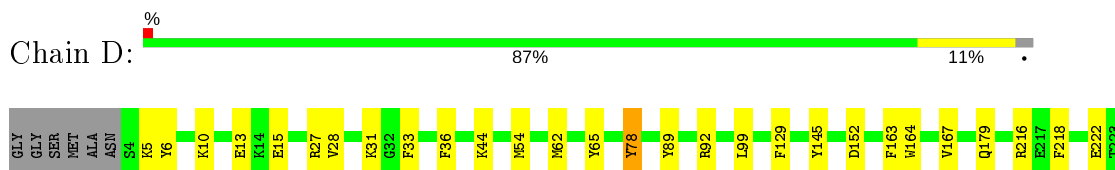
- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase



- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase

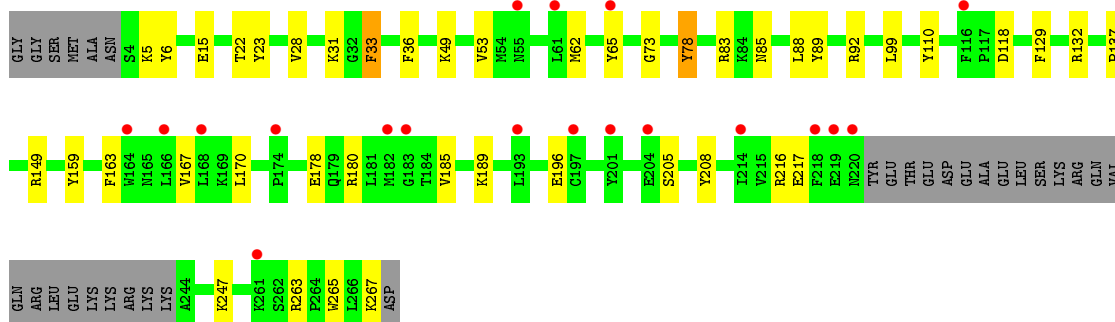
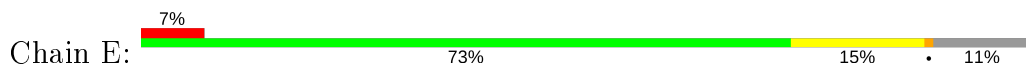


- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase

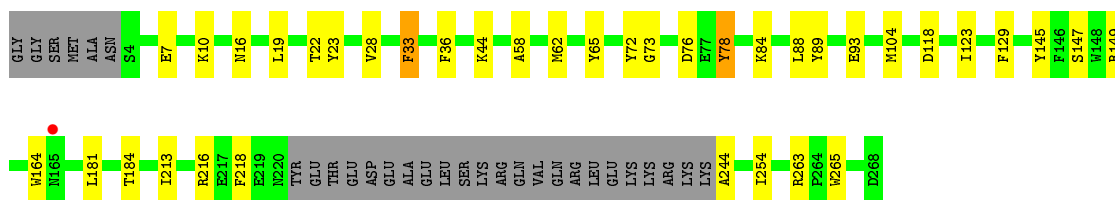




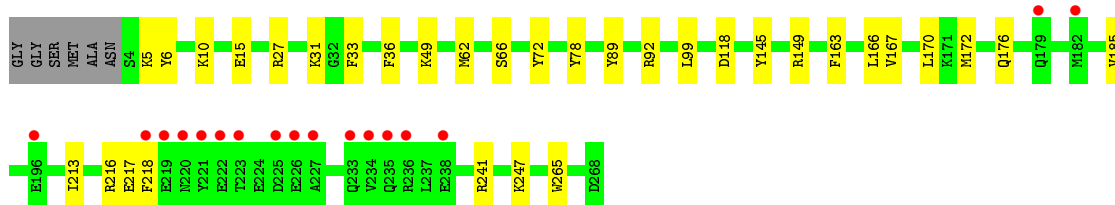
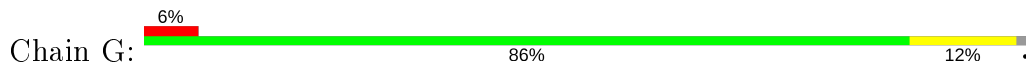
- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase



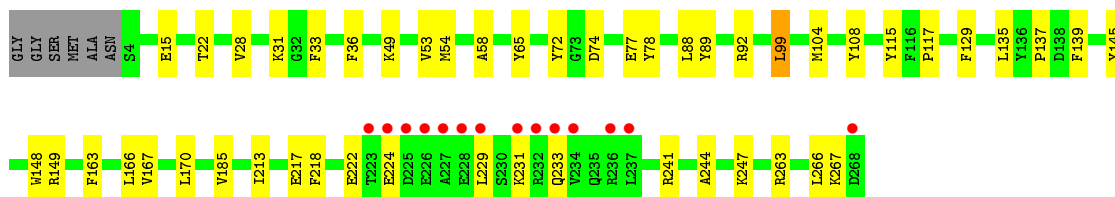
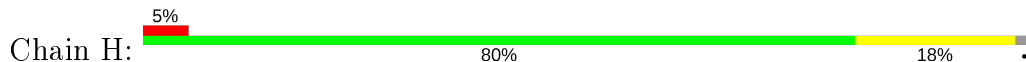
- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase



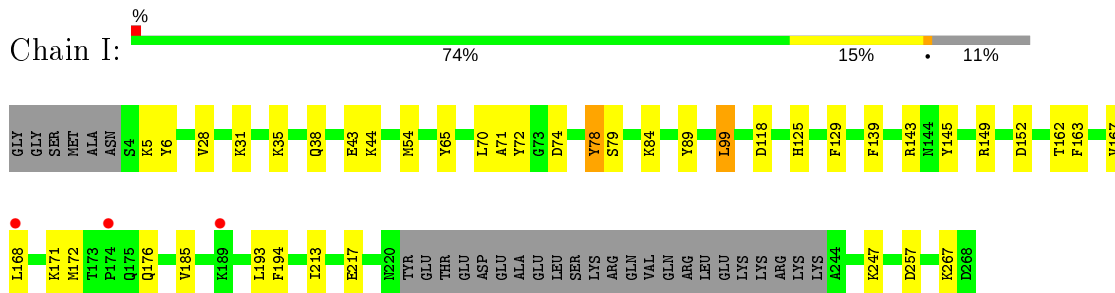
- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase



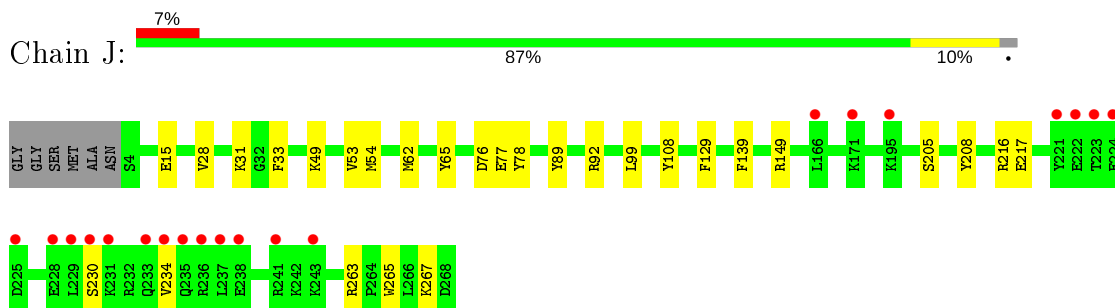
- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase



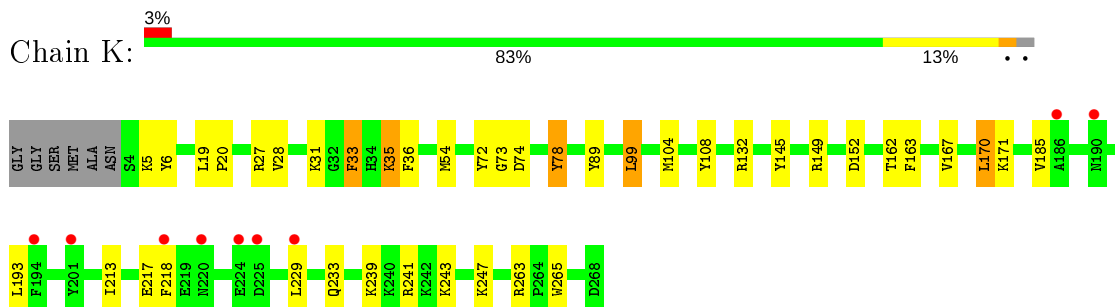
- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase



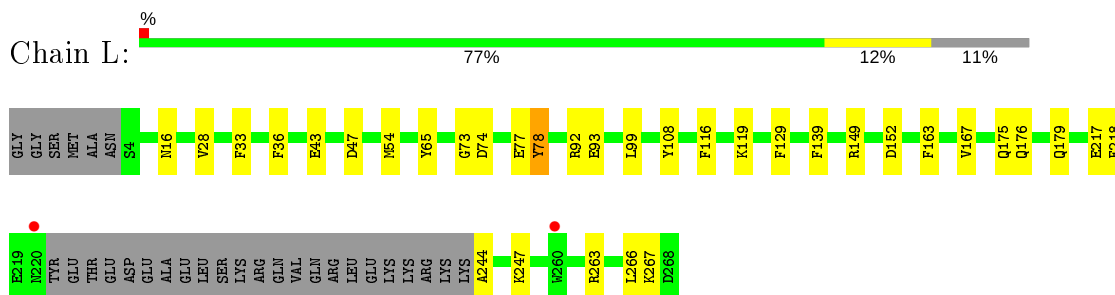
- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase



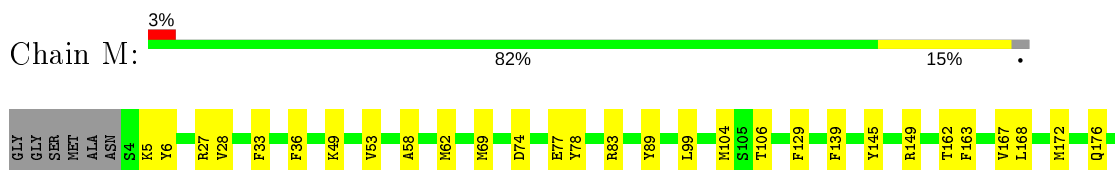
- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase

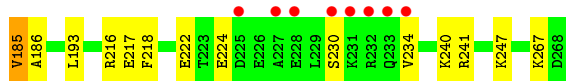


- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase

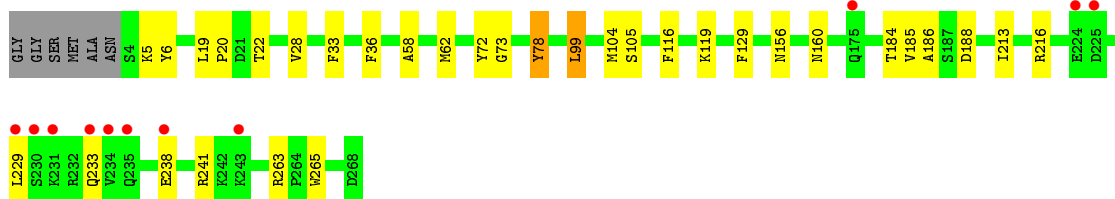
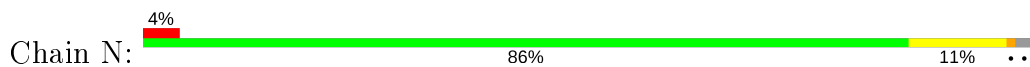


- Molecule 1: Likely histidyl tRNA-specific guanylyltransferase

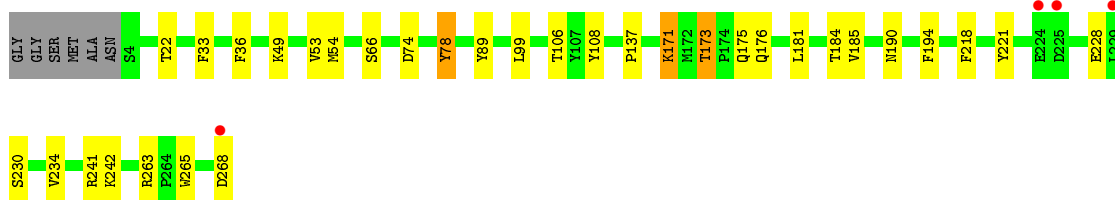
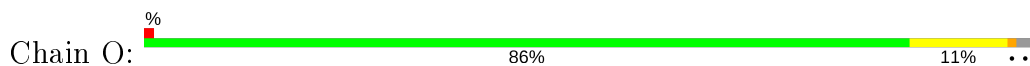




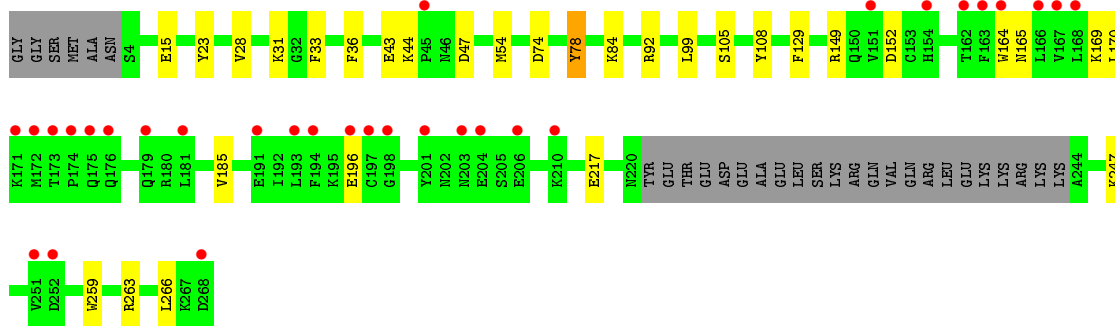
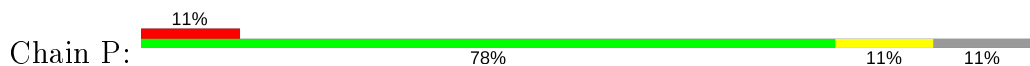
● Molecule 1: Likely histidyl tRNA-specific guanylyltransferase



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● Molecule 1: Likely histidyl tRNA-specific guanylyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	87.43Å 217.56Å 140.91Å 90.00° 102.30° 90.00°	Depositor
Resolution (Å)	44.58 – 3.03 44.58 – 3.03	Depositor EDS
% Data completeness (in resolution range)	98.9 (44.58-3.03) 98.9 (44.58-3.03)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.45 (at 3.01Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, $R_{free}$	0.198 , 0.239 0.199 , 0.233	Depositor DCC
$R_{free}$ test set	7040 reflections (7.14%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	80.9	Xtrriage
Anisotropy	0.337	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 72.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	36341	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	82.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.31% 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: GTP, MG

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	0.28	0/2327	0.47	0/3135
1	B	0.29	0/2327	0.47	0/3135
1	C	0.29	0/2327	0.48	0/3135
1	D	0.30	0/2327	0.48	0/3135
1	E	0.29	0/2115	0.50	0/2854
1	F	0.31	0/2123	0.49	0/2865
1	G	0.30	0/2327	0.50	0/3135
1	H	0.31	0/2327	0.50	0/3135
1	I	0.31	0/2123	0.49	0/2865
1	J	0.29	0/2327	0.49	0/3135
1	K	0.31	0/2327	0.49	0/3135
1	L	0.30	0/2123	0.51	0/2865
1	M	0.32	0/2327	0.51	0/3135
1	N	0.33	0/2327	0.52	1/3135 (0.0%)
1	O	0.33	0/2327	0.52	0/3135
1	P	0.30	0/2123	0.48	0/2865
All	All	0.30	0/36204	0.49	1/48799 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	N	99	LEU	CA-CB-CG	-5.14	103.48	115.30

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	2267	0	2213	23	0
1	B	2267	0	2213	21	0
1	C	2267	0	2213	22	0
1	D	2267	0	2213	20	0
1	E	2057	0	1997	27	0
1	F	2065	0	2001	23	0
1	G	2267	0	2213	25	0
1	H	2267	0	2213	32	0
1	I	2065	0	2001	24	0
1	J	2267	0	2213	16	0
1	K	2267	0	2213	24	0
1	L	2065	0	2001	21	0
1	M	2267	0	2213	24	0
1	N	2267	0	2213	18	0
1	O	2267	0	2213	22	0
1	P	2065	0	2001	16	0
2	A	64	0	24	0	0
2	B	64	0	24	0	0
2	C	64	0	24	0	0
2	D	64	0	24	0	0
2	E	64	0	24	1	0
2	F	64	0	24	4	0
2	G	64	0	24	1	0
2	H	64	0	24	1	0
2	I	64	0	24	0	0
2	J	64	0	24	1	0
2	K	32	0	12	0	0
2	L	96	0	36	2	0
2	M	64	0	24	0	0
2	N	64	0	24	0	0
2	O	96	0	36	1	0
2	P	32	0	12	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
3	D	3	0	0	0	0
3	E	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	4	0	0	0	0
3	G	3	0	0	0	0
3	H	3	0	0	0	0
3	I	3	0	0	0	0
3	J	3	0	0	0	0
3	K	2	0	0	0	0
3	L	4	0	0	0	0
3	M	3	0	0	0	0
3	N	3	0	0	0	0
3	O	3	0	0	0	0
3	P	3	0	0	0	0
4	C	2	0	0	0	0
4	G	1	0	0	0	0
4	H	2	0	0	0	0
4	I	1	0	0	0	0
4	K	1	0	0	0	0
4	M	2	0	0	0	0
4	N	3	0	0	0	0
4	O	3	0	0	0	0
All	All	36341	0	34728	333	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 5.

The worst 5 of 333 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:A:41:GLU:HG2	1:L:175:GLN:HE21	1.34	0.91
1:H:77:GLU:OE1	1:H:149:ARG:NH1	2.14	0.79
1:C:31:LYS:NZ	1:D:65:TYR:OH	2.21	0.74
1:P:263:ARG:HB3	1:P:266:LEU:HD13	1.70	0.73
1:L:92:ARG:NH1	2:L:301:GTP:O2G	2.25	0.70

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	263/271 (97%)	255 (97%)	8 (3%)	0	100	100
1	B	263/271 (97%)	254 (97%)	9 (3%)	0	100	100
1	C	263/271 (97%)	255 (97%)	8 (3%)	0	100	100
1	D	263/271 (97%)	253 (96%)	10 (4%)	0	100	100
1	E	237/271 (88%)	233 (98%)	4 (2%)	0	100	100
1	F	238/271 (88%)	234 (98%)	4 (2%)	0	100	100
1	G	263/271 (97%)	255 (97%)	8 (3%)	0	100	100
1	H	263/271 (97%)	255 (97%)	8 (3%)	0	100	100
1	I	238/271 (88%)	234 (98%)	4 (2%)	0	100	100
1	J	263/271 (97%)	255 (97%)	8 (3%)	0	100	100
1	K	263/271 (97%)	255 (97%)	8 (3%)	0	100	100
1	L	238/271 (88%)	234 (98%)	4 (2%)	0	100	100
1	M	263/271 (97%)	255 (97%)	8 (3%)	0	100	100
1	N	263/271 (97%)	255 (97%)	8 (3%)	0	100	100
1	O	263/271 (97%)	255 (97%)	8 (3%)	0	100	100
1	P	238/271 (88%)	234 (98%)	4 (2%)	0	100	100
All	All	4082/4336 (94%)	3971 (97%)	111 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	249/252 (99%)	248 (100%)	1 (0%)	91	96
1	B	249/252 (99%)	248 (100%)	1 (0%)	91	96
1	C	249/252 (99%)	248 (100%)	1 (0%)	91	96
1	D	249/252 (99%)	247 (99%)	2 (1%)	81	92
1	E	226/252 (90%)	221 (98%)	5 (2%)	52	79
1	F	227/252 (90%)	225 (99%)	2 (1%)	78	91
1	G	249/252 (99%)	247 (99%)	2 (1%)	81	92
1	H	249/252 (99%)	244 (98%)	5 (2%)	55	81
1	I	227/252 (90%)	219 (96%)	8 (4%)	36	69
1	J	249/252 (99%)	247 (99%)	2 (1%)	81	92
1	K	249/252 (99%)	241 (97%)	8 (3%)	39	72
1	L	227/252 (90%)	223 (98%)	4 (2%)	59	83
1	M	249/252 (99%)	244 (98%)	5 (2%)	55	81
1	N	249/252 (99%)	248 (100%)	1 (0%)	91	96
1	O	249/252 (99%)	243 (98%)	6 (2%)	49	78
1	P	227/252 (90%)	221 (97%)	6 (3%)	46	76
All	All	3873/4032 (96%)	3814 (98%)	59 (2%)	65	86

5 of 59 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	194	PHE
1	K	99	LEU
1	P	78	TYR
1	J	33	PHE
1	K	33	PHE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 9 such sidechains are listed below:

Mol	Chain	Res	Type
1	I	38	GLN
1	P	200	ASN
1	N	233	GLN
1	B	46	ASN
1	L	175	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 80 ligands modelled in this entry, 48 are monoatomic - leaving 32 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	GTP	A	401	3	26,34,34	0.98	1 (3%)	33,54,54	1.70	7 (21%)
2	GTP	D	301	3	26,34,34	0.97	1 (3%)	33,54,54	1.99	7 (21%)
2	GTP	B	402	3	26,34,34	0.94	1 (3%)	33,54,54	1.87	7 (21%)
2	GTP	A	402	3	26,34,34	0.93	1 (3%)	33,54,54	1.84	7 (21%)
2	GTP	L	304	3	26,34,34	0.93	1 (3%)	33,54,54	1.89	7 (21%)
2	GTP	J	402	3	26,34,34	0.97	1 (3%)	33,54,54	1.95	7 (21%)
2	GTP	I	402	3	26,34,34	0.95	1 (3%)	33,54,54	1.94	7 (21%)
2	GTP	N	401	3	26,34,34	0.99	2 (7%)	33,54,54	1.70	7 (21%)
2	GTP	E	404	3	26,34,34	1.00	2 (7%)	33,54,54	1.98	8 (24%)
2	GTP	J	401	3	26,34,34	0.96	1 (3%)	33,54,54	1.79	7 (21%)
2	GTP	G	401	3	26,34,34	0.99	1 (3%)	33,54,54	1.71	7 (21%)
2	GTP	F	301	3	26,34,34	0.96	1 (3%)	33,54,54	1.79	7 (21%)
2	GTP	I	401	3	26,34,34	0.99	1 (3%)	33,54,54	1.76	7 (21%)
2	GTP	D	302	3	26,34,34	0.99	1 (3%)	33,54,54	1.73	7 (21%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	GTP	K	401	3	26,34,34	0.97	1 (3%)	33,54,54	1.76	7 (21%)
2	GTP	N	402	3	26,34,34	0.95	1 (3%)	33,54,54	1.90	8 (24%)
2	GTP	M	402	3	26,34,34	0.94	1 (3%)	33,54,54	1.87	9 (27%)
2	GTP	F	303	3	26,34,34	1.00	1 (3%)	33,54,54	1.87	8 (24%)
2	GTP	C	405	3	26,34,34	0.95	1 (3%)	33,54,54	1.92	8 (24%)
2	GTP	O	401	3	26,34,34	0.99	1 (3%)	33,54,54	1.83	9 (27%)
2	GTP	H	402	3	26,34,34	0.94	2 (7%)	33,54,54	1.94	7 (21%)
2	GTP	L	301	3	26,34,34	0.95	1 (3%)	33,54,54	1.89	7 (21%)
2	GTP	E	401	3	26,34,34	0.92	1 (3%)	33,54,54	1.71	6 (18%)
2	GTP	P	401	3	26,34,34	0.99	1 (3%)	33,54,54	1.75	8 (24%)
2	GTP	C	401	3	26,34,34	0.97	1 (3%)	33,54,54	1.74	7 (21%)
2	GTP	M	401	3	26,34,34	1.03	2 (7%)	33,54,54	1.78	8 (24%)
2	GTP	G	402	3	26,34,34	0.93	1 (3%)	33,54,54	1.86	7 (21%)
2	GTP	L	303	3	26,34,34	1.00	1 (3%)	33,54,54	1.81	8 (24%)
2	GTP	O	402	3	26,34,34	0.91	1 (3%)	33,54,54	1.88	7 (21%)
2	GTP	O	406	3	26,34,34	0.96	1 (3%)	33,54,54	1.98	8 (24%)
2	GTP	B	401	3	26,34,34	0.97	1 (3%)	33,54,54	1.80	7 (21%)
2	GTP	H	401	3	26,34,34	0.98	1 (3%)	33,54,54	1.80	9 (27%)

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	GTP	A	401	3	-	8/18/38/38	0/3/3/3
2	GTP	D	301	3	-	9/18/38/38	0/3/3/3
2	GTP	B	402	3	-	11/18/38/38	0/3/3/3
2	GTP	A	402	3	-	9/18/38/38	0/3/3/3
2	GTP	L	304	3	-	10/18/38/38	0/3/3/3
2	GTP	J	402	3	-	7/18/38/38	0/3/3/3
2	GTP	I	402	3	-	10/18/38/38	0/3/3/3
2	GTP	N	401	3	-	2/18/38/38	0/3/3/3
2	GTP	E	404	3	-	5/18/38/38	0/3/3/3
2	GTP	J	401	3	-	8/18/38/38	0/3/3/3
2	GTP	G	401	3	-	6/18/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GTP	F	301	3	-	6/18/38/38	0/3/3/3
2	GTP	I	401	3	-	4/18/38/38	0/3/3/3
2	GTP	D	302	3	-	7/18/38/38	0/3/3/3
2	GTP	K	401	3	-	5/18/38/38	0/3/3/3
2	GTP	N	402	3	-	6/18/38/38	0/3/3/3
2	GTP	M	402	3	-	10/18/38/38	0/3/3/3
2	GTP	F	303	3	-	5/18/38/38	0/3/3/3
2	GTP	C	405	3	-	4/18/38/38	0/3/3/3
2	GTP	O	401	3	-	6/18/38/38	0/3/3/3
2	GTP	H	402	3	-	3/18/38/38	0/3/3/3
2	GTP	L	301	3	-	6/18/38/38	0/3/3/3
2	GTP	E	401	3	-	10/18/38/38	0/3/3/3
2	GTP	P	401	3	-	5/18/38/38	0/3/3/3
2	GTP	C	401	3	-	6/18/38/38	0/3/3/3
2	GTP	M	401	3	-	5/18/38/38	0/3/3/3
2	GTP	G	402	3	-	7/18/38/38	0/3/3/3
2	GTP	L	303	3	-	5/18/38/38	0/3/3/3
2	GTP	O	402	3	-	9/18/38/38	0/3/3/3
2	GTP	O	406	3	-	5/18/38/38	0/3/3/3
2	GTP	B	401	3	-	10/18/38/38	0/3/3/3
2	GTP	H	401	3	-	5/18/38/38	0/3/3/3

The worst 5 of 36 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	303	GTP	C6-N1	3.42	1.39	1.33
2	N	401	GTP	C6-N1	3.36	1.38	1.33
2	F	303	GTP	C6-N1	3.28	1.38	1.33
2	M	401	GTP	C6-N1	3.28	1.38	1.33
2	I	401	GTP	C6-N1	3.28	1.38	1.33

The worst 5 of 237 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	402	GTP	N3-C2-N1	-5.74	119.57	127.22
2	I	402	GTP	N3-C2-N1	-5.73	119.58	127.22
2	D	301	GTP	N3-C2-N1	-5.72	119.60	127.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	401	GTP	N3-C2-N1	-5.66	119.68	127.22
2	E	404	GTP	N3-C2-N1	-5.65	119.69	127.22

There are no chirality outliers.

