



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2021 – 02:18 PM EST

PDB ID : 7LU5  
Title : SAMHD1(113-626) H206R D207N R366H  
Authors : Temple, J.T.; Bowen, N.E.  
Deposited on : 2021-02-20  
Resolution : 3.57 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

---

The 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 : 2.23.2  
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) : 2.23.2

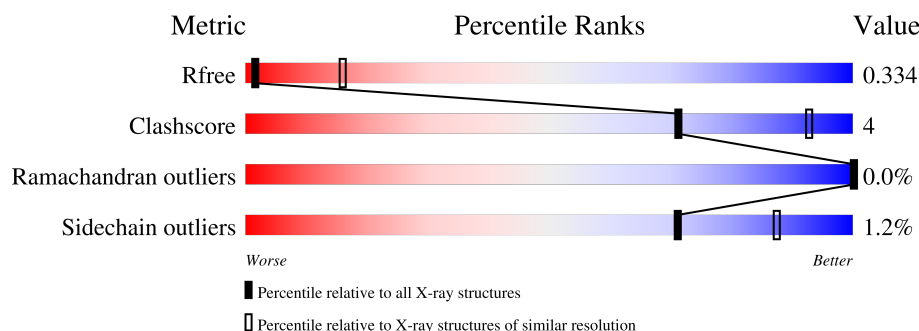
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 3.57 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 <sub>free</sub>	130704	1094 (3.66-3.50)
Clashscore	141614	1181 (3.66-3.50)
Ramachandran outliers	138981	1143 (3.66-3.50)
Sidechain outliers	138945	1143 (3.66-3.50)

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

Mol	Chain	Length	Quality of chain
1	A	535	82% 8% • 10%
1	B	535	81% 8% 10%
1	C	535	82% 8% • 10%
1	D	535	83% 7% 10%
1	E	535	82% 7% 10%
1	F	535	82% 7% • 10%
1	G	535	82% 8% • 10%

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Length	Quality of chain
1	H	535	 83%7%10%
1	I	535	 83%7%10%
1	J	535	 81%8%10%
1	K	535	 81%8%10%
1	L	535	 81%8%10%
1	M	535	 82%7%10%
1	N	535	 82%8%10%
1	O	535	 81%8%10%
1	P	535	 82%7%10%

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 63936 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 Deoxynucleoside triphosphate triphosphohydrolase SAMHD1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	B	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	C	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	D	481	Total	C	N	O	S	0	1	0
			3940	2521	686	713	20			
1	E	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	F	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	G	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	H	481	Total	C	N	O	S	0	1	0
			3940	2521	686	713	20			
1	I	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	J	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	K	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	L	481	Total	C	N	O	S	0	1	0
			3940	2521	686	713	20			
1	M	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	N	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	O	481	Total	C	N	O	S	0	0	0
			3932	2517	684	711	20			
1	P	481	Total	C	N	O	S	0	1	0
			3940	2521	686	713	20			

There are 384 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	92	MET	-	initiating methionine	UNP Q9Y3Z3
A	93	GLY	-	expression tag	UNP Q9Y3Z3
A	94	SER	-	expression tag	UNP Q9Y3Z3
A	95	SER	-	expression tag	UNP Q9Y3Z3
A	96	HIS	-	expression tag	UNP Q9Y3Z3
A	97	HIS	-	expression tag	UNP Q9Y3Z3
A	98	HIS	-	expression tag	UNP Q9Y3Z3
A	99	HIS	-	expression tag	UNP Q9Y3Z3
A	100	HIS	-	expression tag	UNP Q9Y3Z3
A	101	HIS	-	expression tag	UNP Q9Y3Z3
A	102	SER	-	expression tag	UNP Q9Y3Z3
A	103	SER	-	expression tag	UNP Q9Y3Z3
A	104	GLY	-	expression tag	UNP Q9Y3Z3
A	105	LEU	-	expression tag	UNP Q9Y3Z3
A	106	VAL	-	expression tag	UNP Q9Y3Z3
A	107	PRO	-	expression tag	UNP Q9Y3Z3
A	108	ARG	-	expression tag	UNP Q9Y3Z3
A	109	GLY	-	expression tag	UNP Q9Y3Z3
A	110	SER	-	expression tag	UNP Q9Y3Z3
A	111	HIS	-	expression tag	UNP Q9Y3Z3
A	112	MET	-	expression tag	UNP Q9Y3Z3
A	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
A	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
A	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
B	92	MET	-	initiating methionine	UNP Q9Y3Z3
B	93	GLY	-	expression tag	UNP Q9Y3Z3
B	94	SER	-	expression tag	UNP Q9Y3Z3
B	95	SER	-	expression tag	UNP Q9Y3Z3
B	96	HIS	-	expression tag	UNP Q9Y3Z3
B	97	HIS	-	expression tag	UNP Q9Y3Z3
B	98	HIS	-	expression tag	UNP Q9Y3Z3
B	99	HIS	-	expression tag	UNP Q9Y3Z3
B	100	HIS	-	expression tag	UNP Q9Y3Z3
B	101	HIS	-	expression tag	UNP Q9Y3Z3
B	102	SER	-	expression tag	UNP Q9Y3Z3
B	103	SER	-	expression tag	UNP Q9Y3Z3
B	104	GLY	-	expression tag	UNP Q9Y3Z3
B	105	LEU	-	expression tag	UNP Q9Y3Z3
B	106	VAL	-	expression tag	UNP Q9Y3Z3
B	107	PRO	-	expression tag	UNP Q9Y3Z3
B	108	ARG	-	expression tag	UNP Q9Y3Z3
B	109	GLY	-	expression tag	UNP Q9Y3Z3

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
B	110	SER	-	expression tag	UNP Q9Y3Z3
B	111	HIS	-	expression tag	UNP Q9Y3Z3
B	112	MET	-	expression tag	UNP Q9Y3Z3
B	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
B	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
B	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
C	92	MET	-	initiating methionine	UNP Q9Y3Z3
C	93	GLY	-	expression tag	UNP Q9Y3Z3
C	94	SER	-	expression tag	UNP Q9Y3Z3
C	95	SER	-	expression tag	UNP Q9Y3Z3
C	96	HIS	-	expression tag	UNP Q9Y3Z3
C	97	HIS	-	expression tag	UNP Q9Y3Z3
C	98	HIS	-	expression tag	UNP Q9Y3Z3
C	99	HIS	-	expression tag	UNP Q9Y3Z3
C	100	HIS	-	expression tag	UNP Q9Y3Z3
C	101	HIS	-	expression tag	UNP Q9Y3Z3
C	102	SER	-	expression tag	UNP Q9Y3Z3
C	103	SER	-	expression tag	UNP Q9Y3Z3
C	104	GLY	-	expression tag	UNP Q9Y3Z3
C	105	LEU	-	expression tag	UNP Q9Y3Z3
C	106	VAL	-	expression tag	UNP Q9Y3Z3
C	107	PRO	-	expression tag	UNP Q9Y3Z3
C	108	ARG	-	expression tag	UNP Q9Y3Z3
C	109	GLY	-	expression tag	UNP Q9Y3Z3
C	110	SER	-	expression tag	UNP Q9Y3Z3
C	111	HIS	-	expression tag	UNP Q9Y3Z3
C	112	MET	-	expression tag	UNP Q9Y3Z3
C	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
C	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
C	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
D	92	MET	-	initiating methionine	UNP Q9Y3Z3
D	93	GLY	-	expression tag	UNP Q9Y3Z3
D	94	SER	-	expression tag	UNP Q9Y3Z3
D	95	SER	-	expression tag	UNP Q9Y3Z3
D	96	HIS	-	expression tag	UNP Q9Y3Z3
D	97	HIS	-	expression tag	UNP Q9Y3Z3
D	98	HIS	-	expression tag	UNP Q9Y3Z3
D	99	HIS	-	expression tag	UNP Q9Y3Z3
D	100	HIS	-	expression tag	UNP Q9Y3Z3
D	101	HIS	-	expression tag	UNP Q9Y3Z3
D	102	SER	-	expression tag	UNP Q9Y3Z3
D	103	SER	-	expression tag	UNP Q9Y3Z3