5 of 214 torsion outliers are listed below:

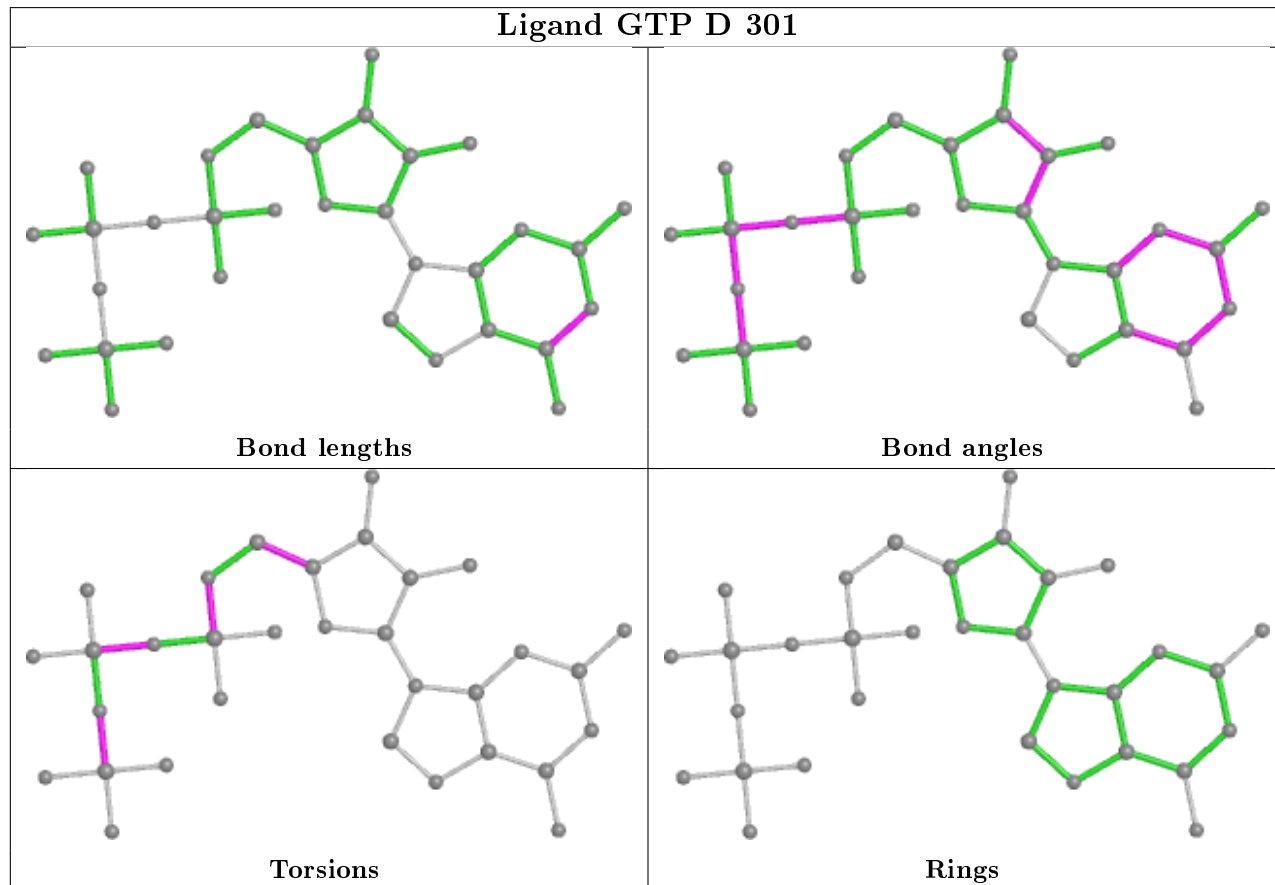
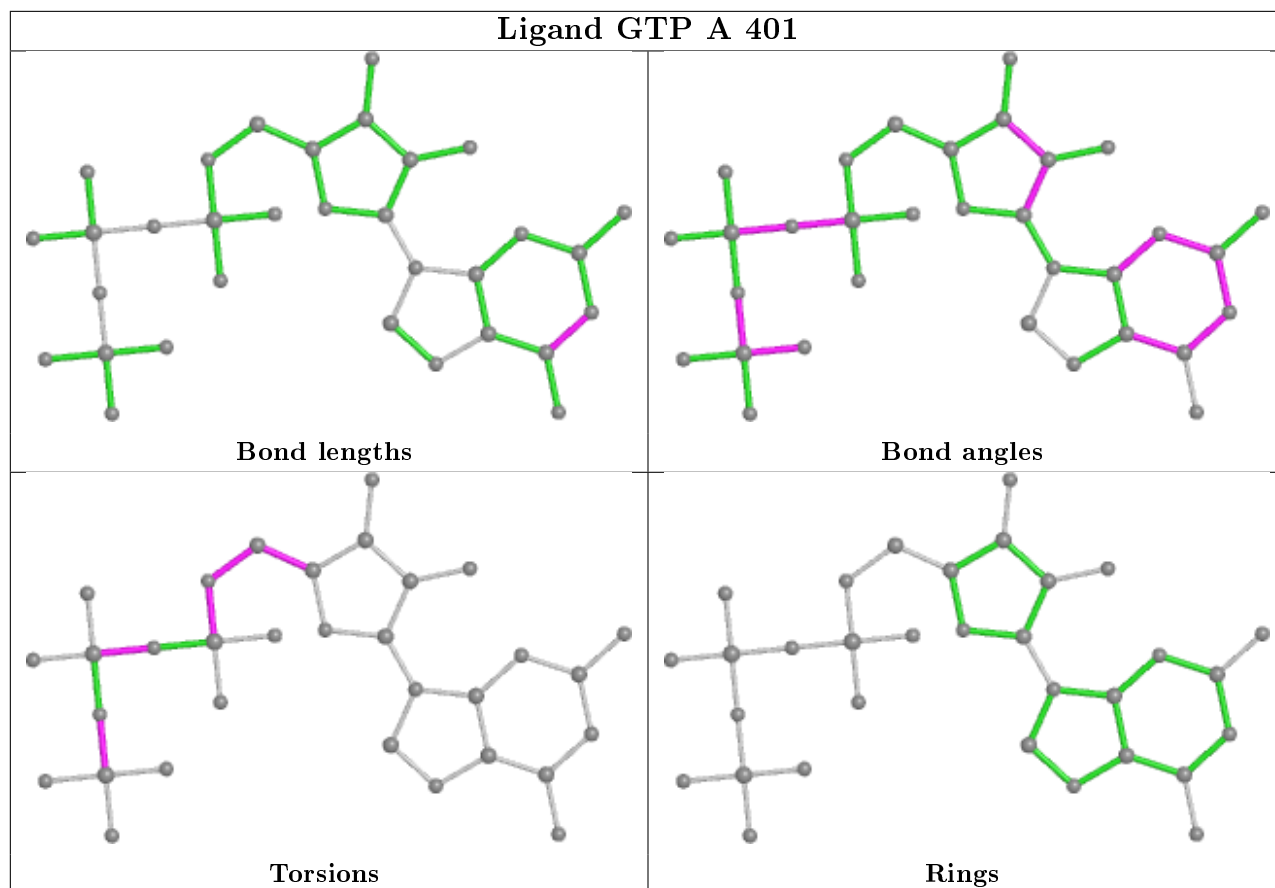
Mol	Chain	Res	Type	Atoms
2	A	401	GTP	C5'-O5'-PA-O1A
2	A	401	GTP	O4'-C4'-C5'-O5'
2	D	301	GTP	PB-O3B-PG-O3G
2	D	301	GTP	C5'-O5'-PA-O1A
2	D	301	GTP	C5'-O5'-PA-O2A

There are no ring outliers.

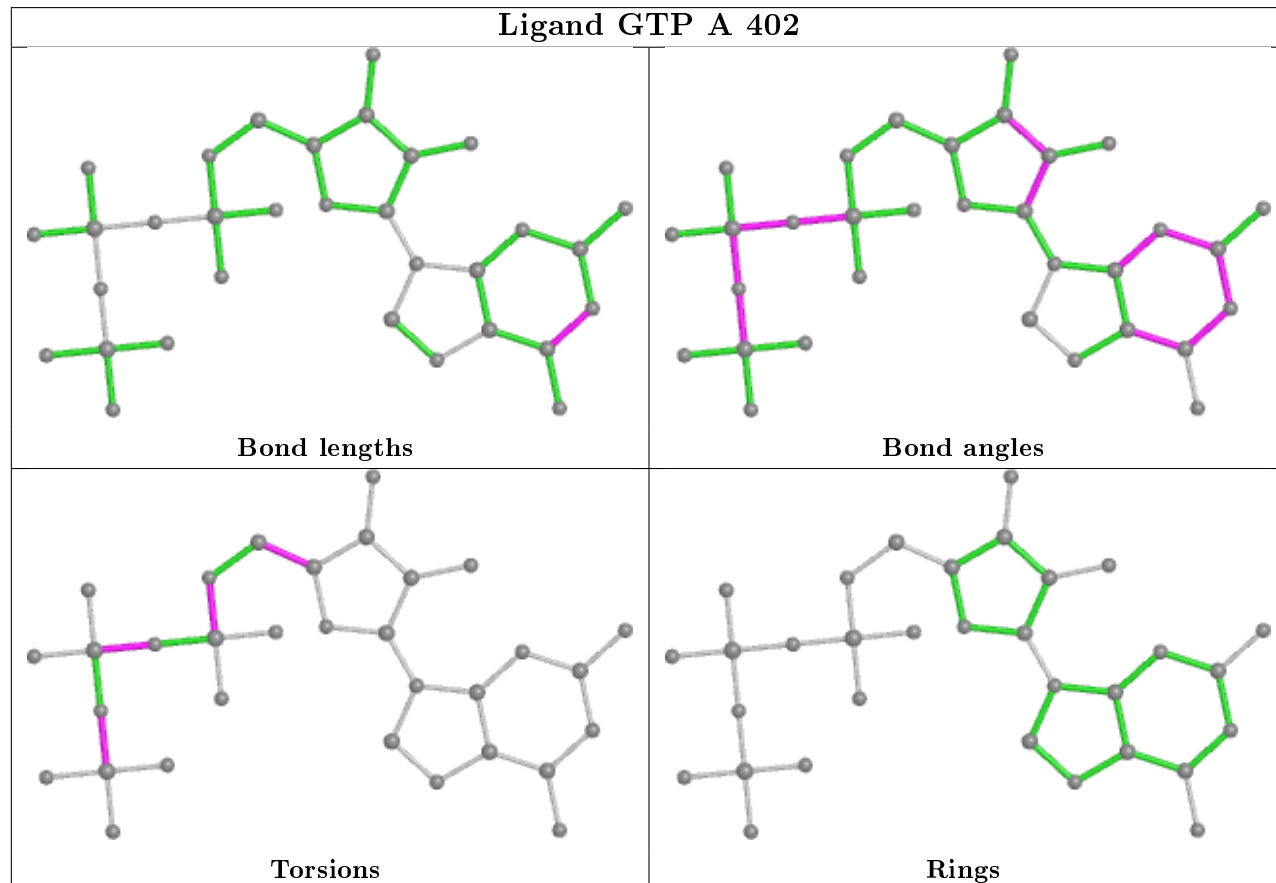
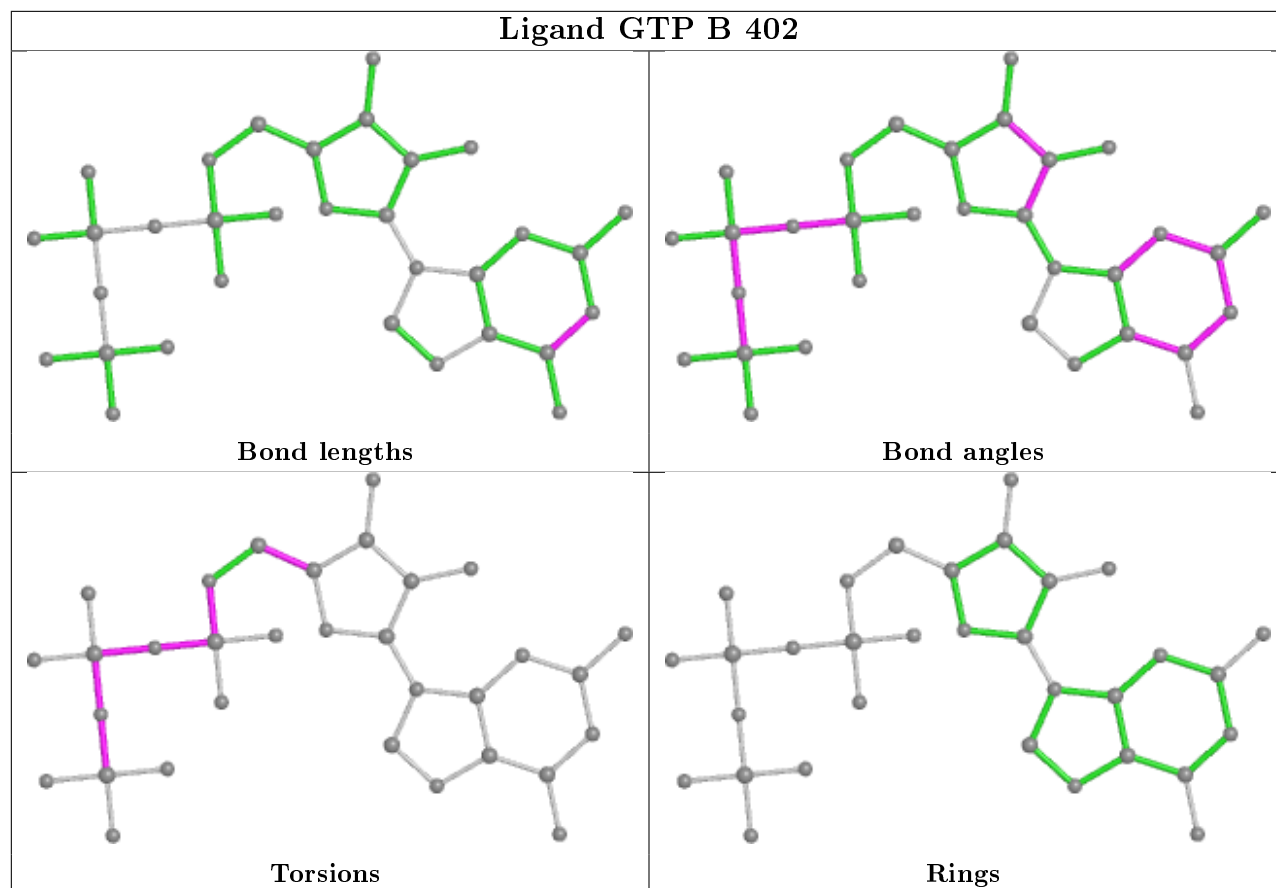
8 monomers are involved in 11 short contacts:

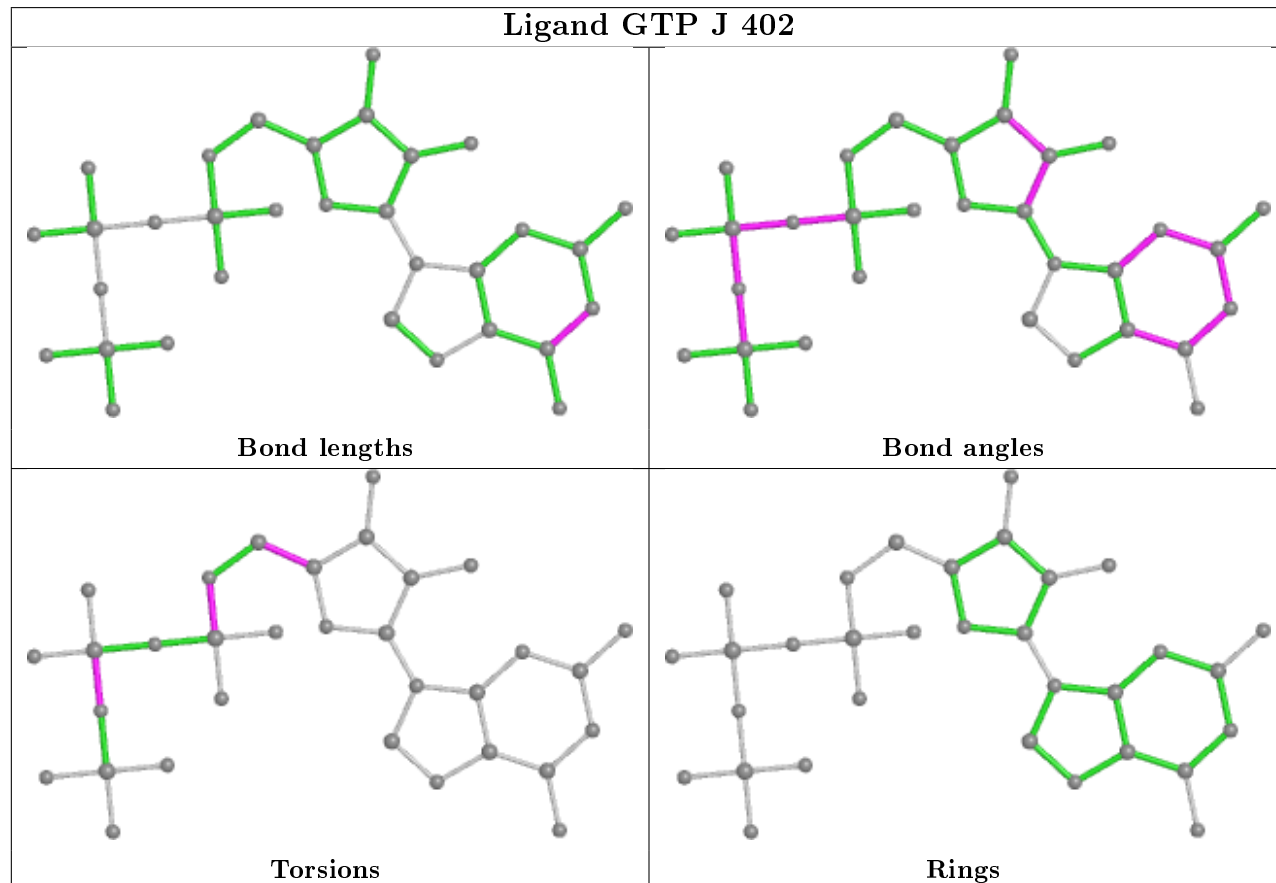
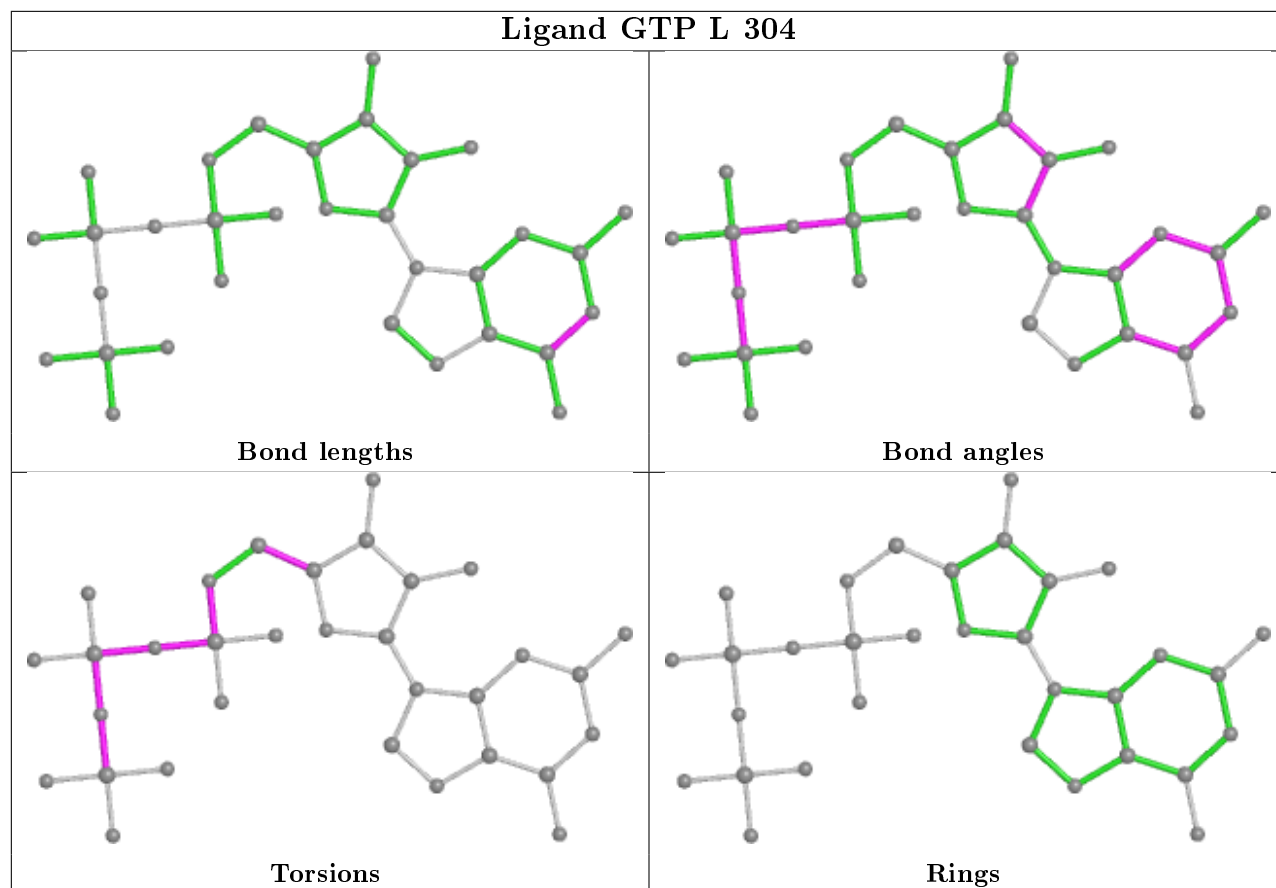
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	404	GTP	1	0
2	J	401	GTP	1	0
2	F	301	GTP	3	0
2	F	303	GTP	1	0
2	H	402	GTP	1	0
2	L	301	GTP	2	0
2	G	402	GTP	1	0
2	O	402	GTP	1	0

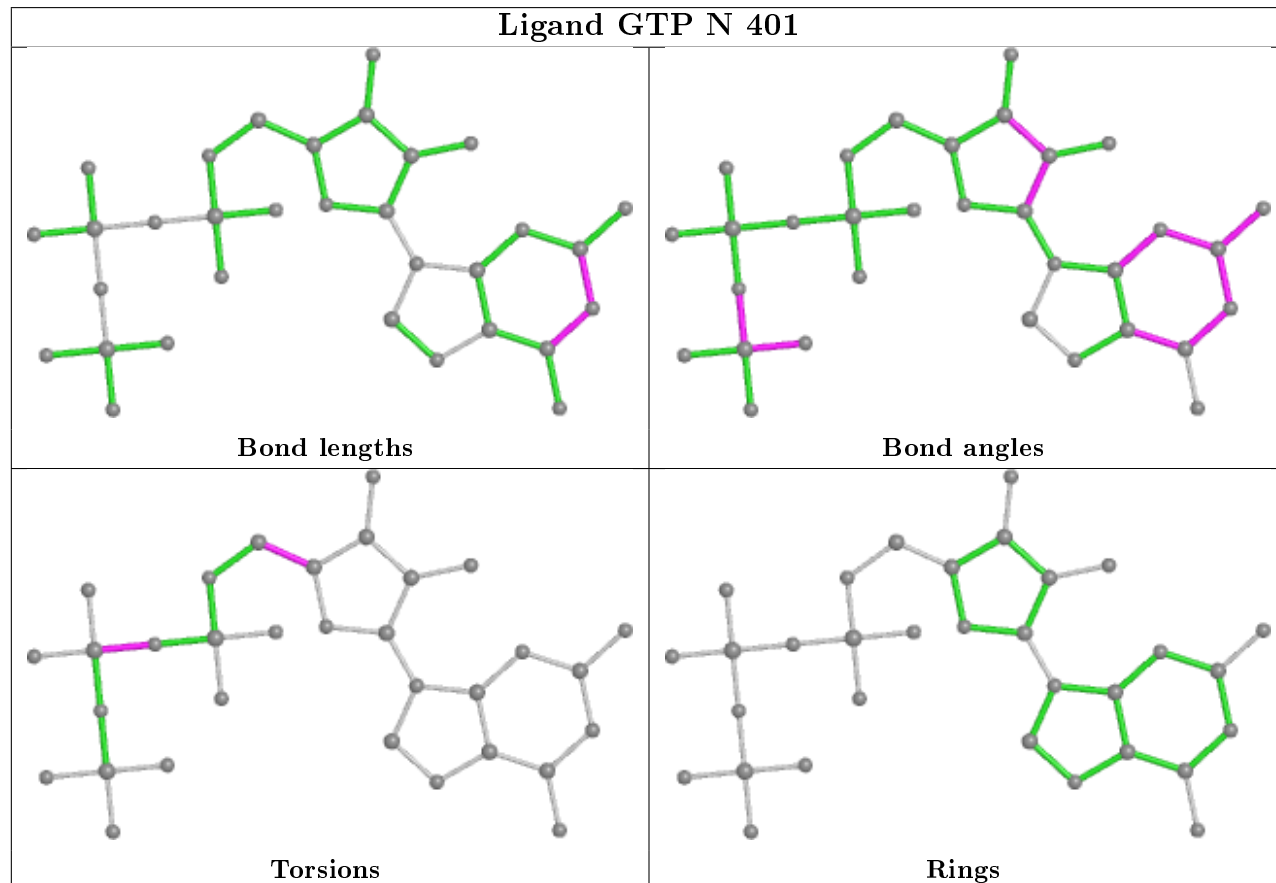
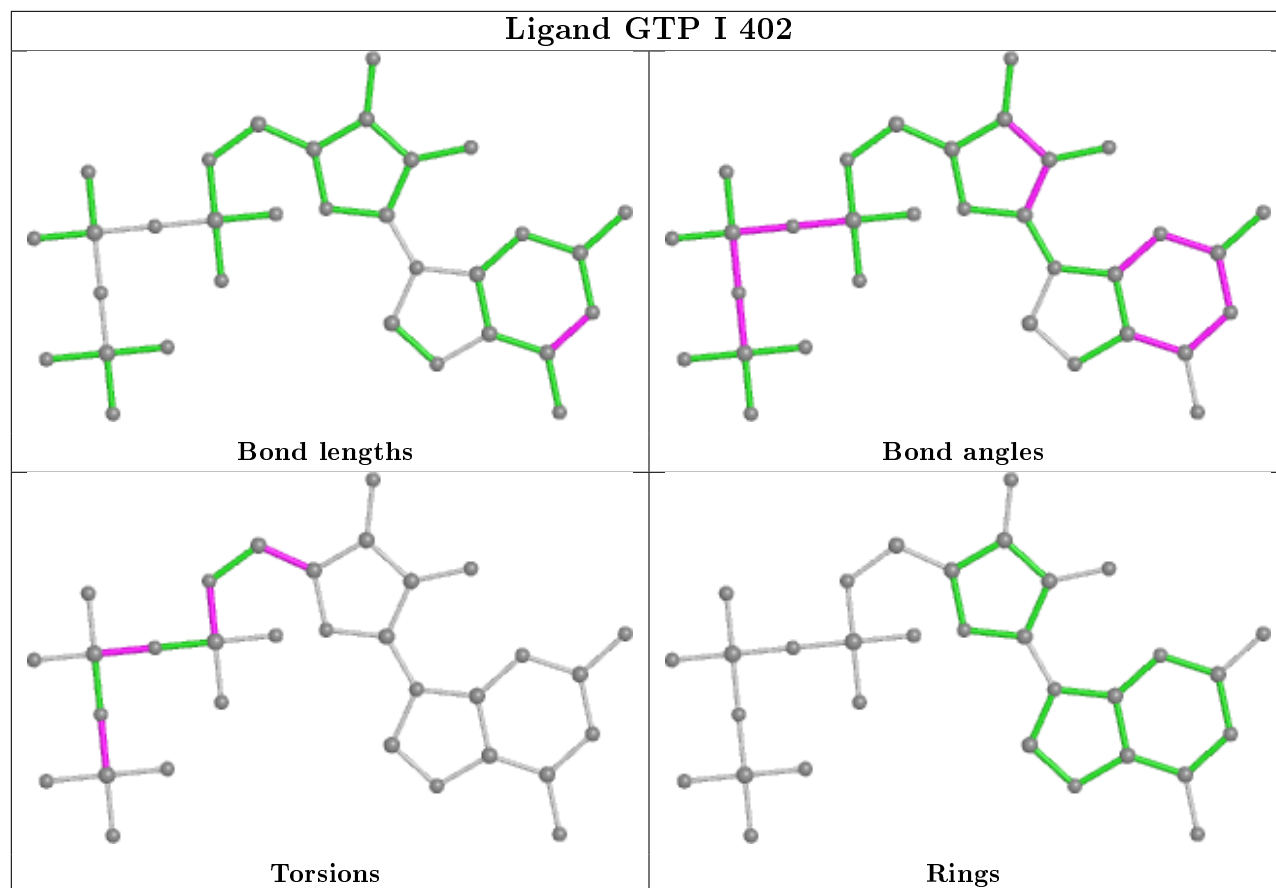
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

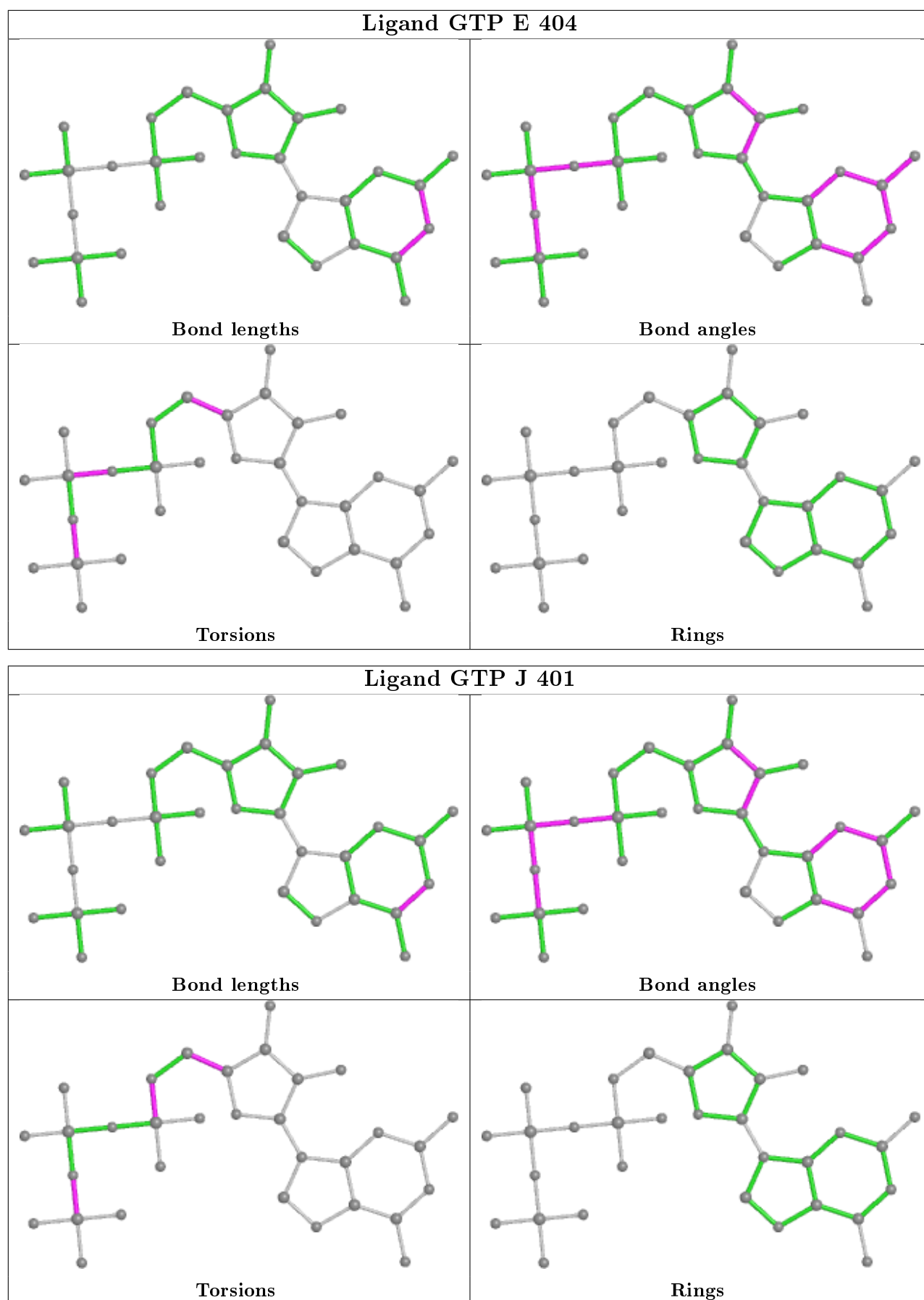


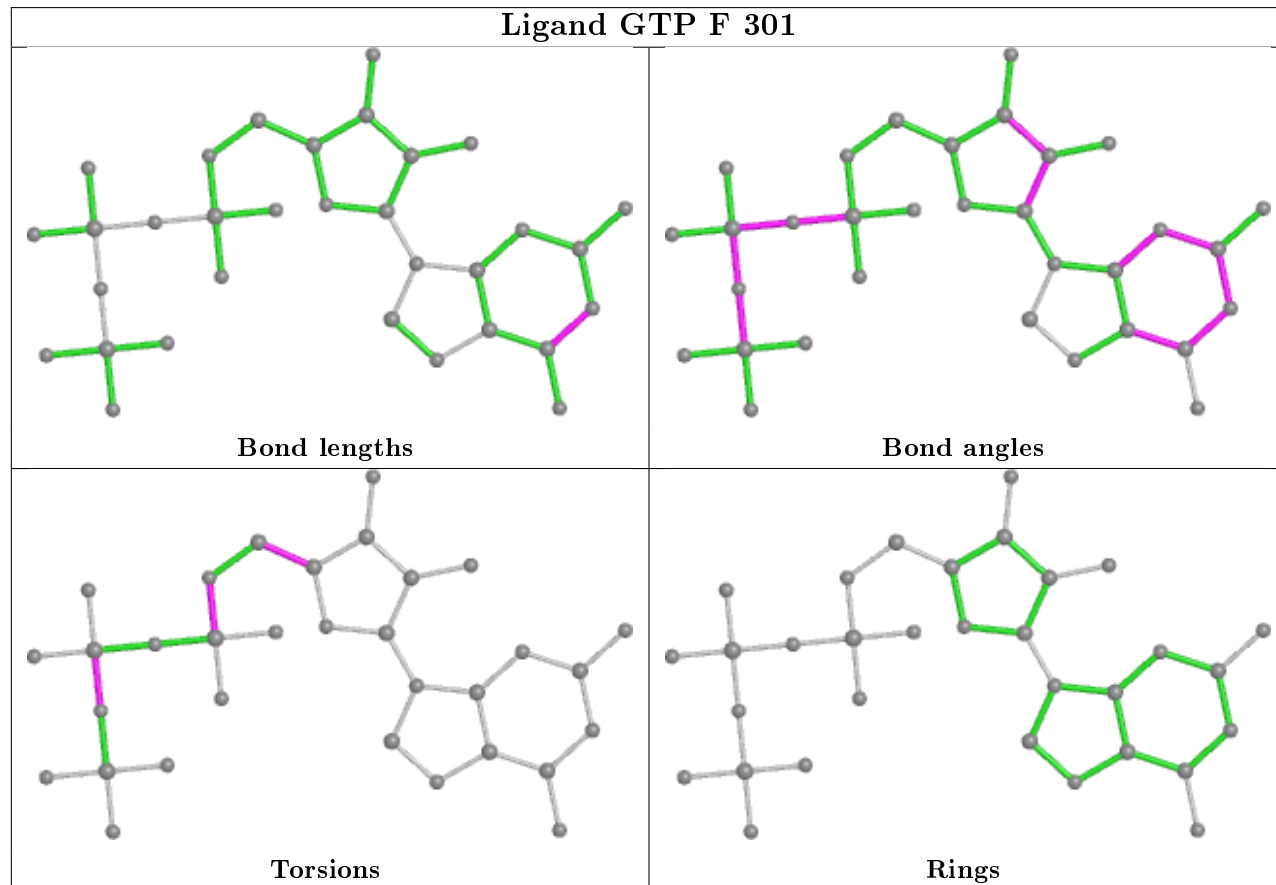
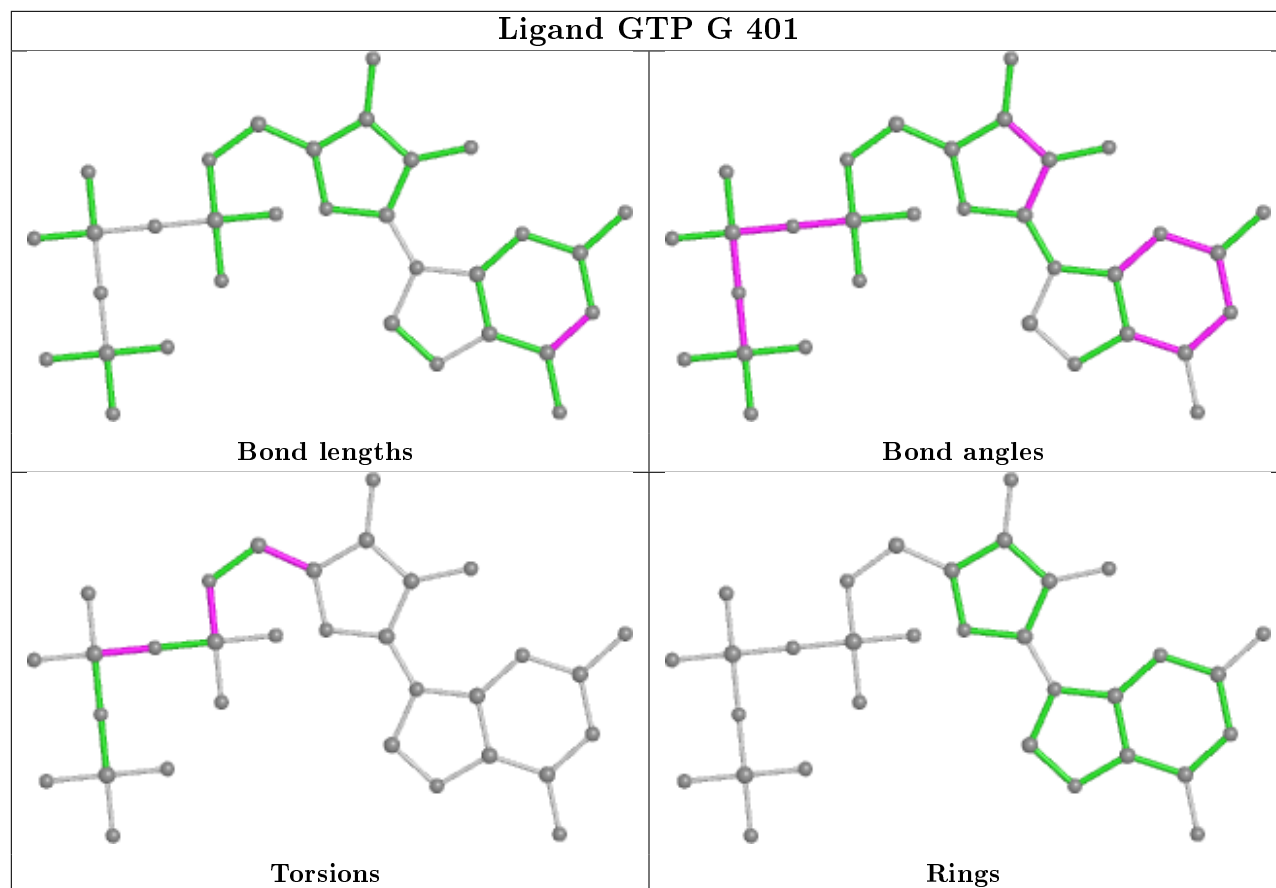


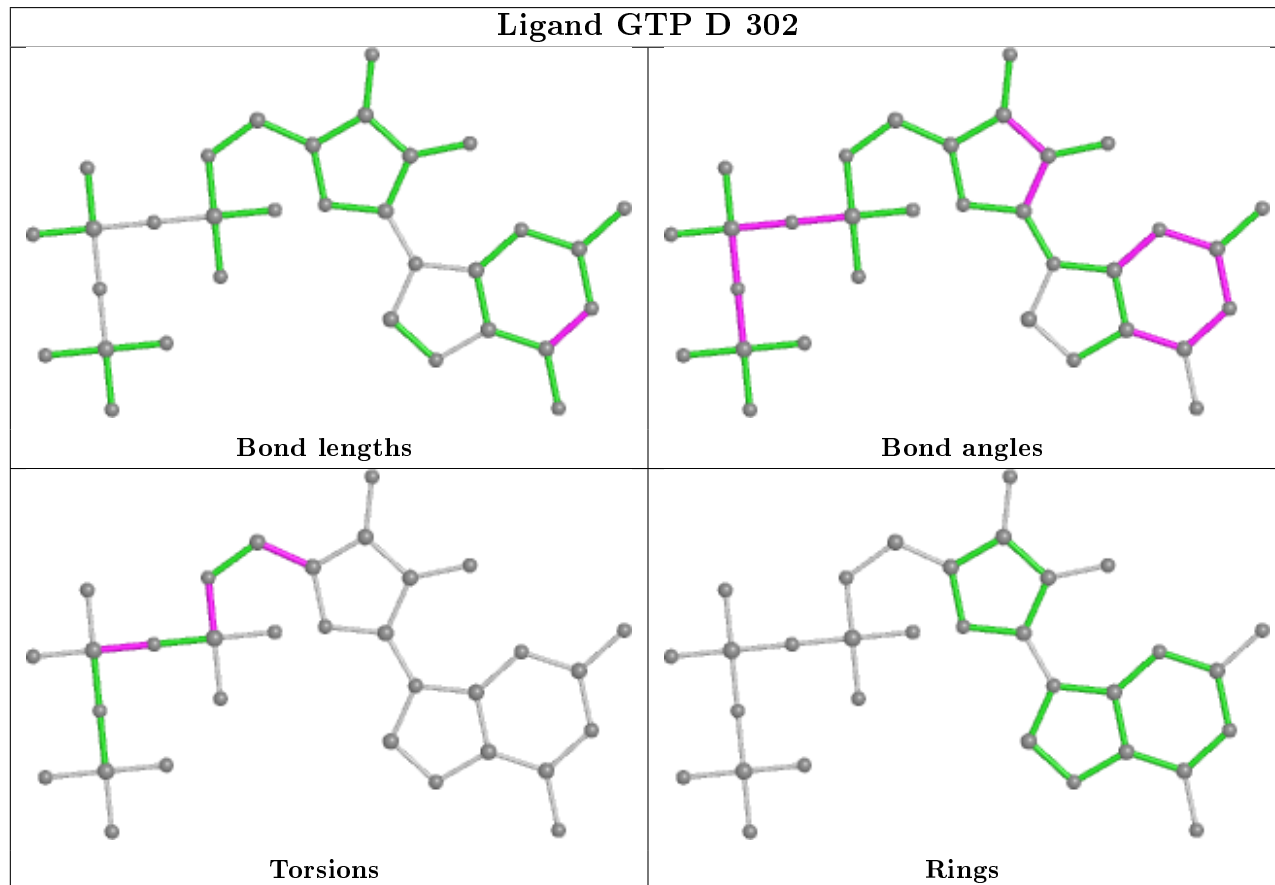
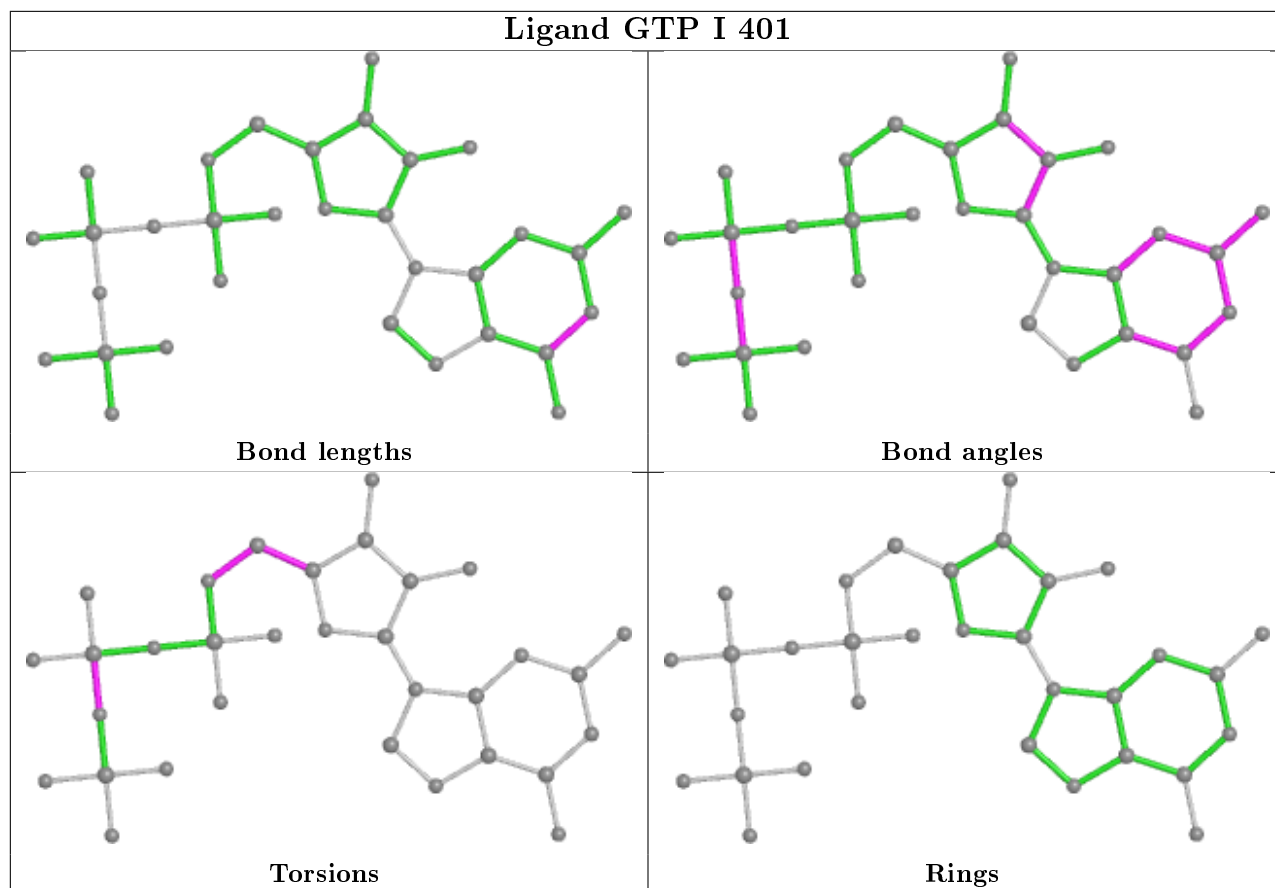