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
D	104	GLY	-	expression tag	UNP Q9Y3Z3
D	105	LEU	-	expression tag	UNP Q9Y3Z3
D	106	VAL	-	expression tag	UNP Q9Y3Z3
D	107	PRO	-	expression tag	UNP Q9Y3Z3
D	108	ARG	-	expression tag	UNP Q9Y3Z3
D	109	GLY	-	expression tag	UNP Q9Y3Z3
D	110	SER	-	expression tag	UNP Q9Y3Z3
D	111	HIS	-	expression tag	UNP Q9Y3Z3
D	112	MET	-	expression tag	UNP Q9Y3Z3
D	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
D	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
D	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
E	92	MET	-	initiating methionine	UNP Q9Y3Z3
E	93	GLY	-	expression tag	UNP Q9Y3Z3
E	94	SER	-	expression tag	UNP Q9Y3Z3
E	95	SER	-	expression tag	UNP Q9Y3Z3
E	96	HIS	-	expression tag	UNP Q9Y3Z3
E	97	HIS	-	expression tag	UNP Q9Y3Z3
E	98	HIS	-	expression tag	UNP Q9Y3Z3
E	99	HIS	-	expression tag	UNP Q9Y3Z3
E	100	HIS	-	expression tag	UNP Q9Y3Z3
E	101	HIS	-	expression tag	UNP Q9Y3Z3
E	102	SER	-	expression tag	UNP Q9Y3Z3
E	103	SER	-	expression tag	UNP Q9Y3Z3
E	104	GLY	-	expression tag	UNP Q9Y3Z3
E	105	LEU	-	expression tag	UNP Q9Y3Z3
E	106	VAL	-	expression tag	UNP Q9Y3Z3
E	107	PRO	-	expression tag	UNP Q9Y3Z3
E	108	ARG	-	expression tag	UNP Q9Y3Z3
E	109	GLY	-	expression tag	UNP Q9Y3Z3
E	110	SER	-	expression tag	UNP Q9Y3Z3
E	111	HIS	-	expression tag	UNP Q9Y3Z3
E	112	MET	-	expression tag	UNP Q9Y3Z3
E	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
E	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
E	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
F	92	MET	-	initiating methionine	UNP Q9Y3Z3
F	93	GLY	-	expression tag	UNP Q9Y3Z3
F	94	SER	-	expression tag	UNP Q9Y3Z3
F	95	SER	-	expression tag	UNP Q9Y3Z3
F	96	HIS	-	expression tag	UNP Q9Y3Z3
F	97	HIS	-	expression tag	UNP Q9Y3Z3