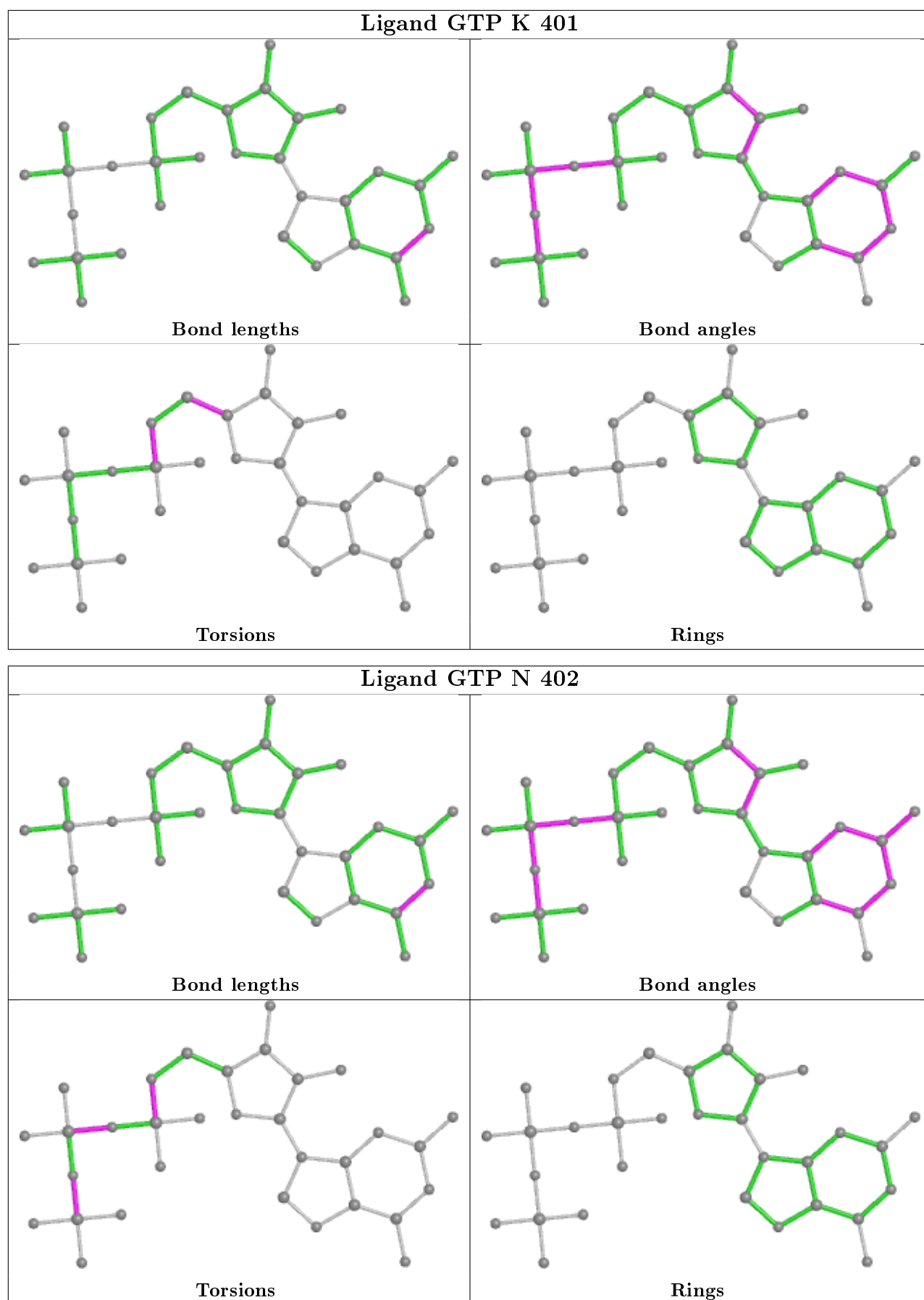


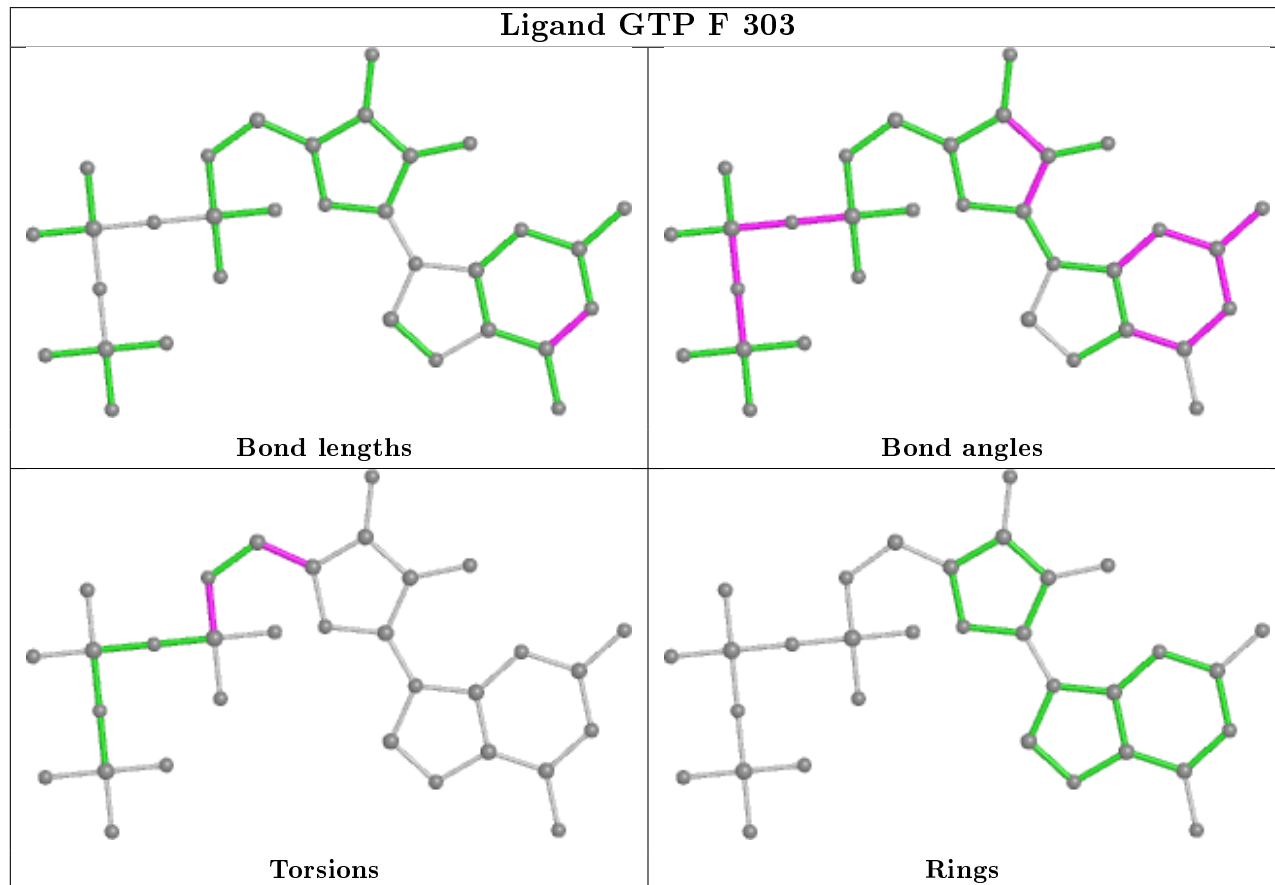
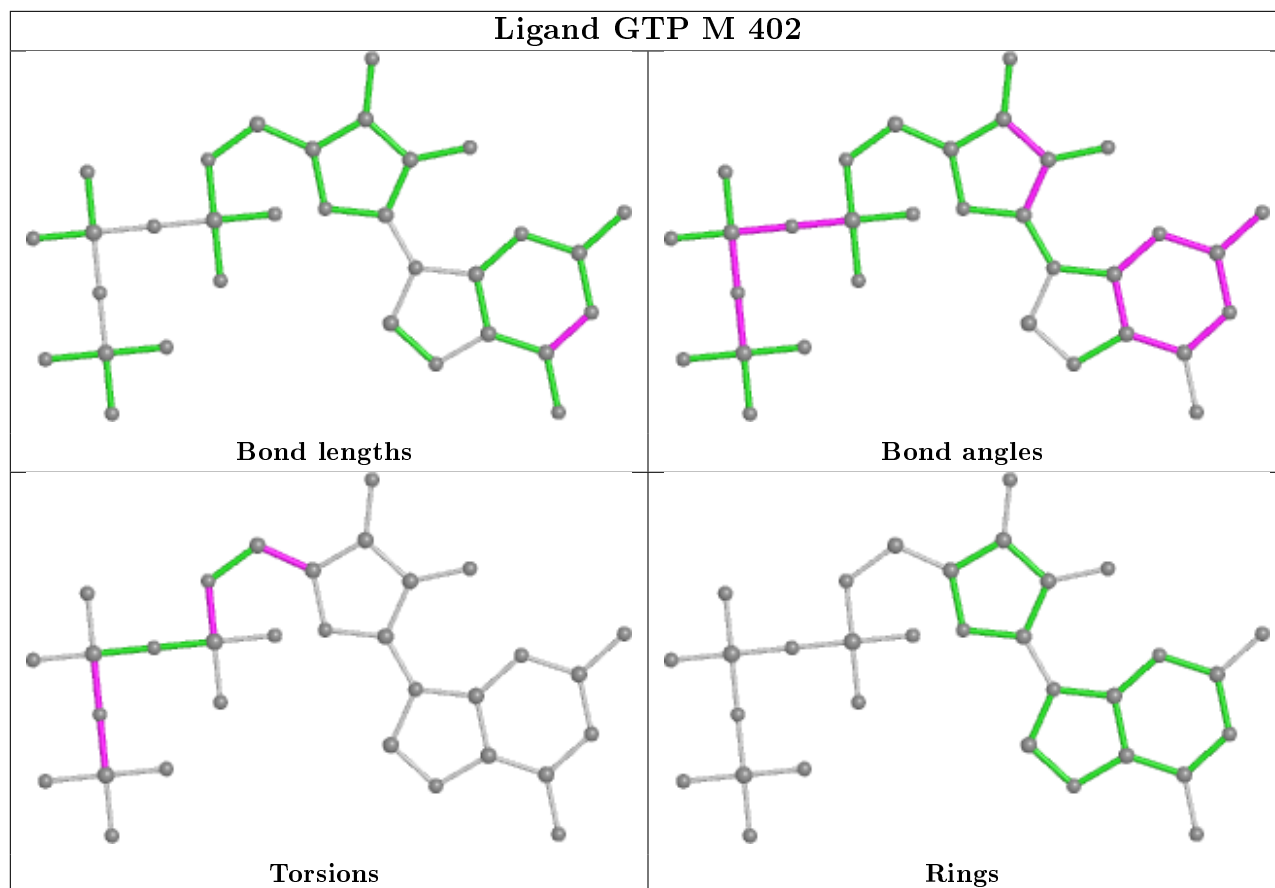




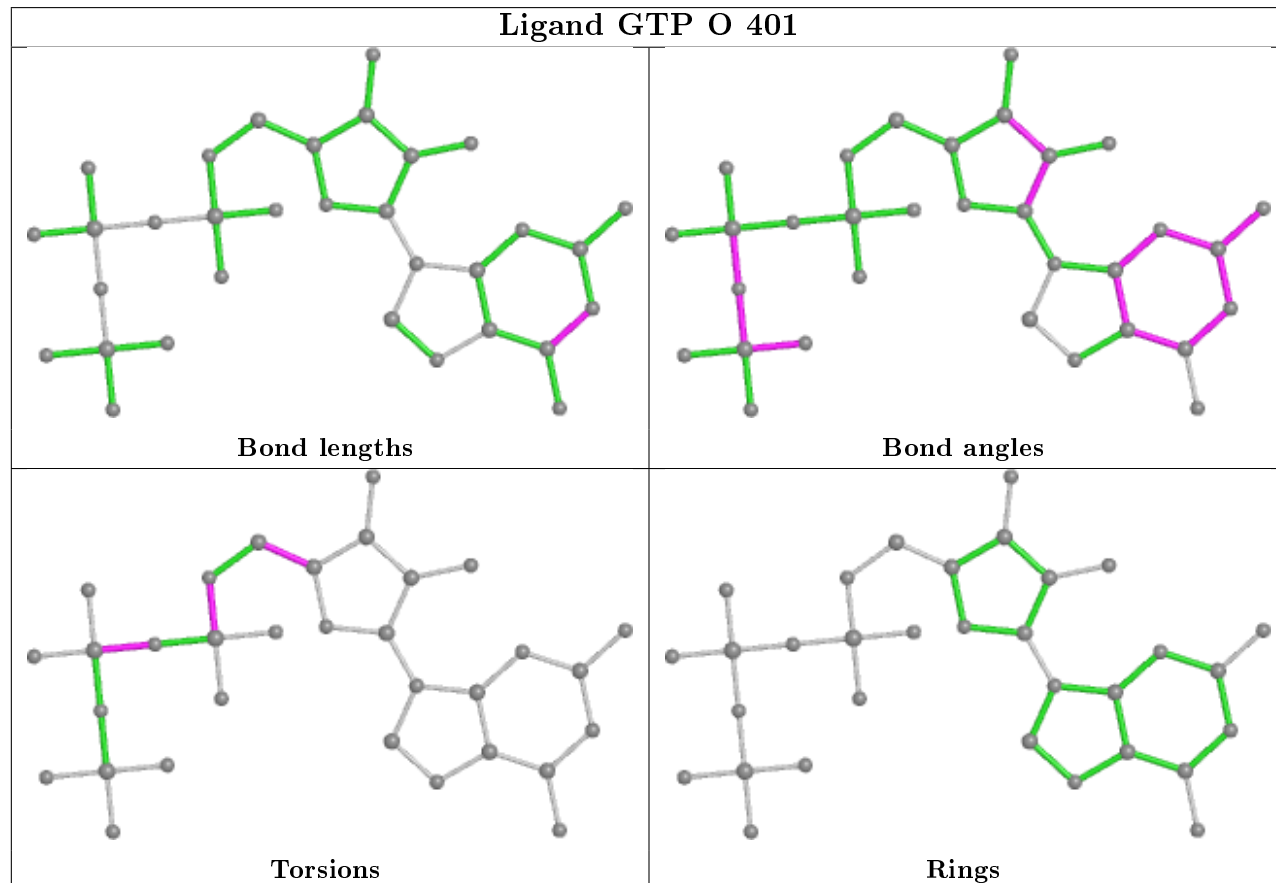
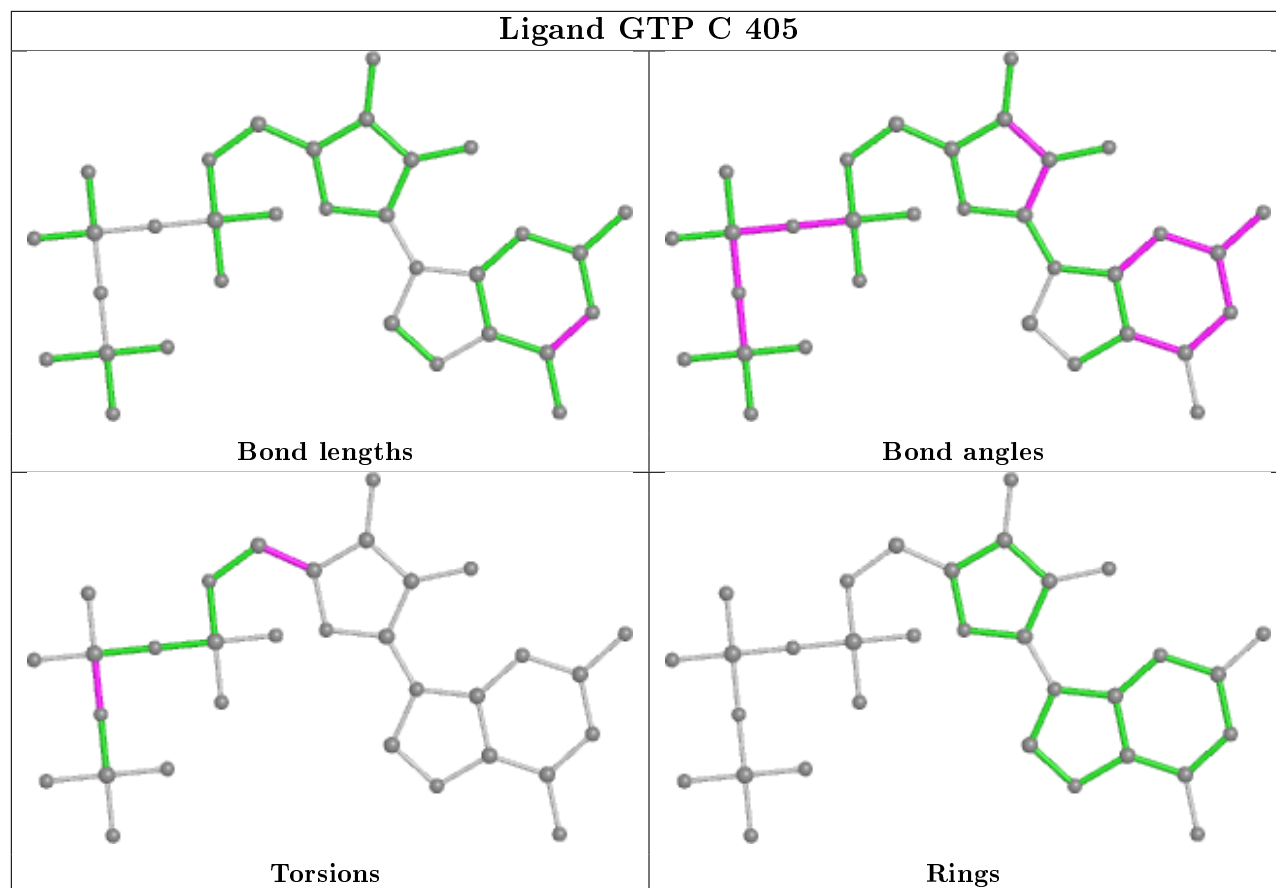


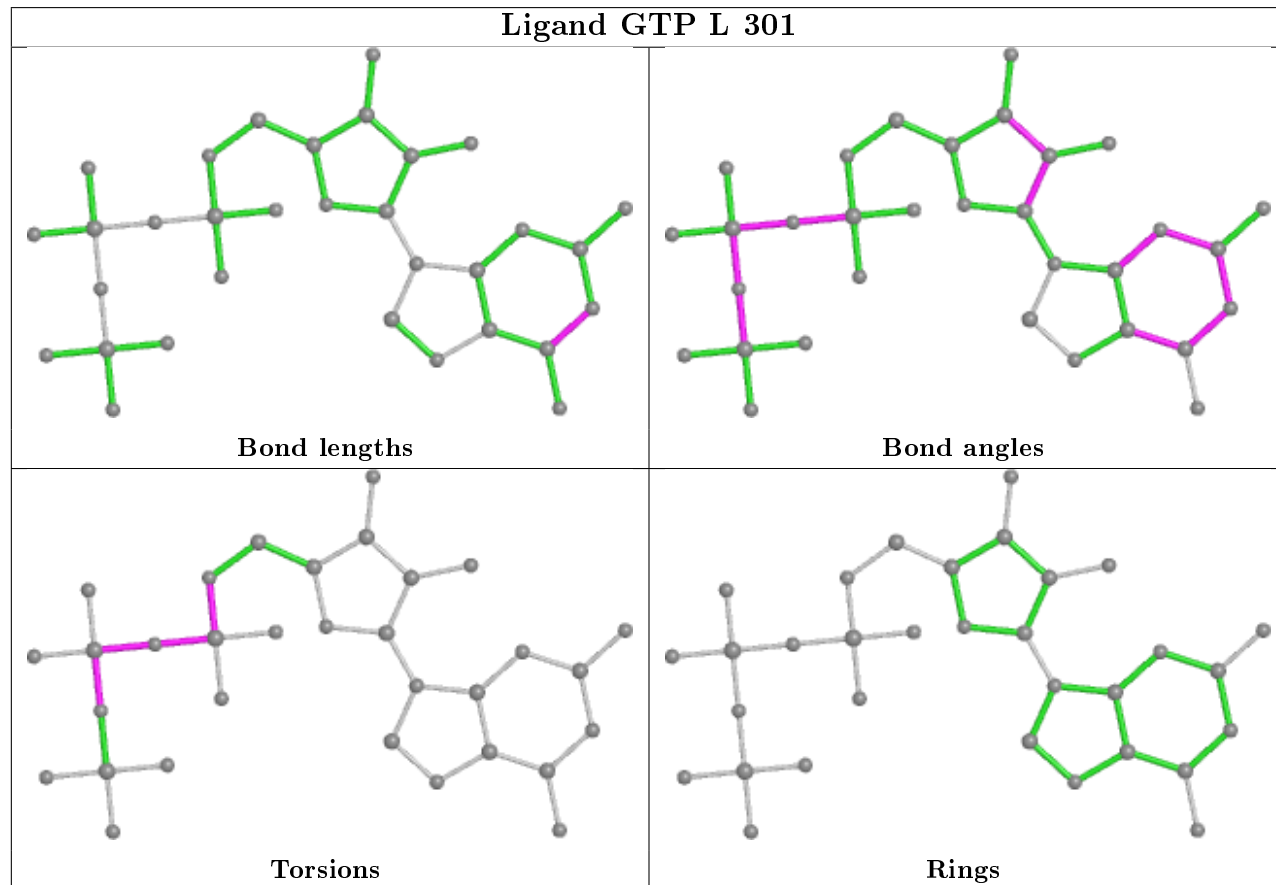
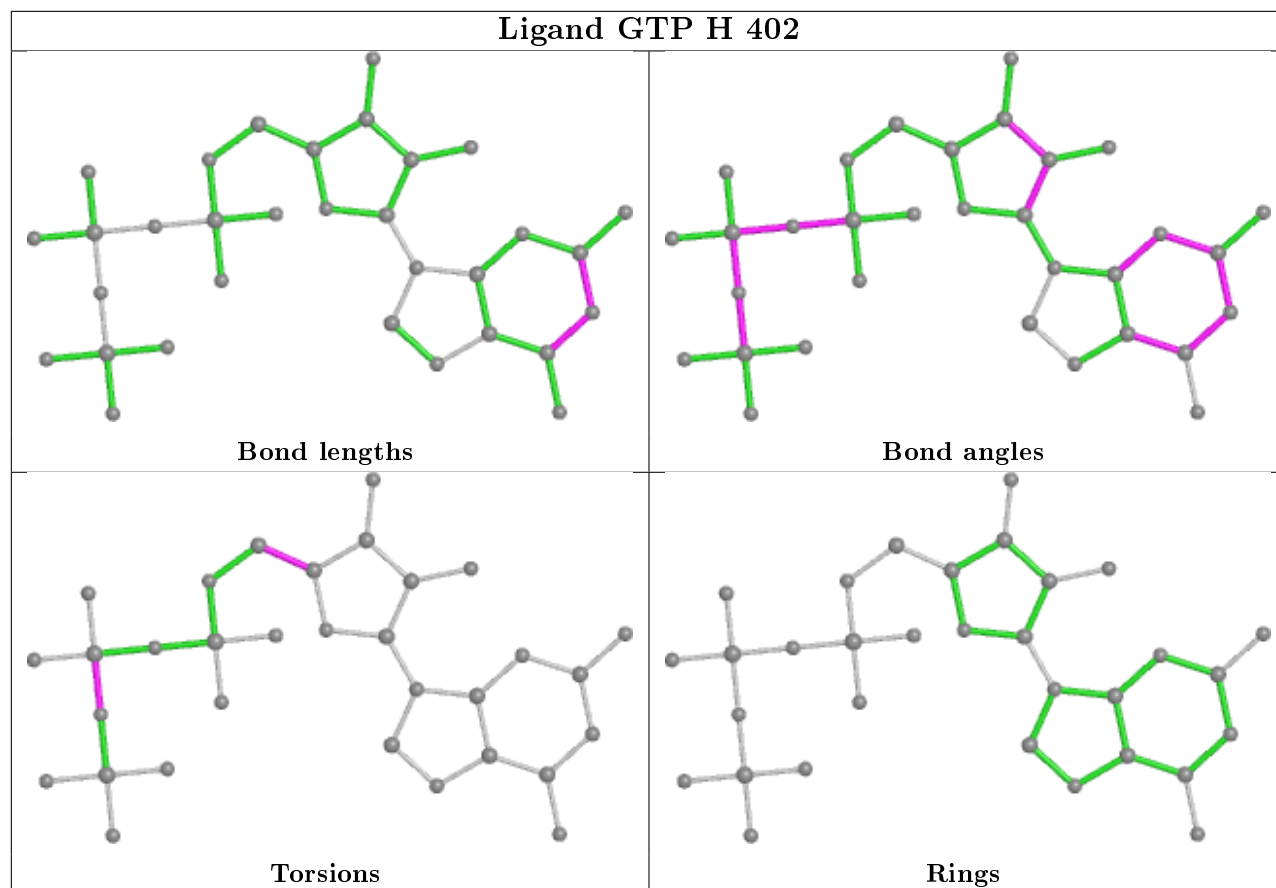