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
F	98	HIS	-	expression tag	UNP Q9Y3Z3
F	99	HIS	-	expression tag	UNP Q9Y3Z3
F	100	HIS	-	expression tag	UNP Q9Y3Z3
F	101	HIS	-	expression tag	UNP Q9Y3Z3
F	102	SER	-	expression tag	UNP Q9Y3Z3
F	103	SER	-	expression tag	UNP Q9Y3Z3
F	104	GLY	-	expression tag	UNP Q9Y3Z3
F	105	LEU	-	expression tag	UNP Q9Y3Z3
F	106	VAL	-	expression tag	UNP Q9Y3Z3
F	107	PRO	-	expression tag	UNP Q9Y3Z3
F	108	ARG	-	expression tag	UNP Q9Y3Z3
F	109	GLY	-	expression tag	UNP Q9Y3Z3
F	110	SER	-	expression tag	UNP Q9Y3Z3
F	111	HIS	-	expression tag	UNP Q9Y3Z3
F	112	MET	-	expression tag	UNP Q9Y3Z3
F	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
F	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
F	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
G	92	MET	-	initiating methionine	UNP Q9Y3Z3
G	93	GLY	-	expression tag	UNP Q9Y3Z3
G	94	SER	-	expression tag	UNP Q9Y3Z3
G	95	SER	-	expression tag	UNP Q9Y3Z3
G	96	HIS	-	expression tag	UNP Q9Y3Z3
G	97	HIS	-	expression tag	UNP Q9Y3Z3
G	98	HIS	-	expression tag	UNP Q9Y3Z3
G	99	HIS	-	expression tag	UNP Q9Y3Z3
G	100	HIS	-	expression tag	UNP Q9Y3Z3
G	101	HIS	-	expression tag	UNP Q9Y3Z3
G	102	SER	-	expression tag	UNP Q9Y3Z3
G	103	SER	-	expression tag	UNP Q9Y3Z3
G	104	GLY	-	expression tag	UNP Q9Y3Z3
G	105	LEU	-	expression tag	UNP Q9Y3Z3
G	106	VAL	-	expression tag	UNP Q9Y3Z3
G	107	PRO	-	expression tag	UNP Q9Y3Z3
G	108	ARG	-	expression tag	UNP Q9Y3Z3
G	109	GLY	-	expression tag	UNP Q9Y3Z3
G	110	SER	-	expression tag	UNP Q9Y3Z3
G	111	HIS	-	expression tag	UNP Q9Y3Z3
G	112	MET	-	expression tag	UNP Q9Y3Z3
G	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
G	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
G	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3

*Continued on next page...*



*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
H	92	MET	-	initiating methionine	UNP Q9Y3Z3
H	93	GLY	-	expression tag	UNP Q9Y3Z3
H	94	SER	-	expression tag	UNP Q9Y3Z3
H	95	SER	-	expression tag	UNP Q9Y3Z3
H	96	HIS	-	expression tag	UNP Q9Y3Z3
H	97	HIS	-	expression tag	UNP Q9Y3Z3
H	98	HIS	-	expression tag	UNP Q9Y3Z3
H	99	HIS	-	expression tag	UNP Q9Y3Z3
H	100	HIS	-	expression tag	UNP Q9Y3Z3
H	101	HIS	-	expression tag	UNP Q9Y3Z3
H	102	SER	-	expression tag	UNP Q9Y3Z3
H	103	SER	-	expression tag	UNP Q9Y3Z3
H	104	GLY	-	expression tag	UNP Q9Y3Z3
H	105	LEU	-	expression tag	UNP Q9Y3Z3
H	106	VAL	-	expression tag	UNP Q9Y3Z3
H	107	PRO	-	expression tag	UNP Q9Y3Z3
H	108	ARG	-	expression tag	UNP Q9Y3Z3
H	109	GLY	-	expression tag	UNP Q9Y3Z3
H	110	SER	-	expression tag	UNP Q9Y3Z3
H	111	HIS	-	expression tag	UNP Q9Y3Z3
H	112	MET	-	expression tag	UNP Q9Y3Z3
H	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
H	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
H	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
I	92	MET	-	initiating methionine	UNP Q9Y3Z3
I	93	GLY	-	expression tag	UNP Q9Y3Z3
I	94	SER	-	expression tag	UNP Q9Y3Z3
I	95	SER	-	expression tag	UNP Q9Y3Z3
I	96	HIS	-	expression tag	UNP Q9Y3Z3
I	97	HIS	-	expression tag	UNP Q9Y3Z3
I	98	HIS	-	expression tag	UNP Q9Y3Z3
I	99	HIS	-	expression tag	UNP Q9Y3Z3
I	100	HIS	-	expression tag	UNP Q9Y3Z3
I	101	HIS	-	expression tag	UNP Q9Y3Z3
I	102	SER	-	expression tag	UNP Q9Y3Z3
I	103	SER	-	expression tag	UNP Q9Y3Z3
I	104	GLY	-	expression tag	UNP Q9Y3Z3
I	105	LEU	-	expression tag	UNP Q9Y3Z3
I	106	VAL	-	expression tag	UNP Q9Y3Z3
I	107	PRO	-	expression tag	UNP Q9Y3Z3
I	108	ARG	-	expression tag	UNP Q9Y3Z3
I	109	GLY	-	expression tag	UNP Q9Y3Z3