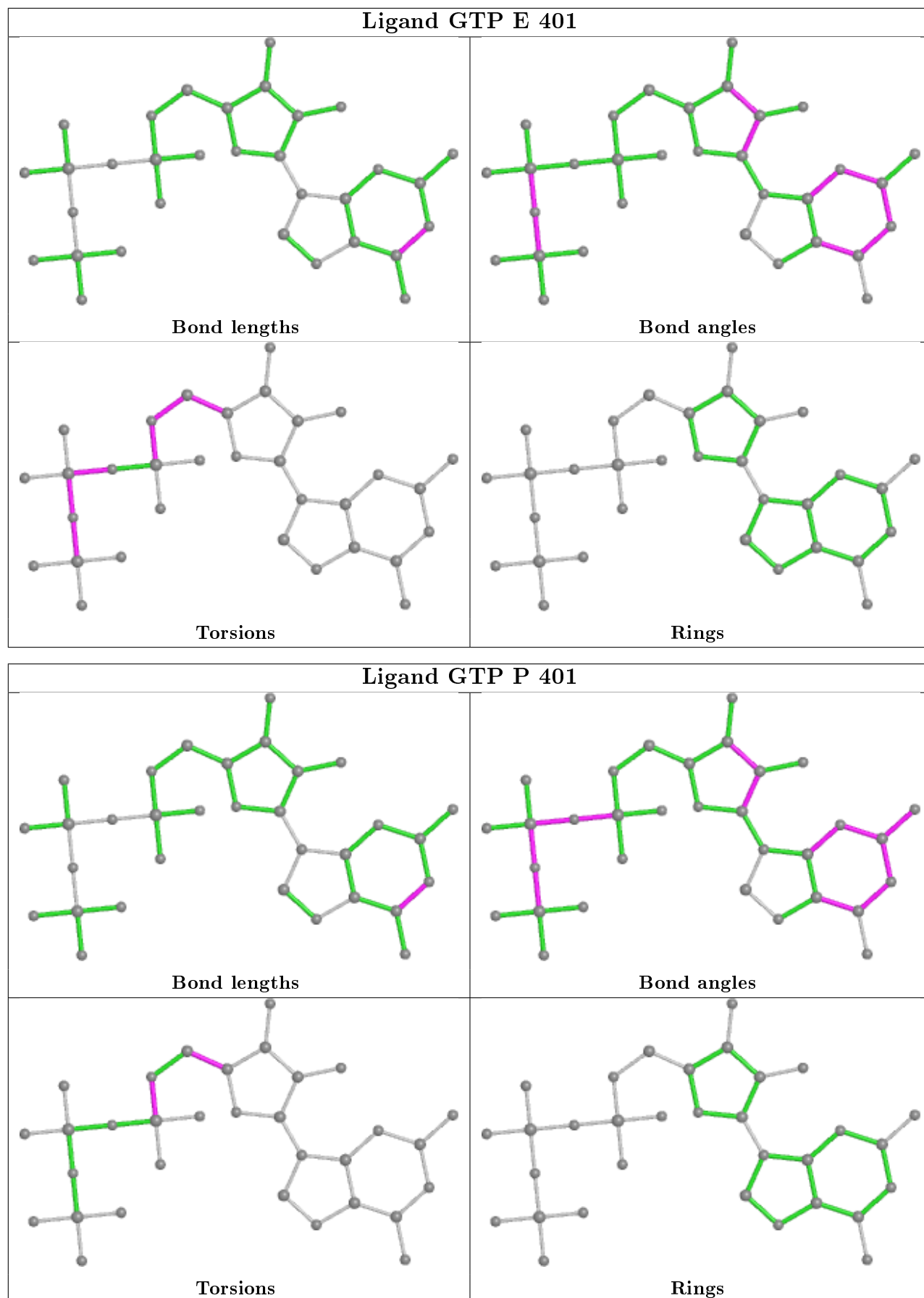


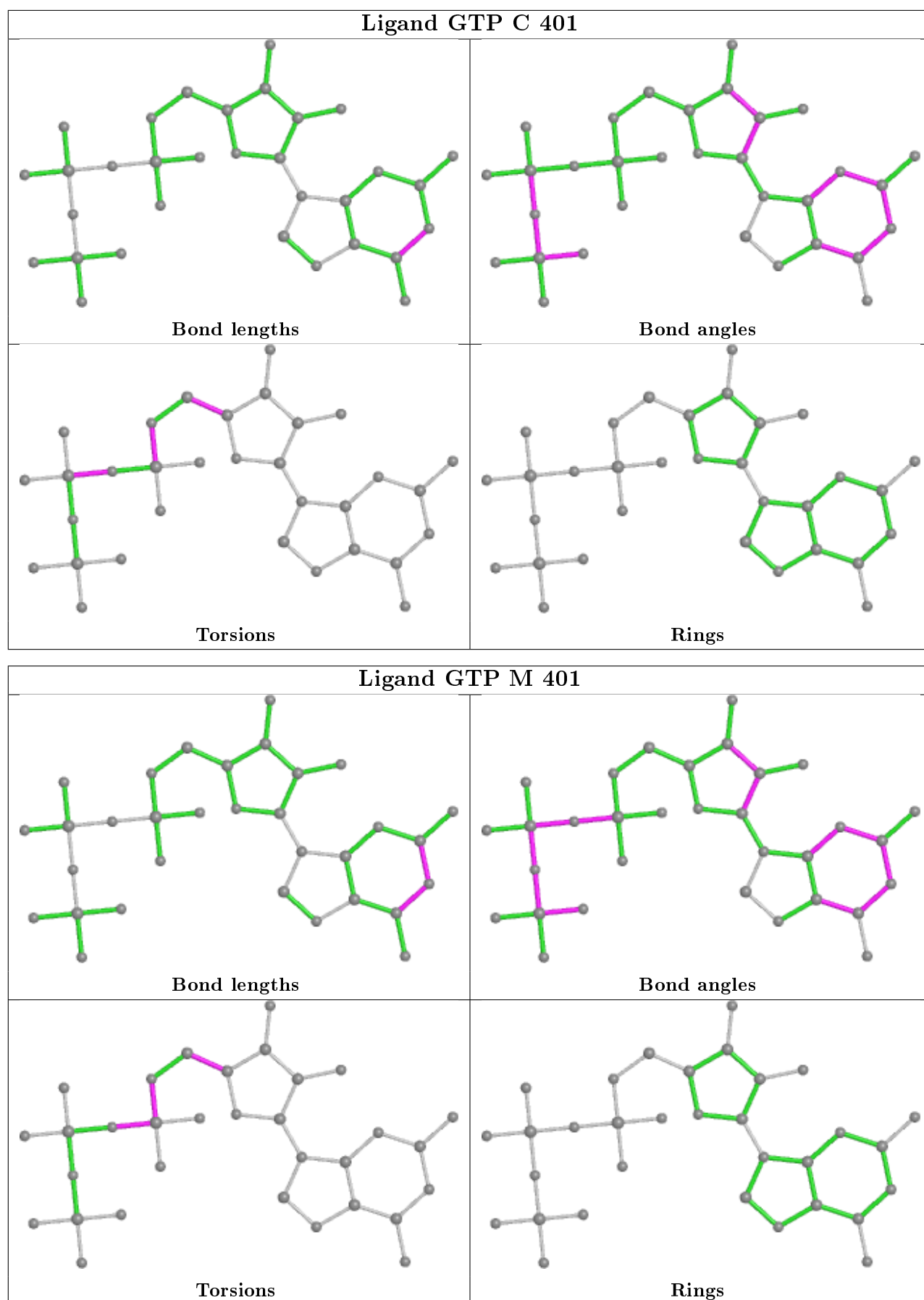


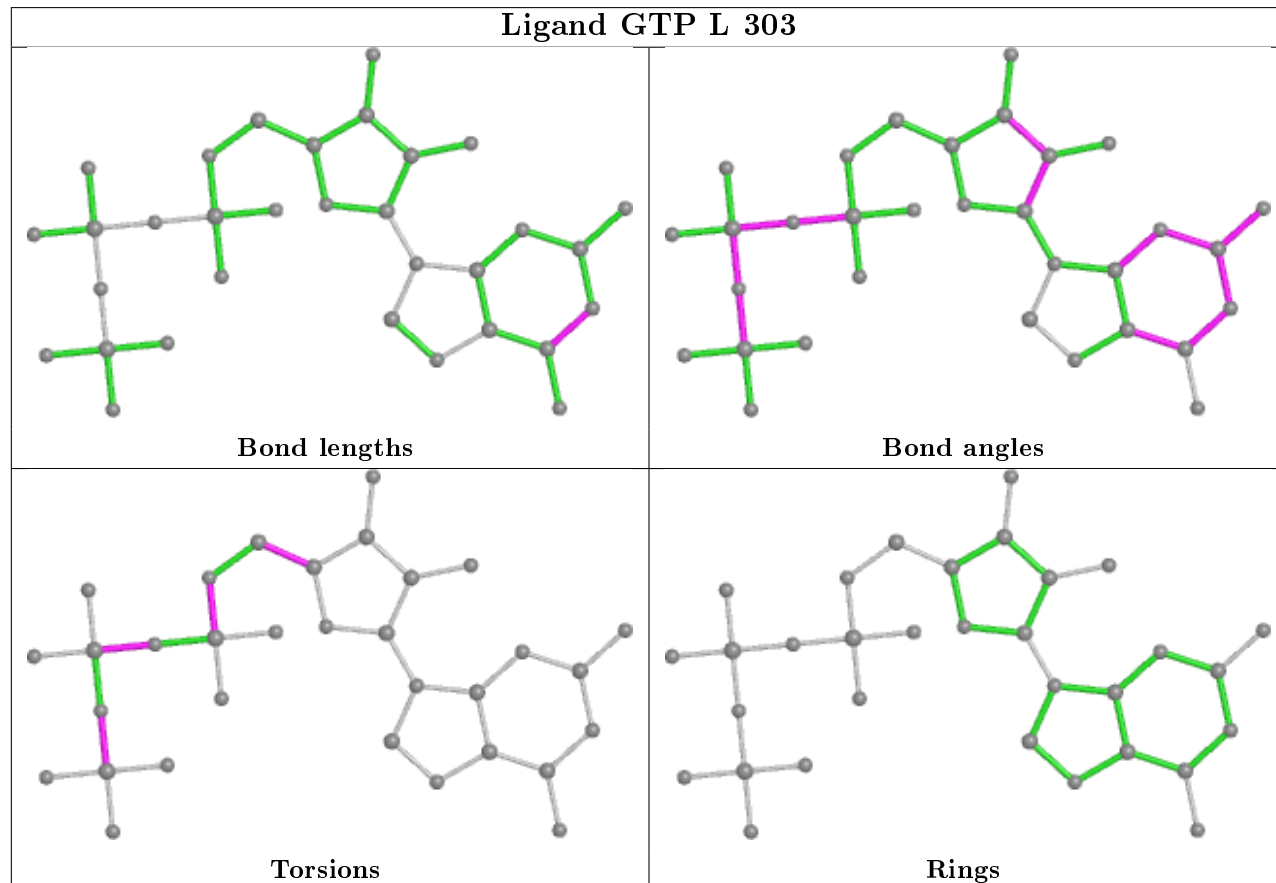
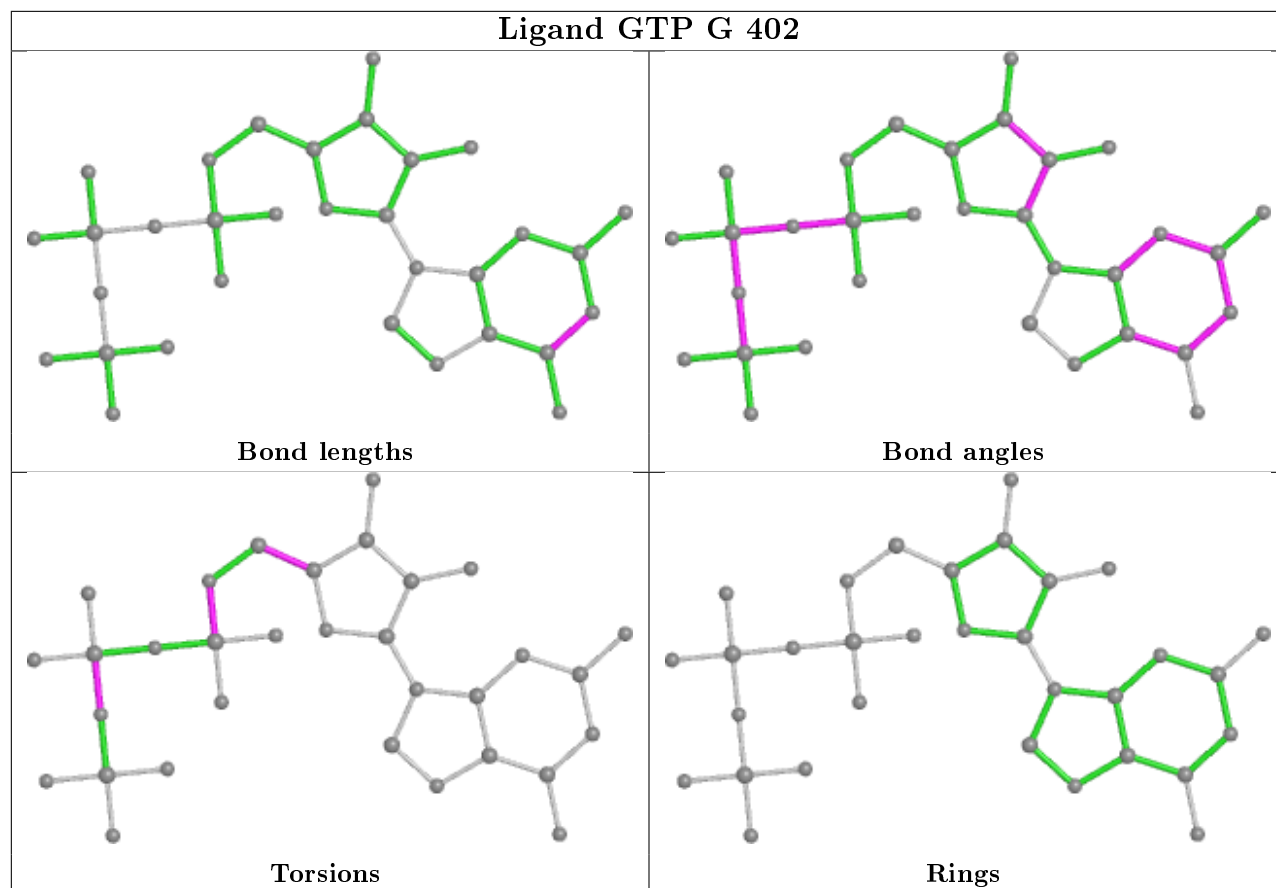


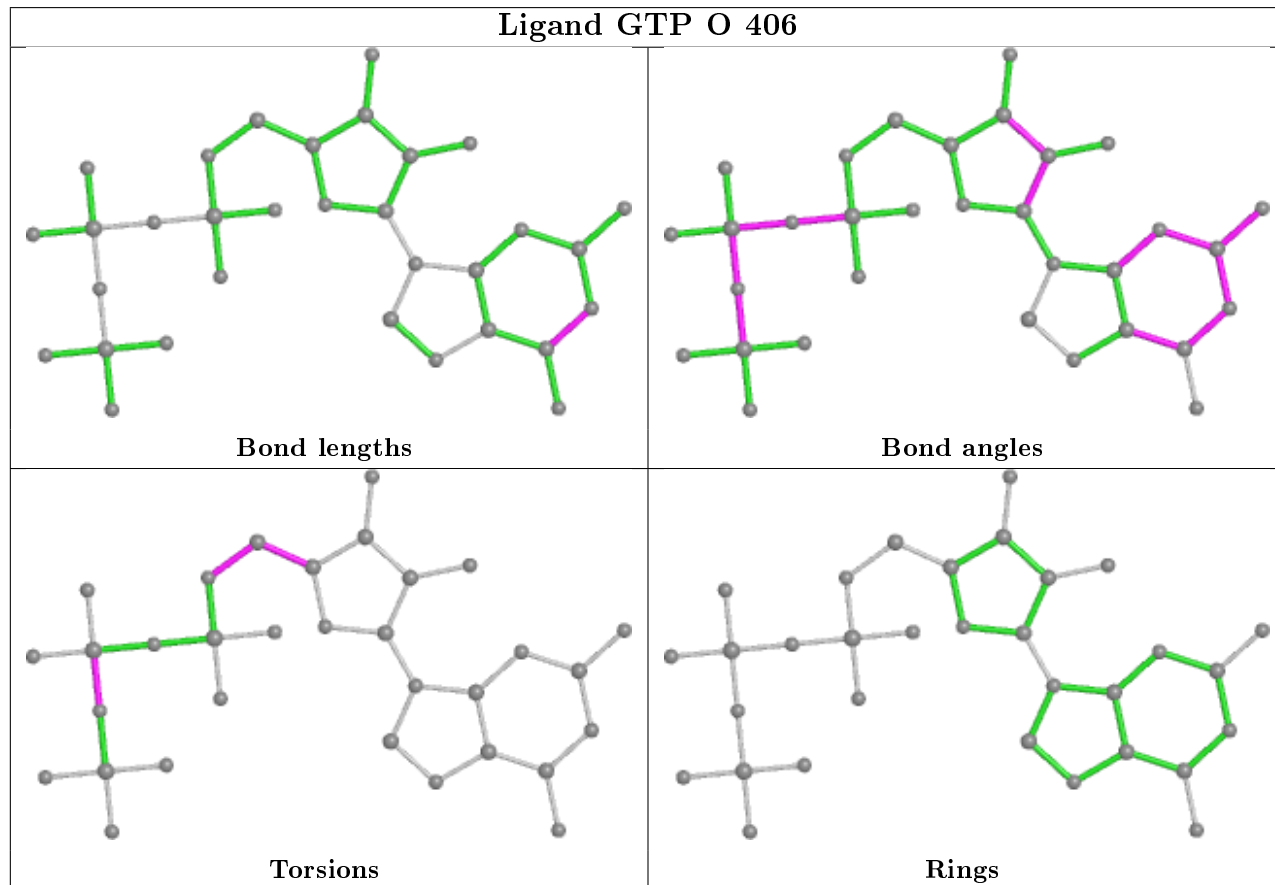
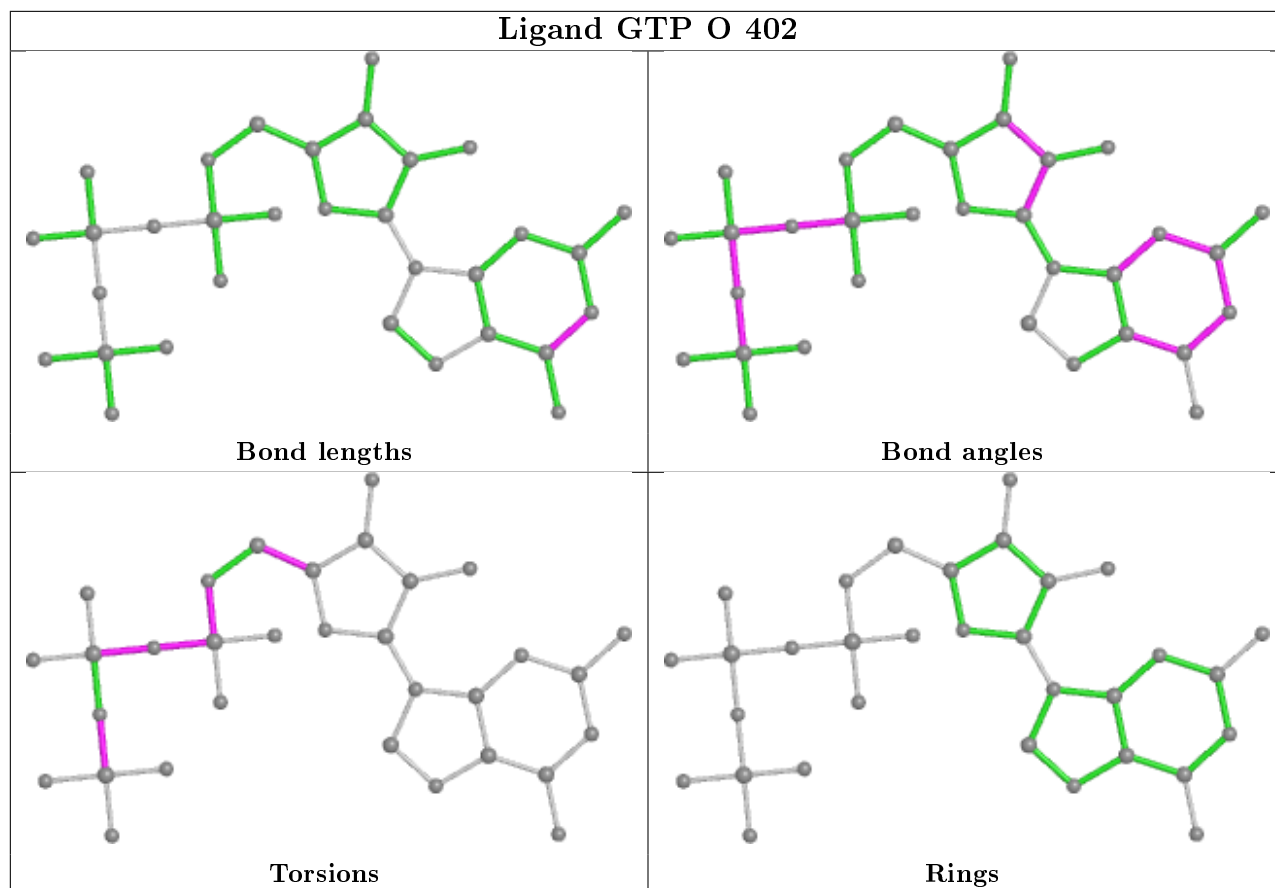


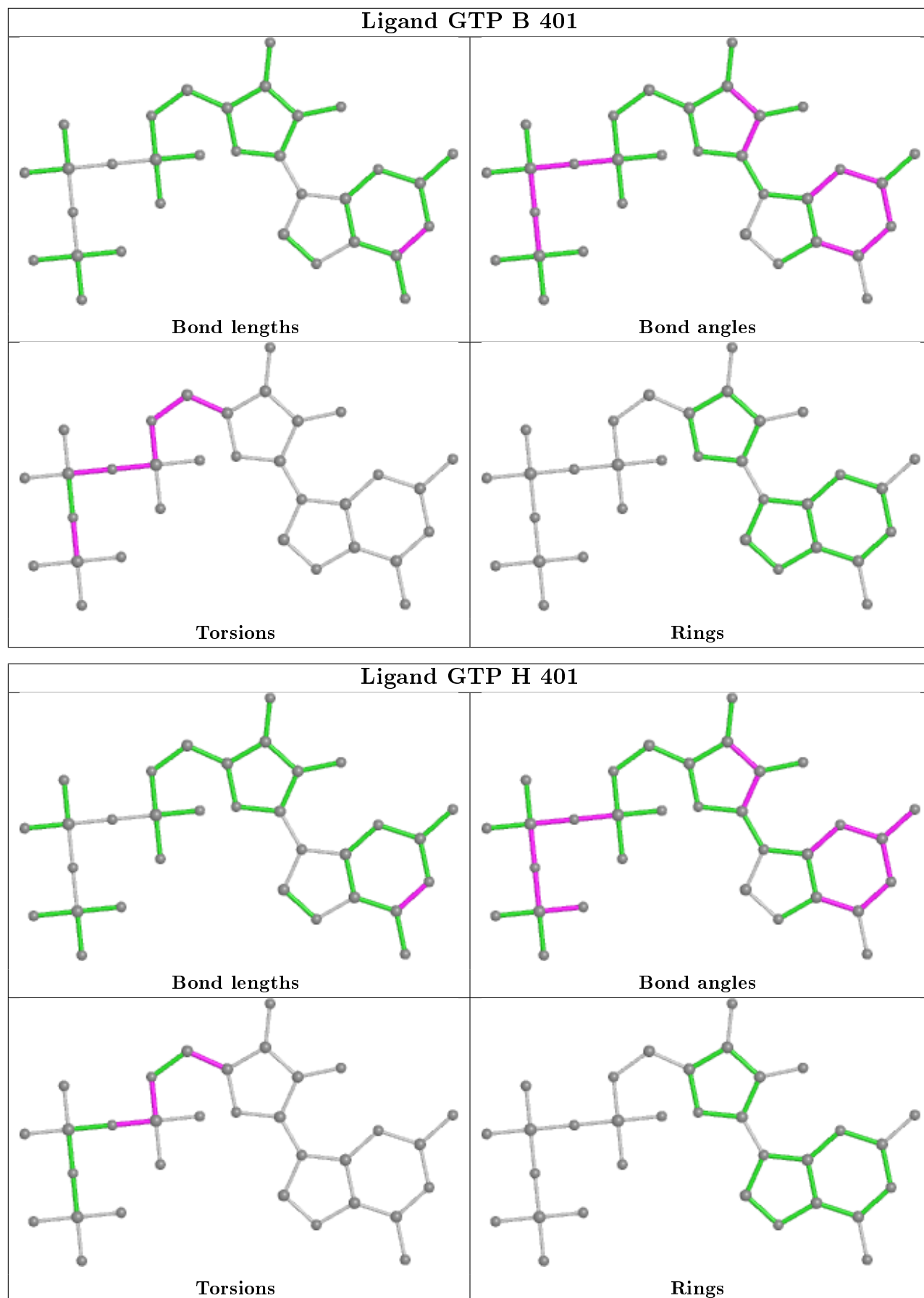












## 5.7 Other polymers

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	265/271 (97%)	0.27	17 (6%) 19 6	56, 90, 141, 156	0
1	B	265/271 (97%)	0.49	30 (11%) 5 1	50, 82, 139, 162	0
1	C	265/271 (97%)	0.09	6 (2%) 60 31	48, 92, 130, 138	0
1	D	265/271 (97%)	-0.14	2 (0%) 86 65	45, 69, 108, 128	0
1	E	241/271 (88%)	0.52	19 (7%) 12 4	44, 106, 147, 162	0
1	F	242/271 (89%)	0.01	1 (0%) 92 79	38, 67, 107, 126	0
1	G	265/271 (97%)	0.22	17 (6%) 19 6	38, 81, 139, 150	0
1	H	265/271 (97%)	0.05	14 (5%) 26 9	34, 65, 126, 146	0
1	I	242/271 (89%)	-0.01	3 (1%) 79 53	40, 72, 119, 135	0
1	J	265/271 (97%)	0.32	20 (7%) 14 4	41, 82, 141, 162	0
1	K	265/271 (97%)	0.20	9 (3%) 45 19	36, 83, 142, 168	0
1	L	242/271 (89%)	-0.09	2 (0%) 86 65	48, 78, 110, 135	0
1	M	265/271 (97%)	0.03	8 (3%) 50 22	31, 59, 122, 139	0
1	N	265/271 (97%)	0.09	11 (4%) 36 14	28, 58, 120, 144	0
1	O	265/271 (97%)	-0.10	4 (1%) 73 46	40, 60, 118, 148	0
1	P	242/271 (89%)	0.74	31 (12%) 3 1	40, 93, 179, 188	0
All	All	4124/4336 (95%)	0.17	194 (4%) 31 12	28, 77, 137, 188	0

The worst 5 of 194 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	N	234	VAL	8.9
1	P	167	VAL	7.7
1	O	225	ASP	7.6
1	G	225	ASP	7.3
1	P	168	LEU	6.6

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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
3	MG	K	402	1/1	0.46	0.45	100,100,100,100	0
3	MG	O	403	1/1	0.77	0.22	48,48,48,48	0
3	MG	L	305	1/1	0.80	0.26	67,67,67,67	0
3	MG	G	403	1/1	0.81	0.17	69,69,69,69	0
3	MG	N	403	1/1	0.81	0.17	51,51,51,51	0
3	MG	D	303	1/1	0.82	0.10	77,77,77,77	0
3	MG	D	304	1/1	0.83	0.14	46,46,46,46	0
3	MG	H	403	1/1	0.84	0.16	67,67,67,67	0
3	MG	O	405	1/1	0.87	0.23	25,25,25,25	0
3	MG	C	402	1/1	0.87	0.10	82,82,82,82	0
2	GTP	F	301	32/32	0.87	0.30	64,129,135,137	0
2	GTP	L	301	32/32	0.88	0.21	123,138,167,192	0
3	MG	O	404	1/1	0.88	0.20	27,27,27,27	0
3	MG	B	403	1/1	0.88	0.16	71,71,71,71	0
3	MG	H	404	1/1	0.88	0.12	29,29,29,29	0
3	MG	K	403	1/1	0.88	0.13	220,220,220,220	0
3	MG	P	402	1/1	0.89	0.10	65,65,65,65	0
3	MG	F	304	1/1	0.89	0.26	60,60,60,60	0
2	GTP	K	401	32/32	0.89	0.21	129,146,177,177	0
3	MG	E	403	1/1	0.90	0.19	68,68,68,68	0
3	MG	M	403	1/1	0.90	0.10	54,54,54,54	0
3	MG	J	403	1/1	0.90	0.19	54,54,54,54	0
3	MG	C	404	1/1	0.91	0.16	54,54,54,54	0
2	GTP	E	401	32/32	0.91	0.19	62,102,120,156	0
2	GTP	P	401	32/32	0.91	0.21	58,101,113,116	0
2	GTP	C	405	32/32	0.92	0.22	40,103,115,121	0
2	GTP	O	406	32/32	0.92	0.21	39,129,133,134	0

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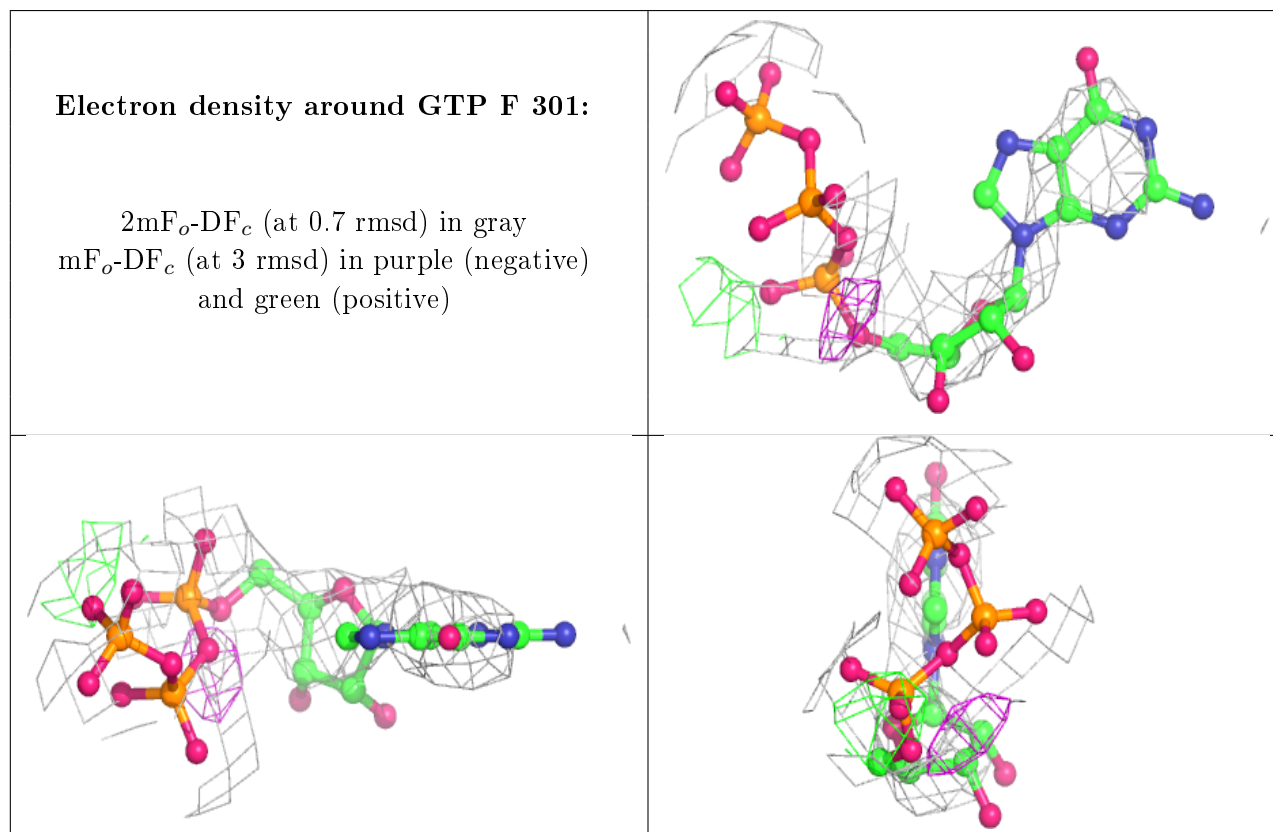
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
3	MG	F	305	1/1	0.92	0.19	46,46,46,46	0
3	MG	I	405	1/1	0.92	0.24	34,34,34,34	0
2	GTP	C	401	32/32	0.93	0.23	64,104,111,111	0
3	MG	A	403	1/1	0.93	0.13	73,73,73,73	0
3	MG	J	405	1/1	0.93	0.29	43,43,43,43	0
2	GTP	B	401	32/32	0.93	0.18	63,86,95,115	0
3	MG	L	302	1/1	0.93	0.15	74,74,74,74	0
3	MG	P	403	1/1	0.93	0.21	42,42,42,42	0
3	MG	M	404	1/1	0.94	0.08	16,16,16,16	0
3	MG	F	302	1/1	0.94	0.37	61,61,61,61	0
2	GTP	A	402	32/32	0.94	0.19	43,122,128,129	0
2	GTP	L	303	32/32	0.94	0.19	39,69,78,80	0
2	GTP	H	402	32/32	0.94	0.23	26,92,99,101	0
3	MG	H	405	1/1	0.94	0.26	31,31,31,31	0
2	GTP	O	402	32/32	0.94	0.24	22,113,118,118	0
2	GTP	D	301	32/32	0.94	0.20	70,113,118,119	0
3	MG	G	404	1/1	0.94	0.19	39,39,39,39	0
2	GTP	A	401	32/32	0.94	0.15	60,81,91,96	0
3	MG	F	306	1/1	0.94	0.33	45,45,45,45	0
2	GTP	H	401	32/32	0.94	0.17	30,57,66,66	0
2	GTP	O	401	32/32	0.94	0.21	21,57,70,85	0
3	MG	L	306	1/1	0.94	0.17	59,59,59,59	0
2	GTP	J	401	32/32	0.95	0.19	43,69,74,75	0
2	GTP	I	402	32/32	0.95	0.18	39,94,103,104	0
3	MG	L	307	1/1	0.95	0.16	27,27,27,27	0
3	MG	A	405	1/1	0.95	0.21	45,45,45,45	0
2	GTP	E	404	32/32	0.95	0.21	20,101,105,107	0
2	GTP	G	402	32/32	0.95	0.21	44,143,155,155	0
3	MG	I	403	1/1	0.95	0.25	54,54,54,54	0
2	GTP	B	402	32/32	0.95	0.18	59,123,126,126	0
2	GTP	F	303	32/32	0.95	0.21	37,76,89,95	0
3	MG	N	405	1/1	0.96	0.29	23,23,23,23	0
2	GTP	M	402	32/32	0.96	0.17	31,95,104,106	0
2	GTP	L	304	32/32	0.96	0.19	28,120,123,124	0
3	MG	D	305	1/1	0.96	0.41	53,53,53,53	0
3	MG	G	405	1/1	0.96	0.32	50,50,50,50	0
2	GTP	I	401	32/32	0.96	0.18	10,93,102,127	0
2	GTP	D	302	32/32	0.96	0.17	25,66,71,73	0
2	GTP	G	401	32/32	0.96	0.23	45,84,90,92	0
3	MG	J	404	1/1	0.96	0.15	47,47,47,47	0
2	GTP	M	401	32/32	0.96	0.16	25,55,68,72	0
3	MG	B	404	1/1	0.96	0.12	42,42,42,42	0

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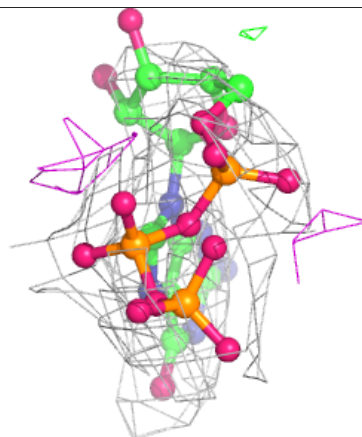
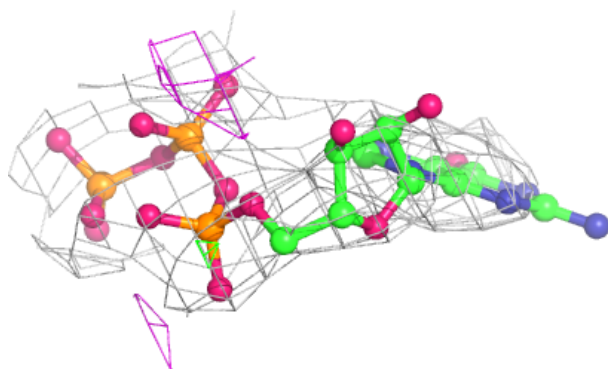
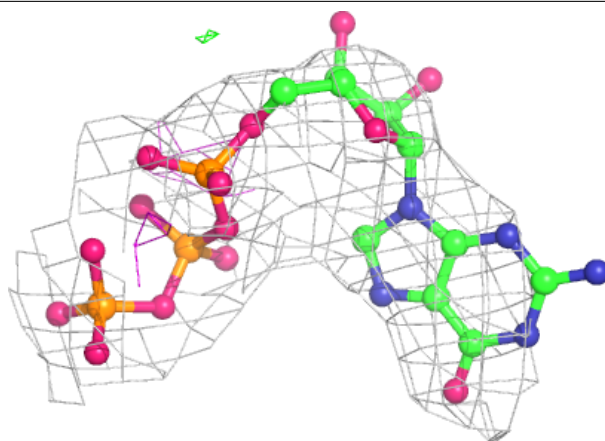
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	GTP	N	402	32/32	0.96	0.18	18,114,121,123	0
3	MG	M	405	1/1	0.97	0.17	21,21,21,21	0
2	GTP	N	401	32/32	0.97	0.17	0,63,68,70	0
3	MG	P	404	1/1	0.97	0.26	42,42,42,42	0
2	GTP	J	402	32/32	0.97	0.19	37,100,121,122	0
3	MG	E	402	1/1	0.98	0.24	91,91,91,91	0
3	MG	A	404	1/1	0.98	0.15	39,39,39,39	0
3	MG	B	405	1/1	0.98	0.23	56,56,56,56	0
3	MG	N	404	1/1	0.99	0.16	38,38,38,38	0
3	MG	I	404	1/1	0.99	0.16	61,61,61,61	0
3	MG	C	403	1/1	0.99	0.07	53,53,53,53	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.

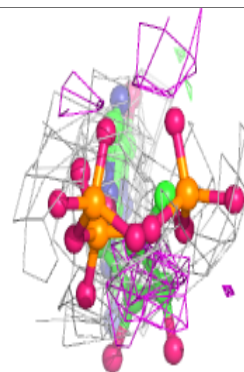
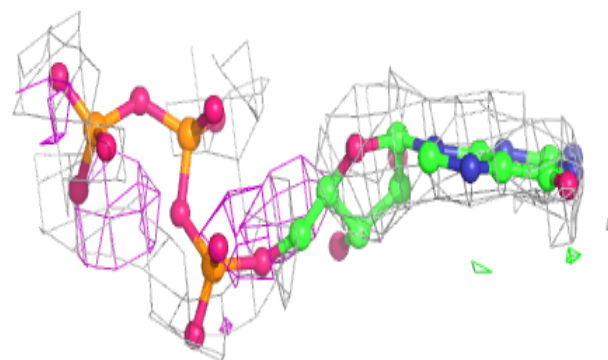
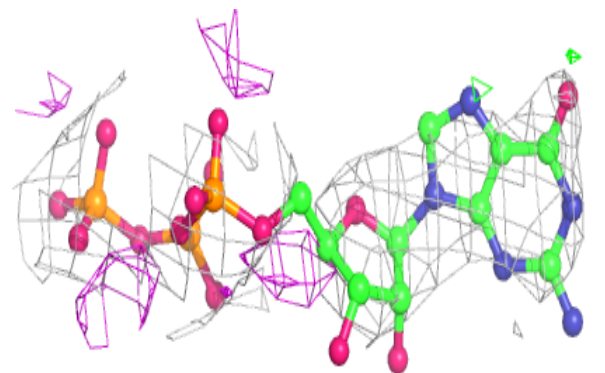


**Electron density around GTP L 301:**

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

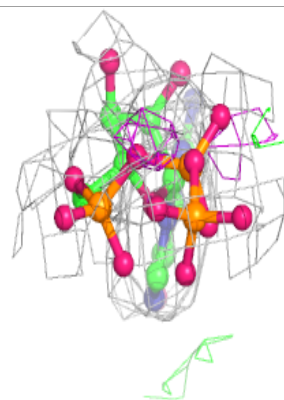
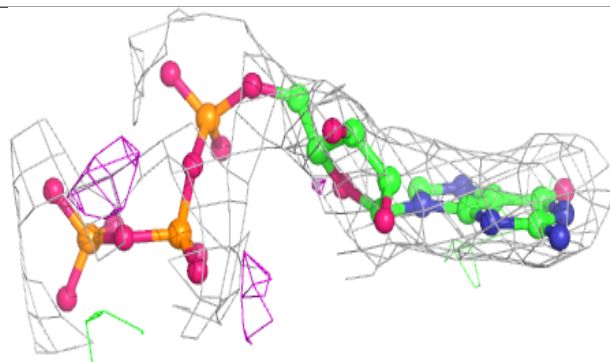
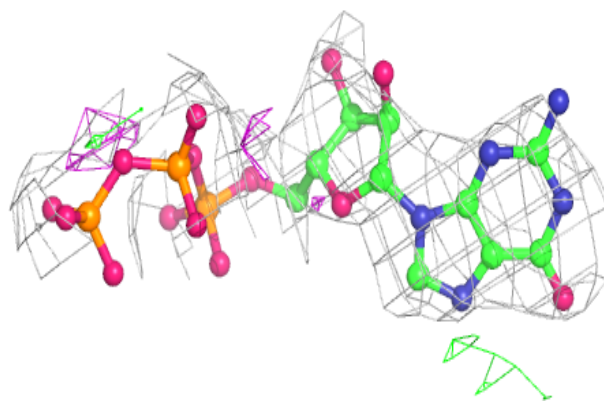
**Electron density around GTP K 401:**

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

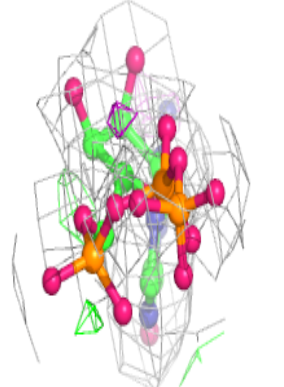
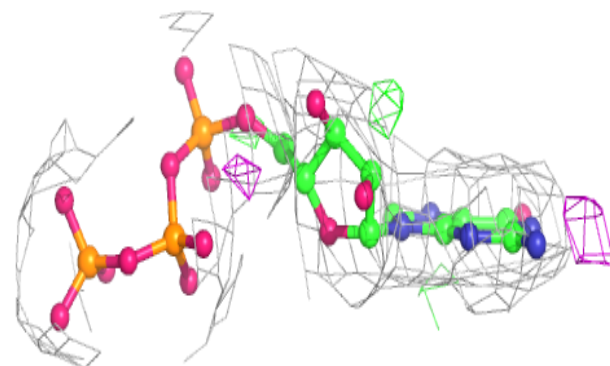
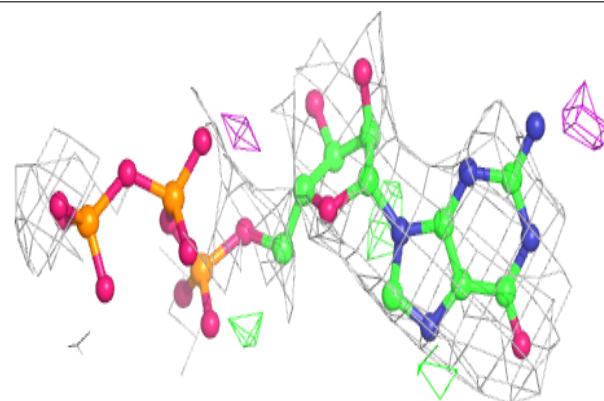


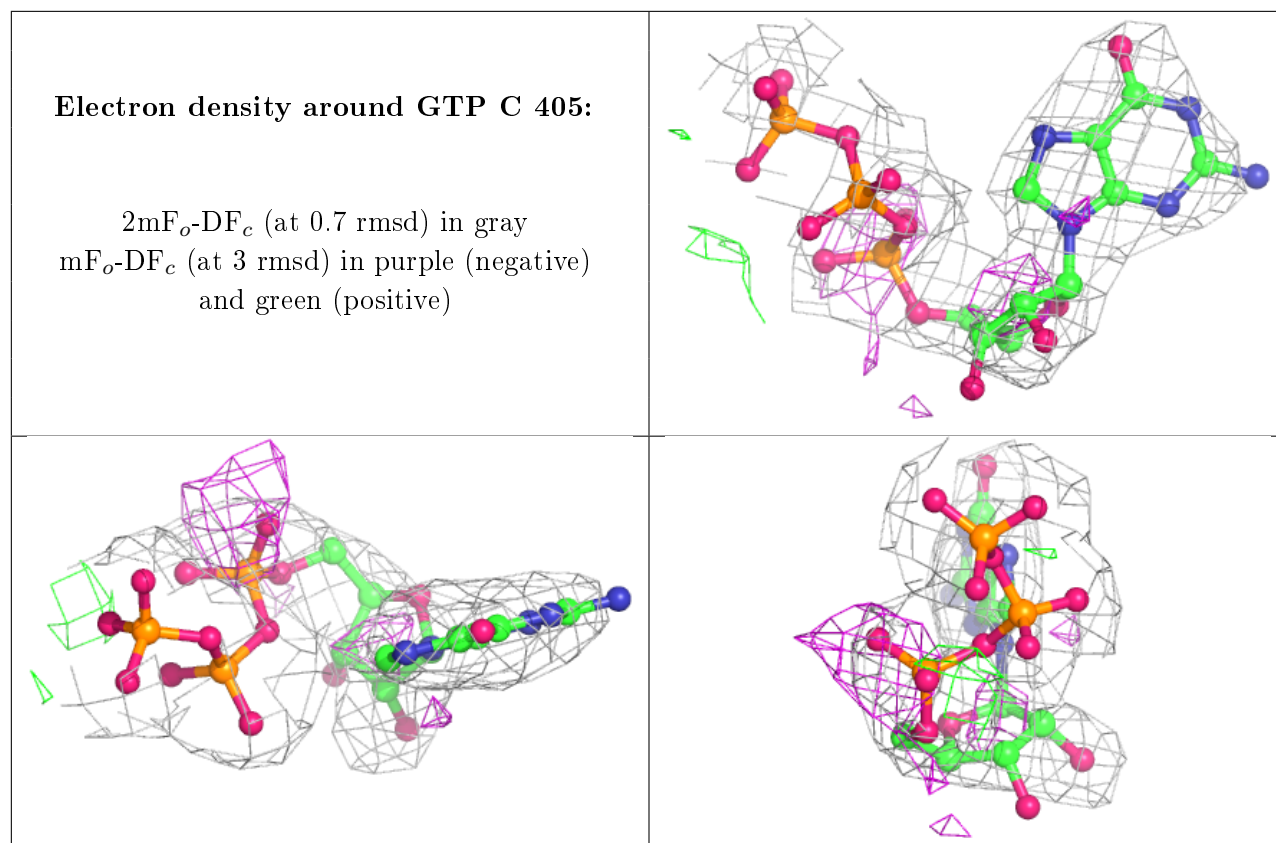
**Electron density around GTP E 401:**

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

**Electron density around GTP P 401:**

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

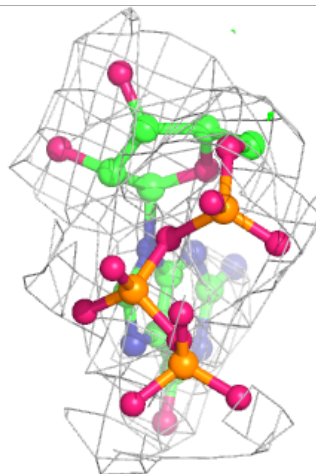
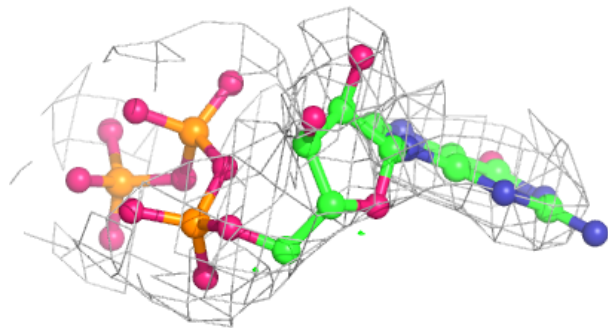
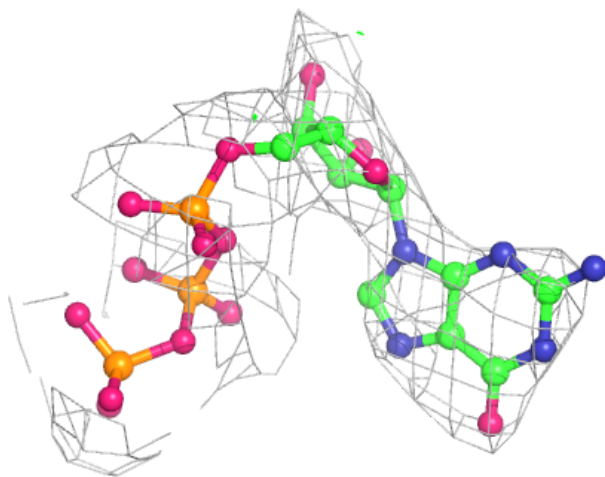