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
I	110	SER	-	expression tag	UNP Q9Y3Z3
I	111	HIS	-	expression tag	UNP Q9Y3Z3
I	112	MET	-	expression tag	UNP Q9Y3Z3
I	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
I	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
I	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
J	92	MET	-	initiating methionine	UNP Q9Y3Z3
J	93	GLY	-	expression tag	UNP Q9Y3Z3
J	94	SER	-	expression tag	UNP Q9Y3Z3
J	95	SER	-	expression tag	UNP Q9Y3Z3
J	96	HIS	-	expression tag	UNP Q9Y3Z3
J	97	HIS	-	expression tag	UNP Q9Y3Z3
J	98	HIS	-	expression tag	UNP Q9Y3Z3
J	99	HIS	-	expression tag	UNP Q9Y3Z3
J	100	HIS	-	expression tag	UNP Q9Y3Z3
J	101	HIS	-	expression tag	UNP Q9Y3Z3
J	102	SER	-	expression tag	UNP Q9Y3Z3
J	103	SER	-	expression tag	UNP Q9Y3Z3
J	104	GLY	-	expression tag	UNP Q9Y3Z3
J	105	LEU	-	expression tag	UNP Q9Y3Z3
J	106	VAL	-	expression tag	UNP Q9Y3Z3
J	107	PRO	-	expression tag	UNP Q9Y3Z3
J	108	ARG	-	expression tag	UNP Q9Y3Z3
J	109	GLY	-	expression tag	UNP Q9Y3Z3
J	110	SER	-	expression tag	UNP Q9Y3Z3
J	111	HIS	-	expression tag	UNP Q9Y3Z3
J	112	MET	-	expression tag	UNP Q9Y3Z3
J	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
J	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
J	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
K	92	MET	-	initiating methionine	UNP Q9Y3Z3
K	93	GLY	-	expression tag	UNP Q9Y3Z3
K	94	SER	-	expression tag	UNP Q9Y3Z3
K	95	SER	-	expression tag	UNP Q9Y3Z3
K	96	HIS	-	expression tag	UNP Q9Y3Z3
K	97	HIS	-	expression tag	UNP Q9Y3Z3
K	98	HIS	-	expression tag	UNP Q9Y3Z3
K	99	HIS	-	expression tag	UNP Q9Y3Z3
K	100	HIS	-	expression tag	UNP Q9Y3Z3
K	101	HIS	-	expression tag	UNP Q9Y3Z3
K	102	SER	-	expression tag	UNP Q9Y3Z3
K	103	SER	-	expression tag	UNP Q9Y3Z3