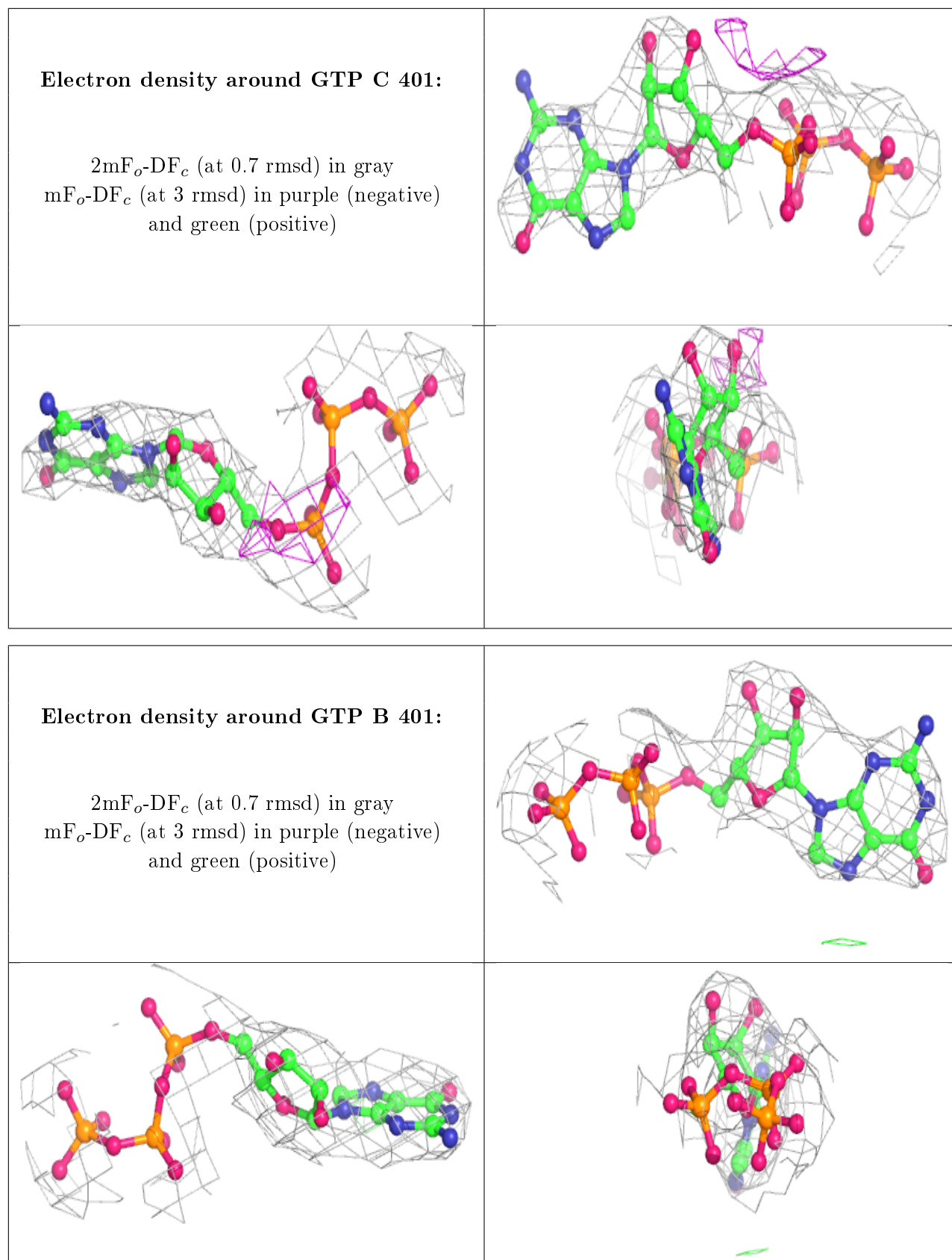


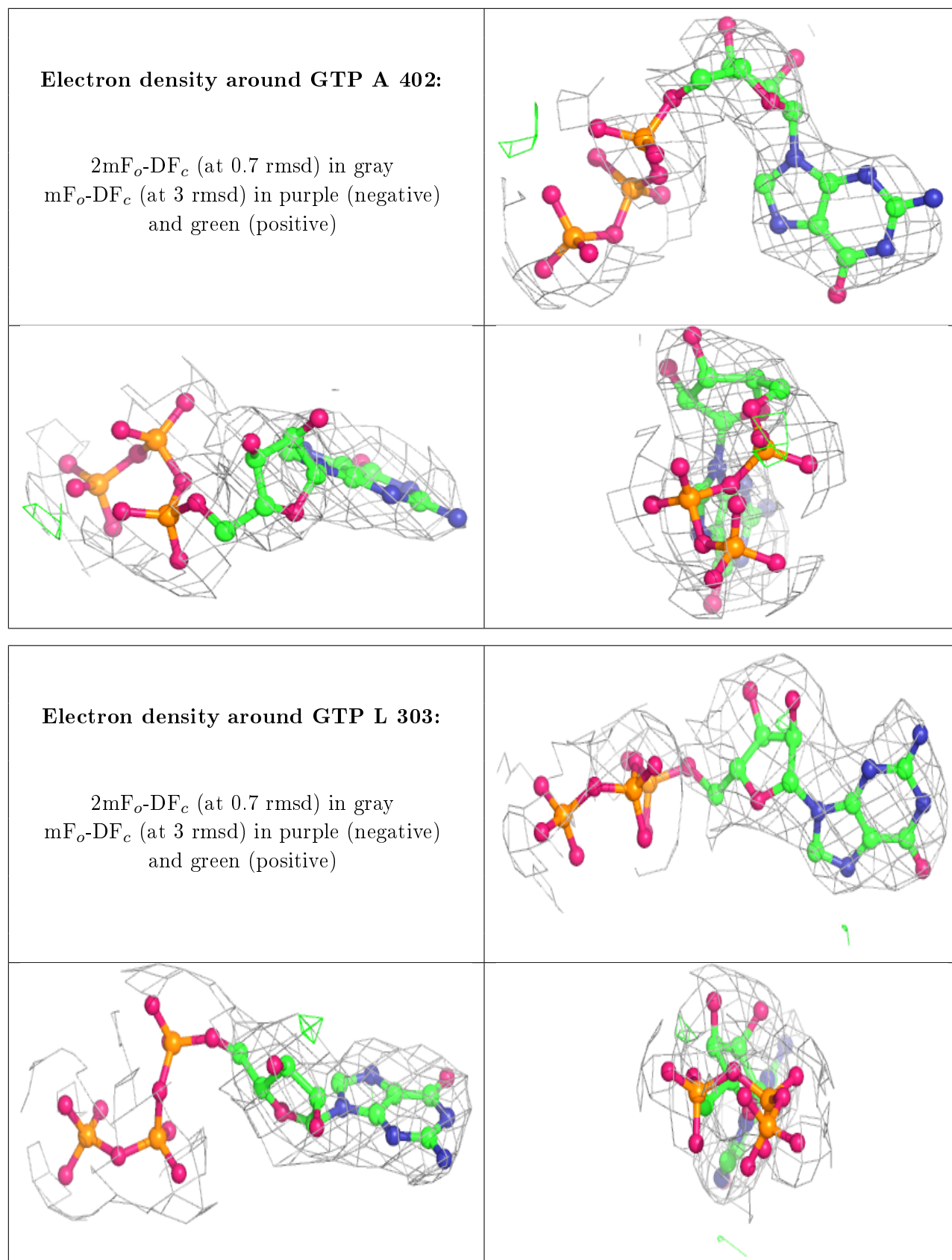
**Electron density around GTP O 406:**

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and green (positive)



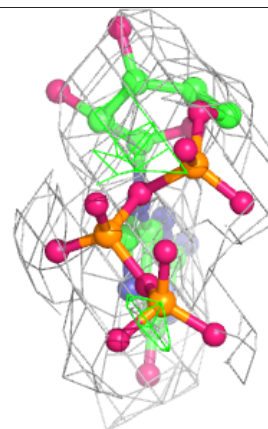
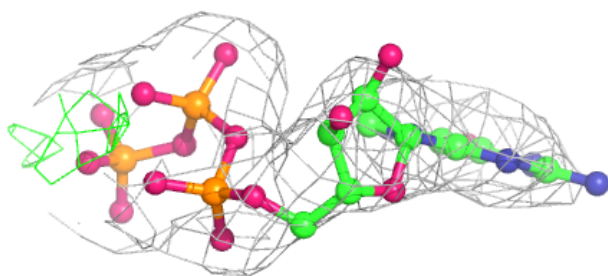
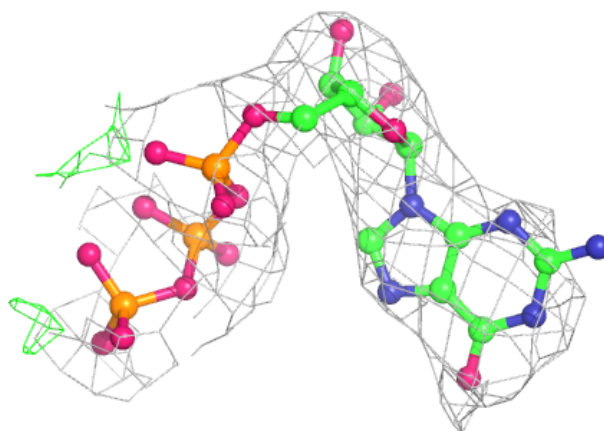


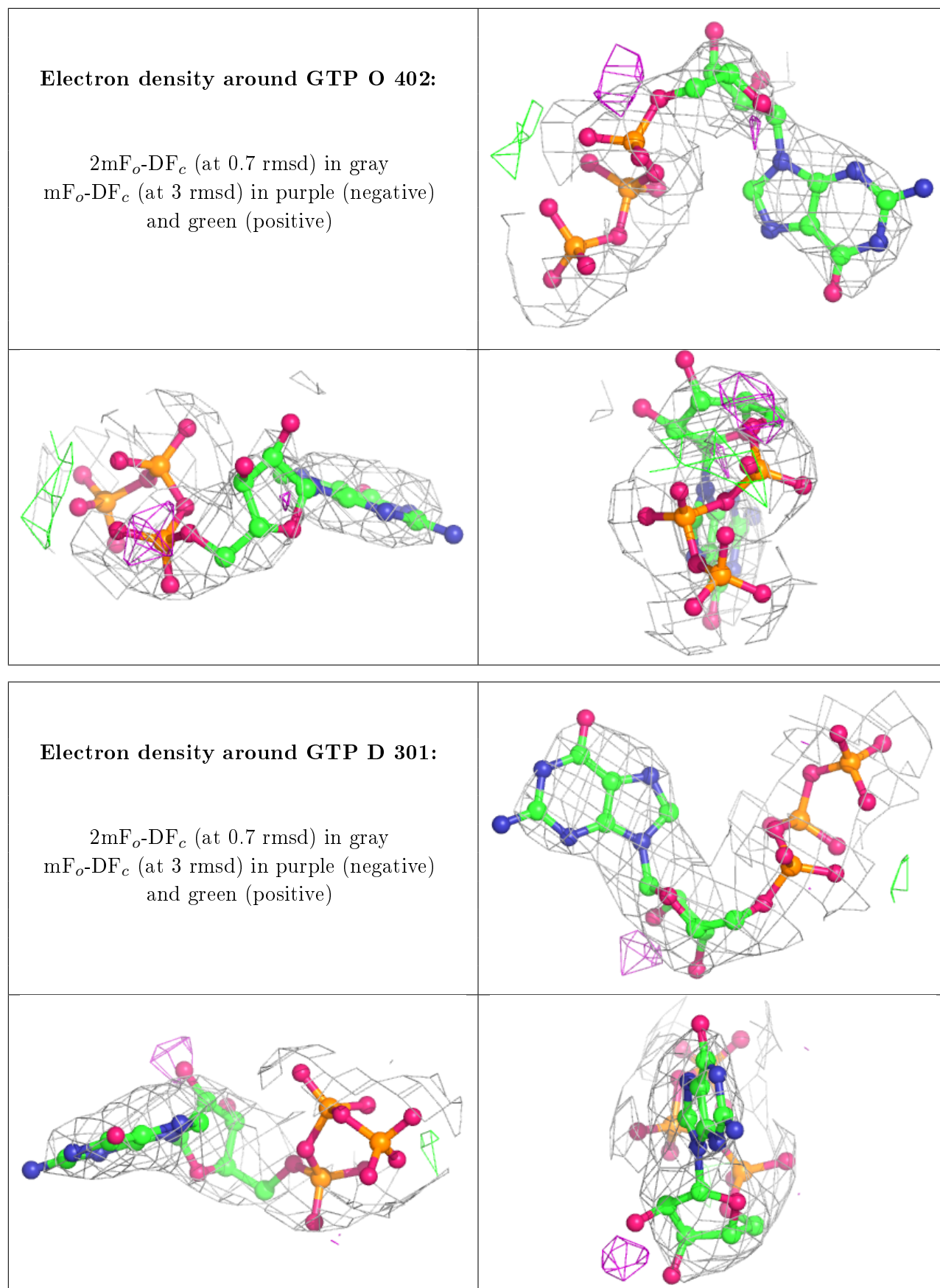


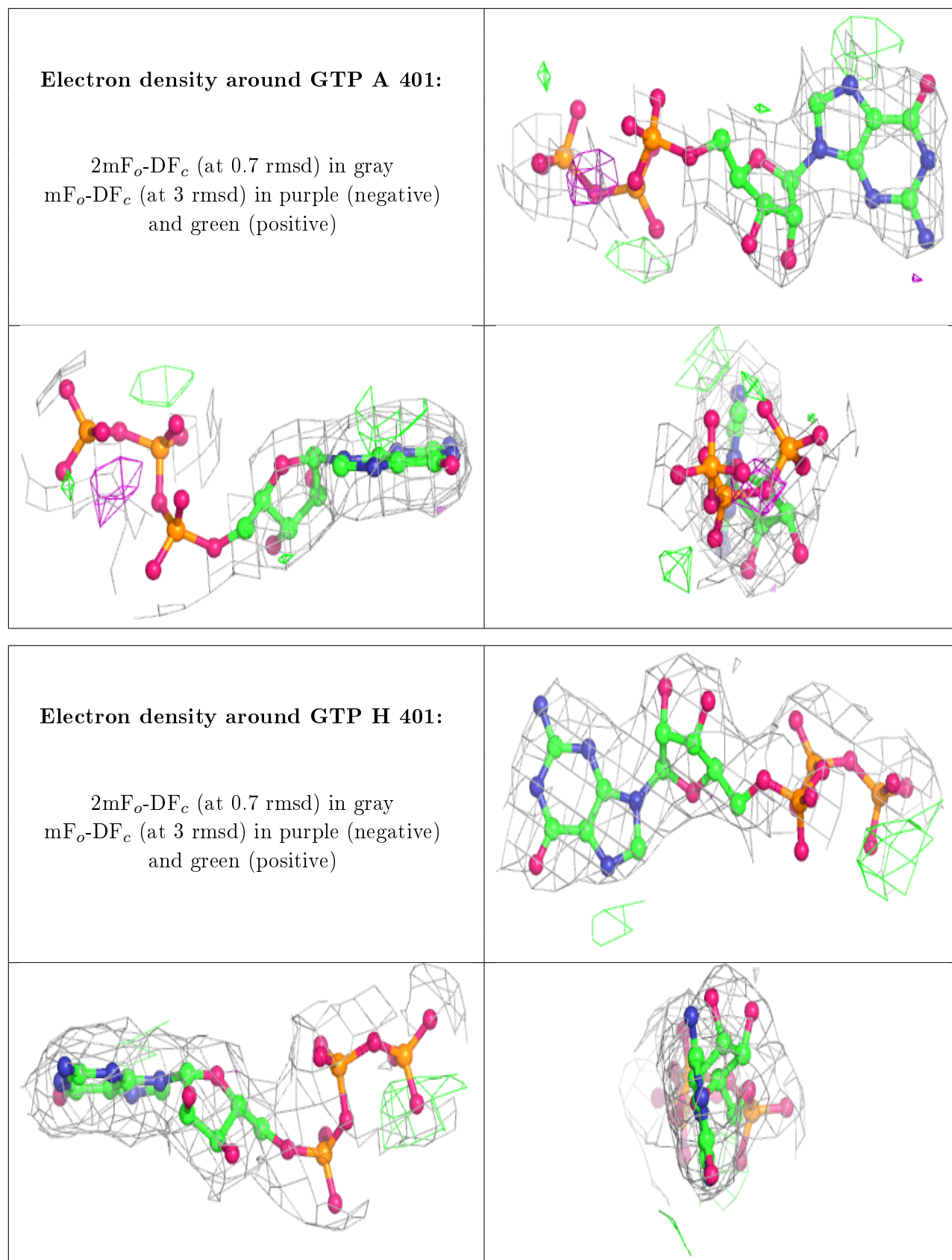


**Electron density around GTP H 402:**

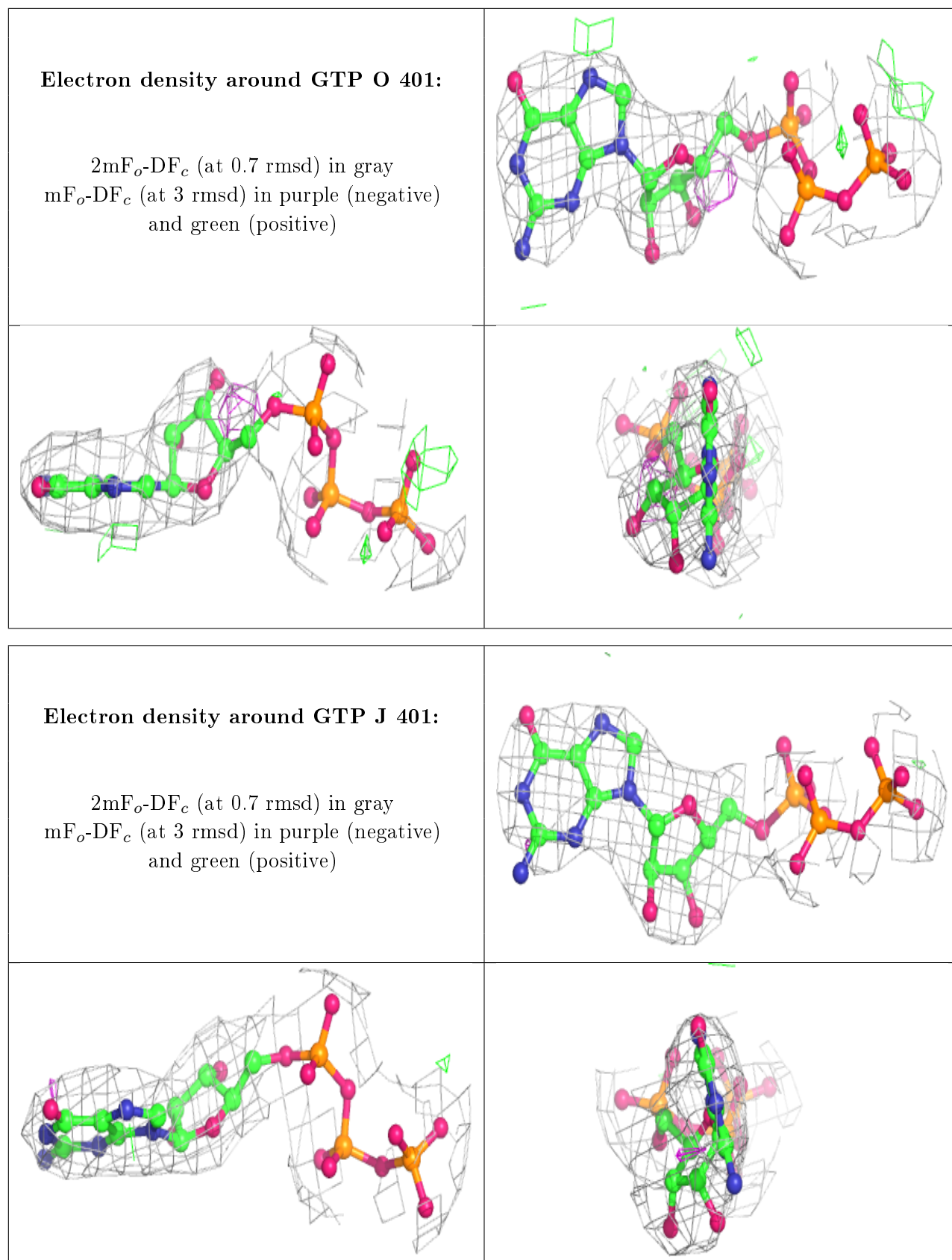
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





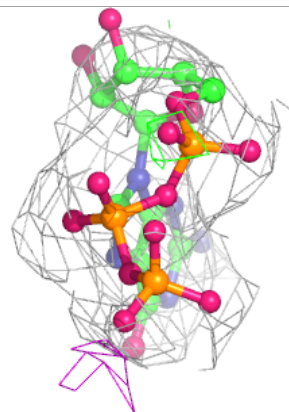
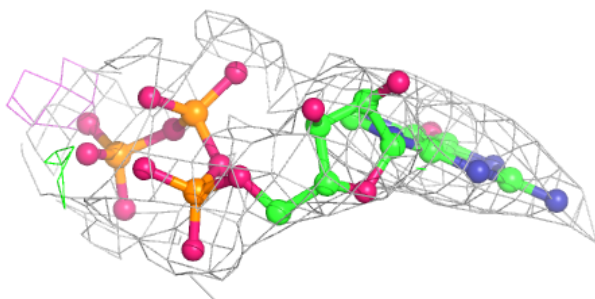
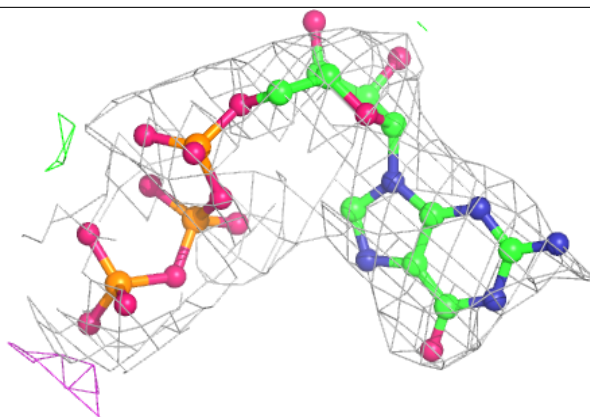




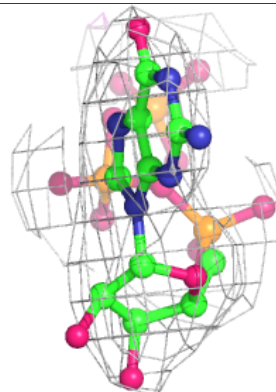
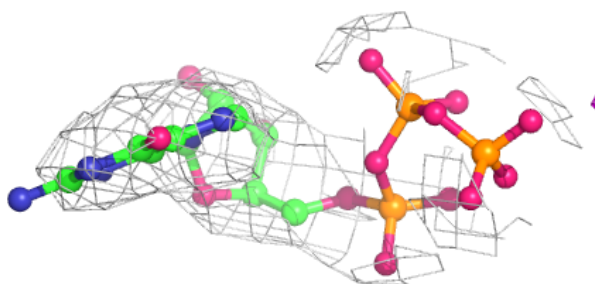
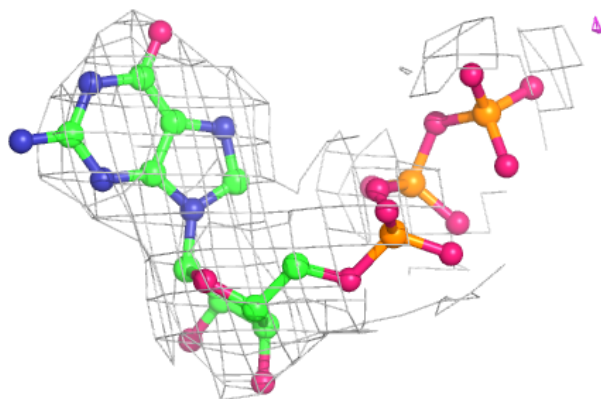


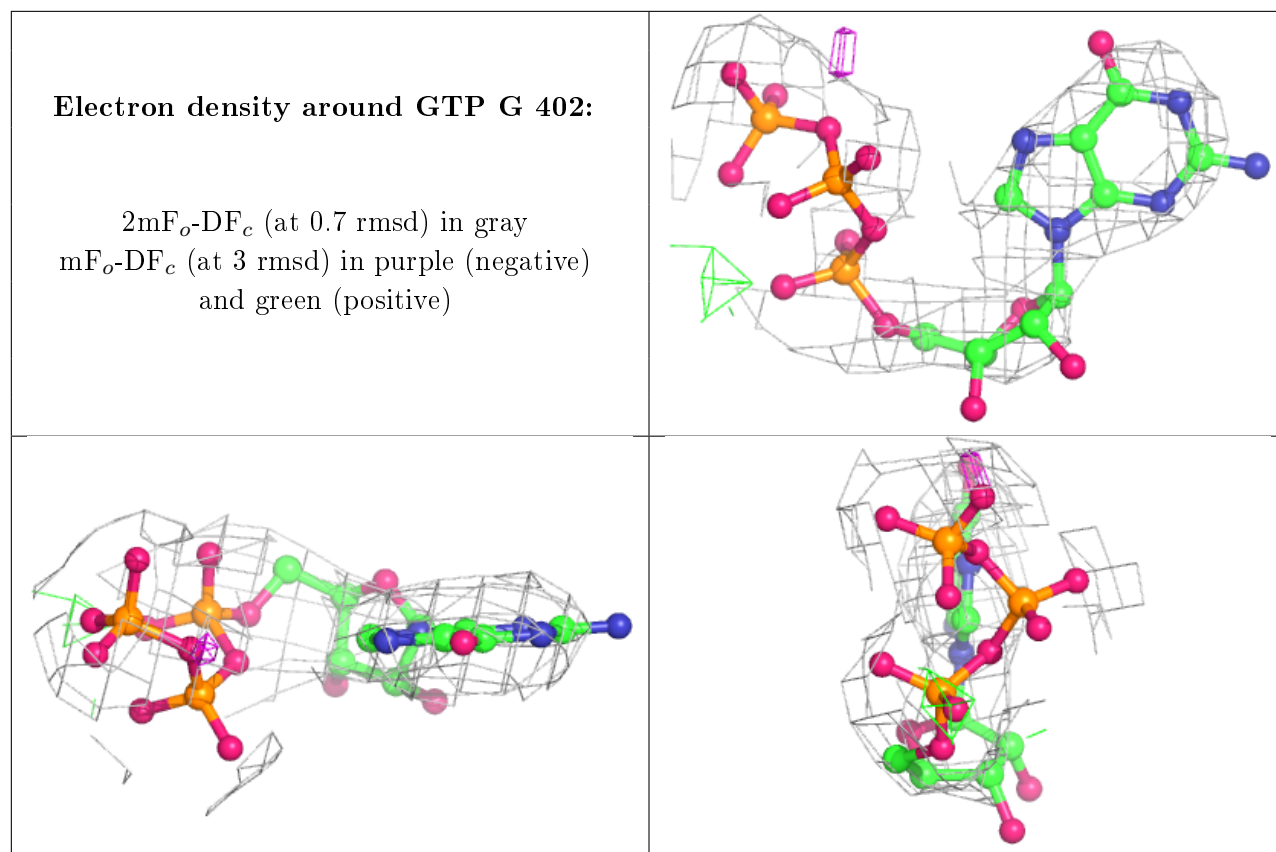
**Electron density around GTP I 402:**

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

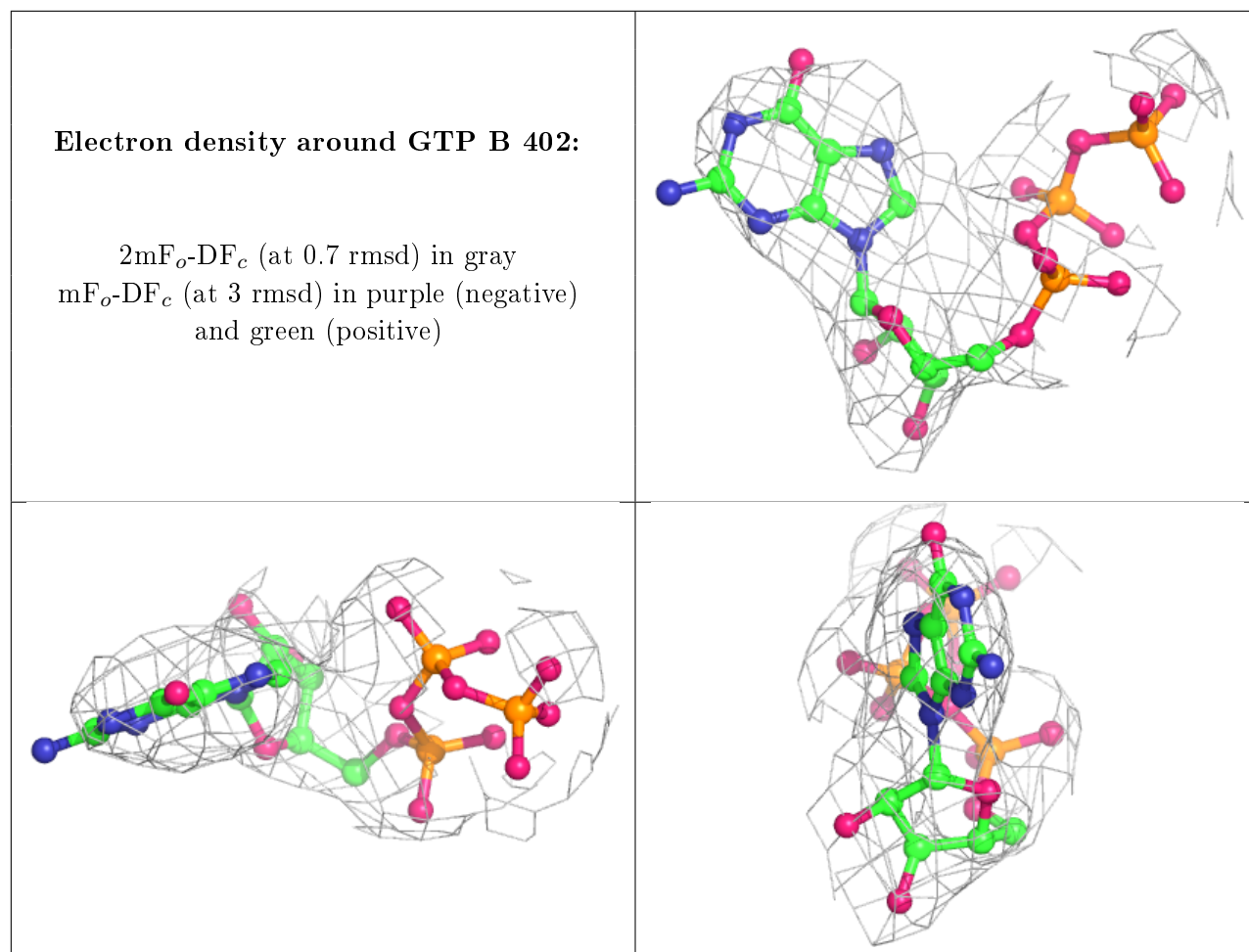
**Electron density around GTP E 404:**

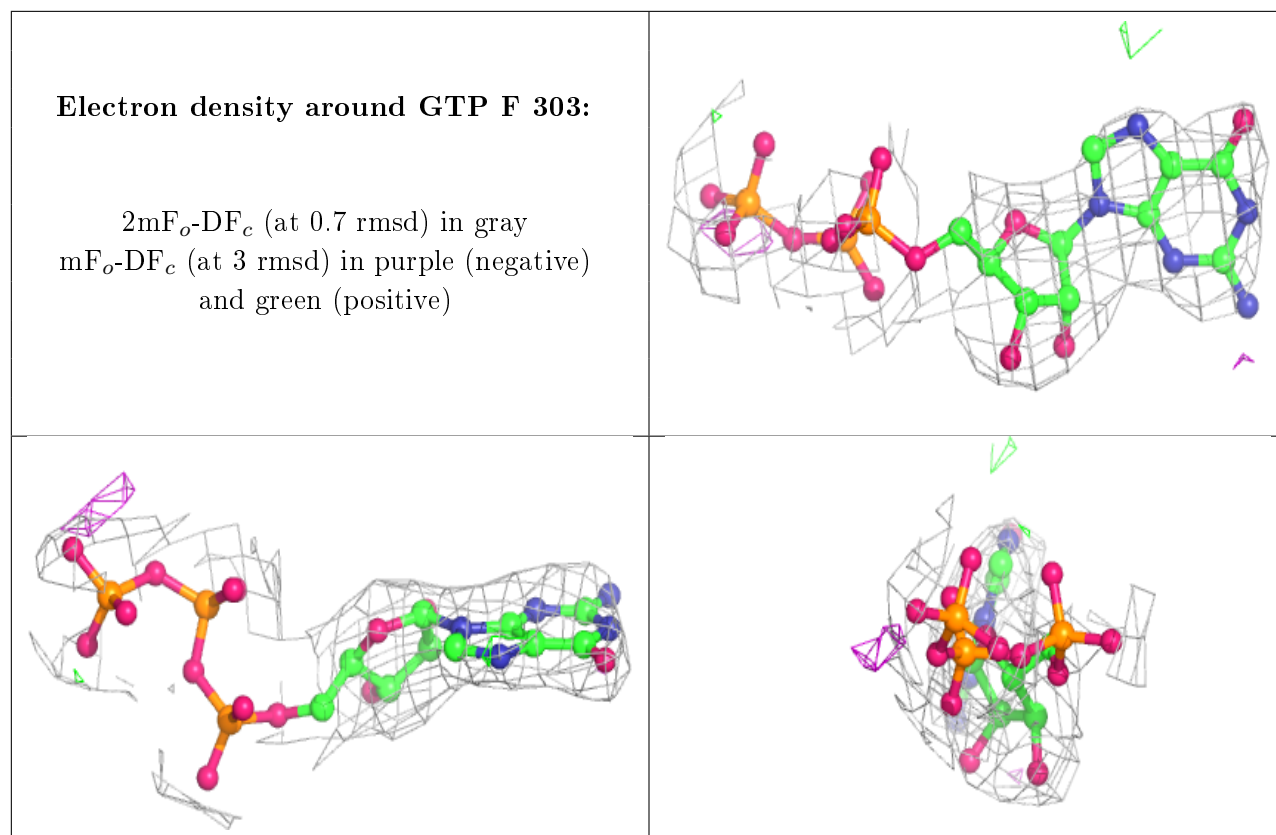
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





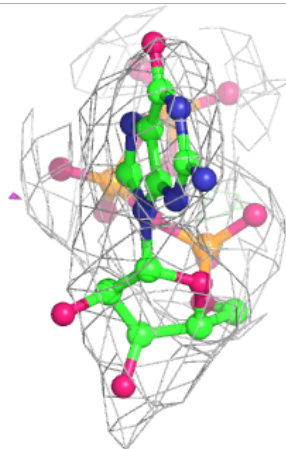
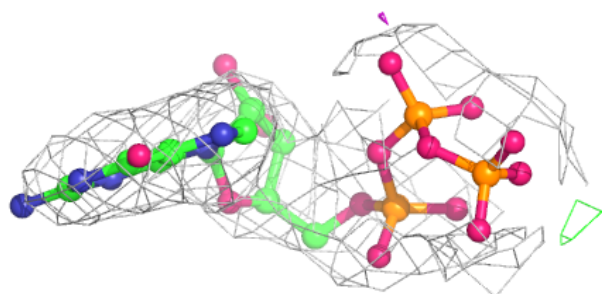
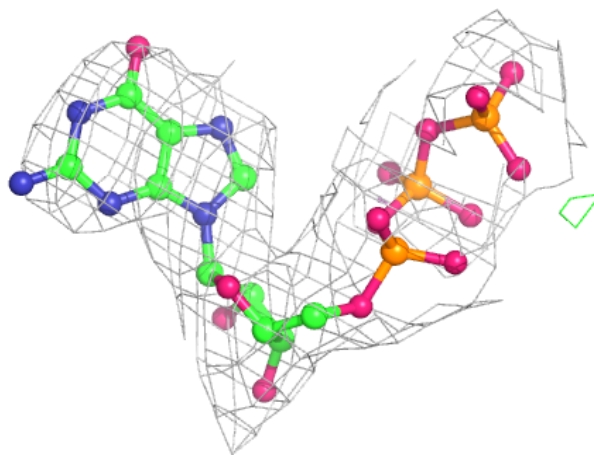






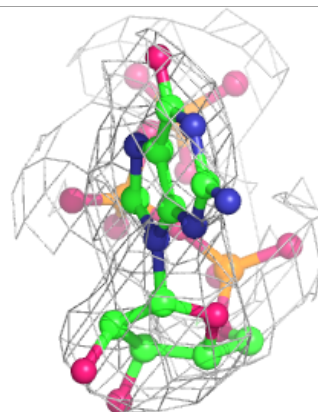
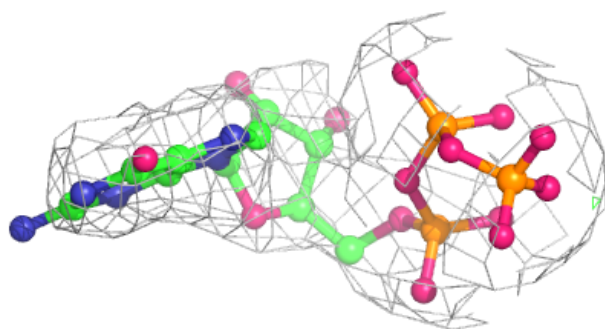
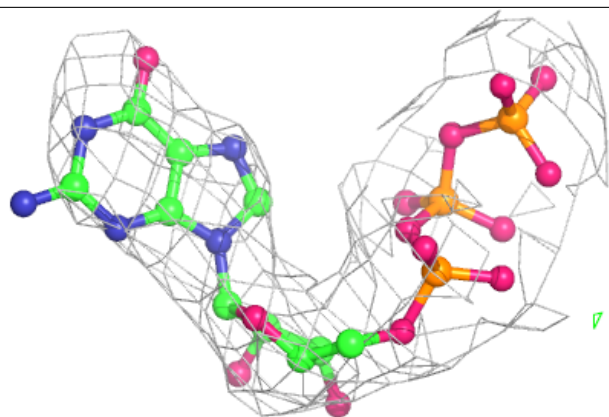
**Electron density around GTP M 402:**

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

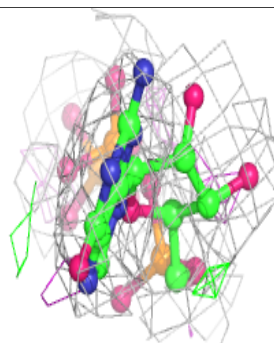
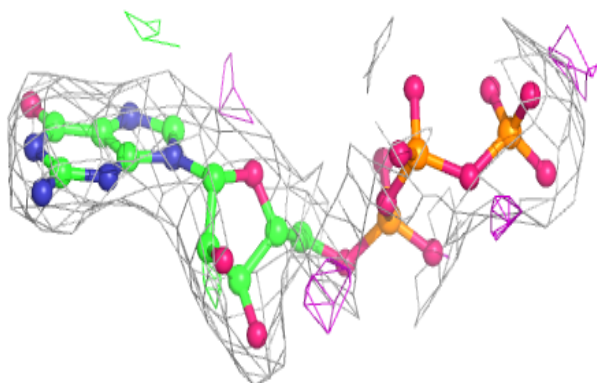
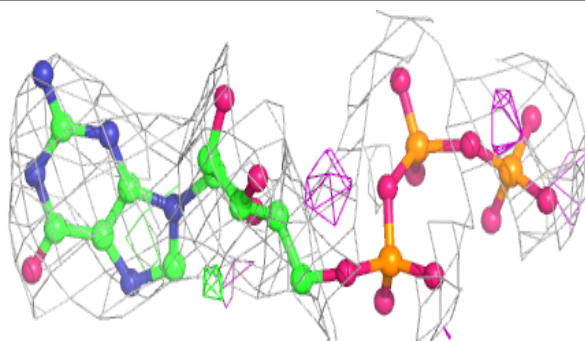


**Electron density around GTP L 304:**

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

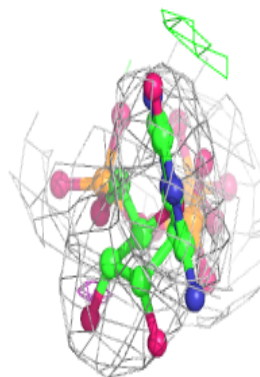
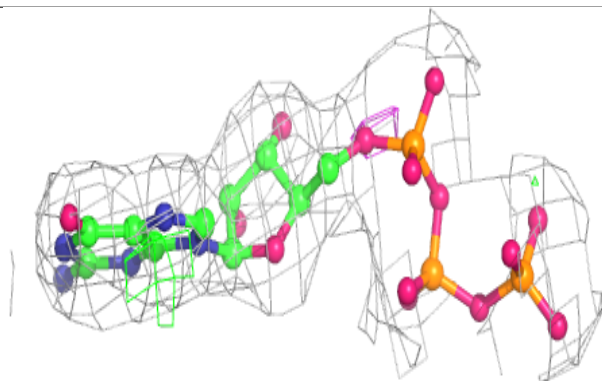
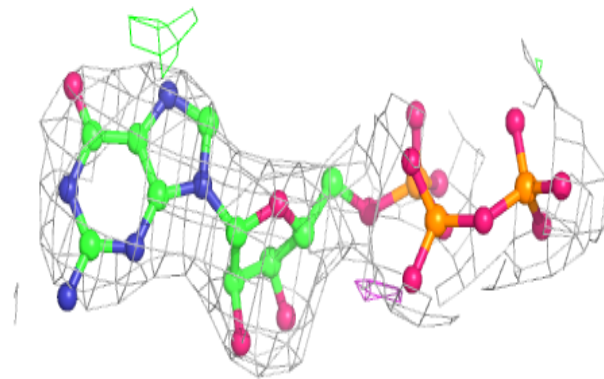
**Electron density around GTP I 401:**

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

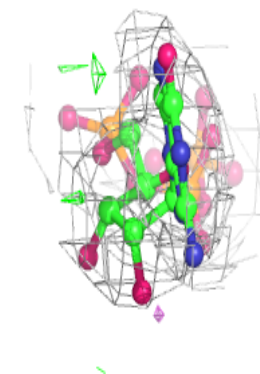
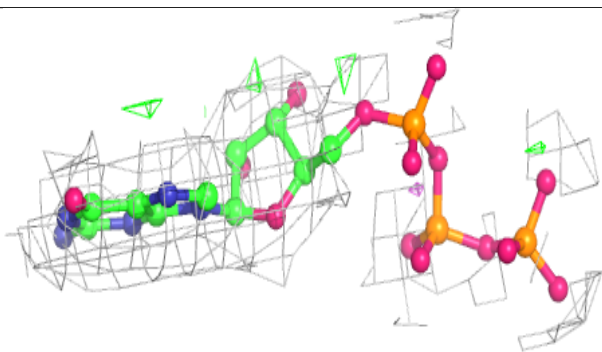
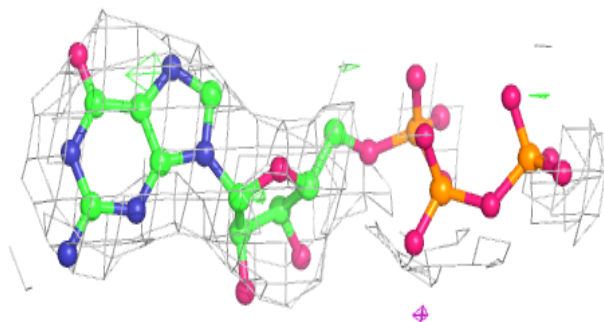


**Electron density around GTP D 302:**

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

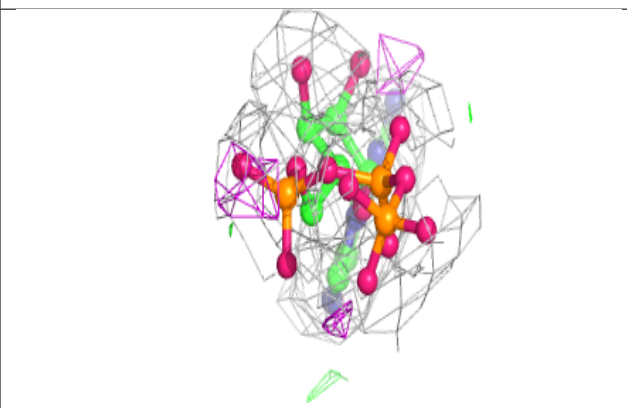
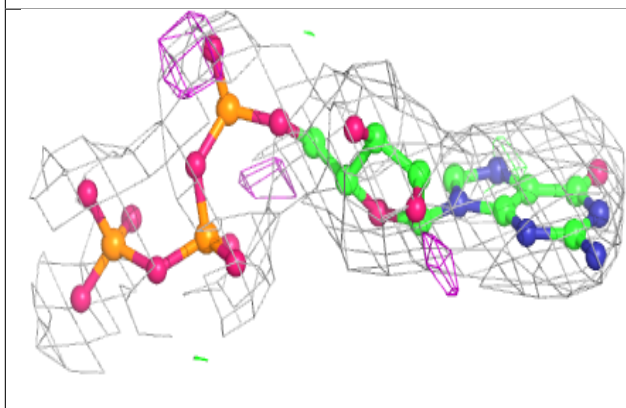
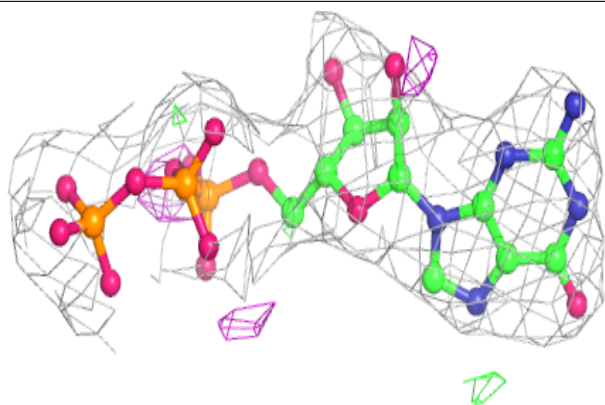
**Electron density around GTP G 401:**

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

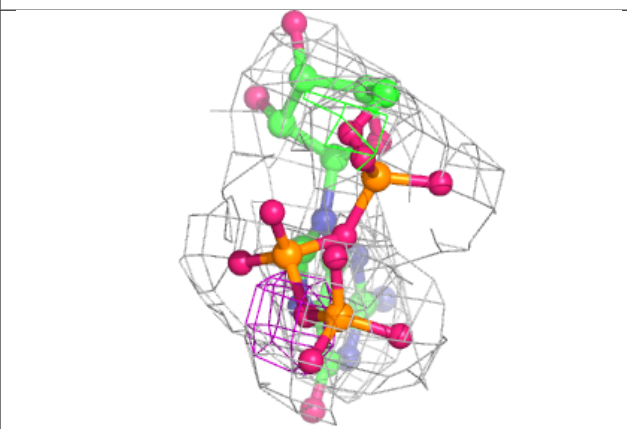
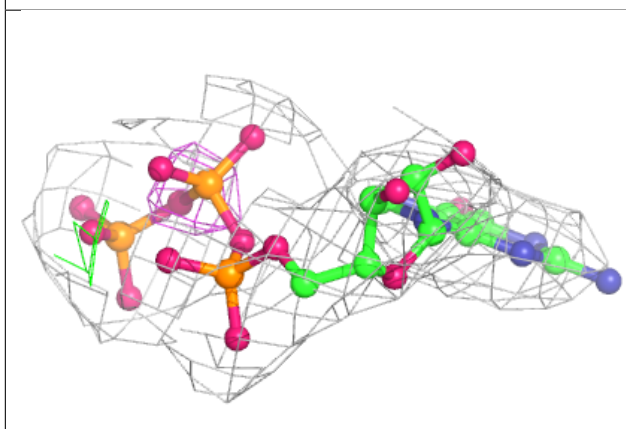
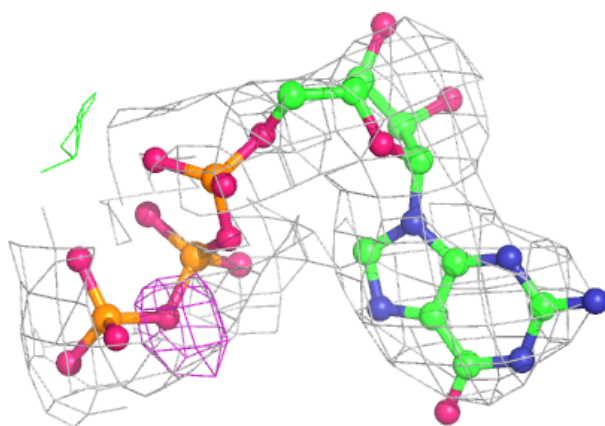


**Electron density around GTP M 401:**

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

**Electron density around GTP N 402:**

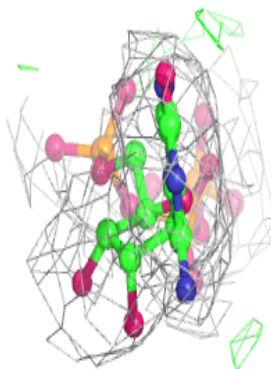
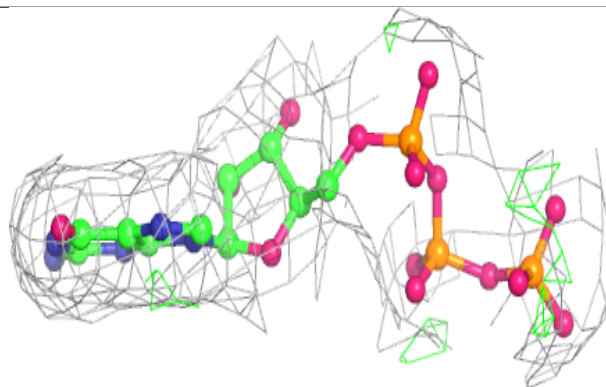
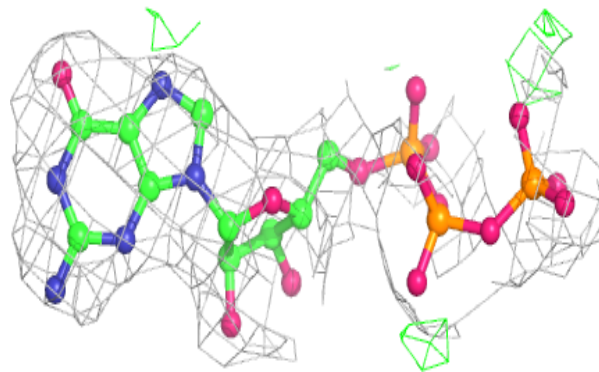
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



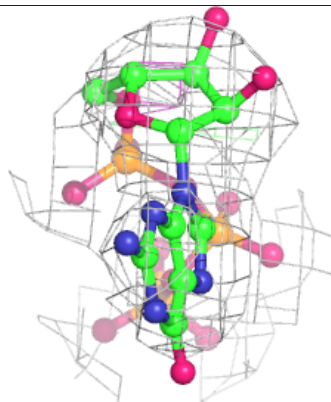
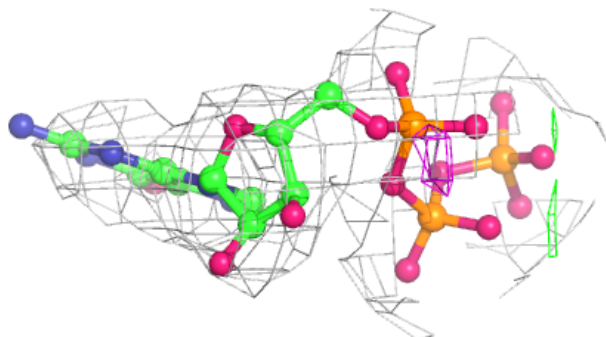
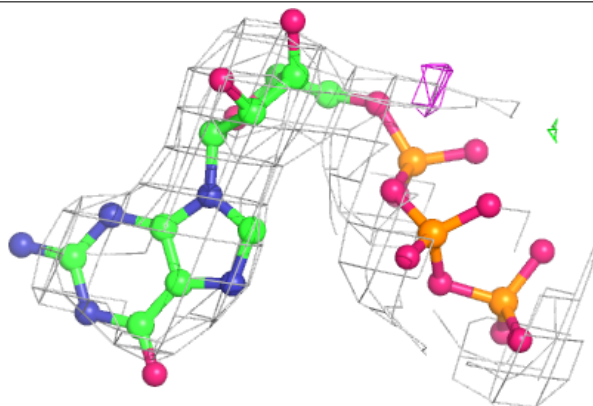


**Electron density around GTP N 401:**

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

**Electron density around GTP J 402:**

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



## 6.5 Other polymers

There are no such residues in this entry.