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
K	104	GLY	-	expression tag	UNP Q9Y3Z3
K	105	LEU	-	expression tag	UNP Q9Y3Z3
K	106	VAL	-	expression tag	UNP Q9Y3Z3
K	107	PRO	-	expression tag	UNP Q9Y3Z3
K	108	ARG	-	expression tag	UNP Q9Y3Z3
K	109	GLY	-	expression tag	UNP Q9Y3Z3
K	110	SER	-	expression tag	UNP Q9Y3Z3
K	111	HIS	-	expression tag	UNP Q9Y3Z3
K	112	MET	-	expression tag	UNP Q9Y3Z3
K	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
K	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
K	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
L	92	MET	-	initiating methionine	UNP Q9Y3Z3
L	93	GLY	-	expression tag	UNP Q9Y3Z3
L	94	SER	-	expression tag	UNP Q9Y3Z3
L	95	SER	-	expression tag	UNP Q9Y3Z3
L	96	HIS	-	expression tag	UNP Q9Y3Z3
L	97	HIS	-	expression tag	UNP Q9Y3Z3
L	98	HIS	-	expression tag	UNP Q9Y3Z3
L	99	HIS	-	expression tag	UNP Q9Y3Z3
L	100	HIS	-	expression tag	UNP Q9Y3Z3
L	101	HIS	-	expression tag	UNP Q9Y3Z3
L	102	SER	-	expression tag	UNP Q9Y3Z3
L	103	SER	-	expression tag	UNP Q9Y3Z3
L	104	GLY	-	expression tag	UNP Q9Y3Z3
L	105	LEU	-	expression tag	UNP Q9Y3Z3
L	106	VAL	-	expression tag	UNP Q9Y3Z3
L	107	PRO	-	expression tag	UNP Q9Y3Z3
L	108	ARG	-	expression tag	UNP Q9Y3Z3
L	109	GLY	-	expression tag	UNP Q9Y3Z3
L	110	SER	-	expression tag	UNP Q9Y3Z3
L	111	HIS	-	expression tag	UNP Q9Y3Z3
L	112	MET	-	expression tag	UNP Q9Y3Z3
L	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
L	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
L	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
M	92	MET	-	initiating methionine	UNP Q9Y3Z3
M	93	GLY	-	expression tag	UNP Q9Y3Z3
M	94	SER	-	expression tag	UNP Q9Y3Z3
M	95	SER	-	expression tag	UNP Q9Y3Z3
M	96	HIS	-	expression tag	UNP Q9Y3Z3
M	97	HIS	-	expression tag	UNP Q9Y3Z3

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
M	98	HIS	-	expression tag	UNP Q9Y3Z3
M	99	HIS	-	expression tag	UNP Q9Y3Z3
M	100	HIS	-	expression tag	UNP Q9Y3Z3
M	101	HIS	-	expression tag	UNP Q9Y3Z3
M	102	SER	-	expression tag	UNP Q9Y3Z3
M	103	SER	-	expression tag	UNP Q9Y3Z3
M	104	GLY	-	expression tag	UNP Q9Y3Z3
M	105	LEU	-	expression tag	UNP Q9Y3Z3
M	106	VAL	-	expression tag	UNP Q9Y3Z3
M	107	PRO	-	expression tag	UNP Q9Y3Z3
M	108	ARG	-	expression tag	UNP Q9Y3Z3
M	109	GLY	-	expression tag	UNP Q9Y3Z3
M	110	SER	-	expression tag	UNP Q9Y3Z3
M	111	HIS	-	expression tag	UNP Q9Y3Z3
M	112	MET	-	expression tag	UNP Q9Y3Z3
M	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
M	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
M	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
N	92	MET	-	initiating methionine	UNP Q9Y3Z3
N	93	GLY	-	expression tag	UNP Q9Y3Z3
N	94	SER	-	expression tag	UNP Q9Y3Z3
N	95	SER	-	expression tag	UNP Q9Y3Z3
N	96	HIS	-	expression tag	UNP Q9Y3Z3
N	97	HIS	-	expression tag	UNP Q9Y3Z3
N	98	HIS	-	expression tag	UNP Q9Y3Z3
N	99	HIS	-	expression tag	UNP Q9Y3Z3
N	100	HIS	-	expression tag	UNP Q9Y3Z3
N	101	HIS	-	expression tag	UNP Q9Y3Z3
N	102	SER	-	expression tag	UNP Q9Y3Z3
N	103	SER	-	expression tag	UNP Q9Y3Z3
N	104	GLY	-	expression tag	UNP Q9Y3Z3
N	105	LEU	-	expression tag	UNP Q9Y3Z3
N	106	VAL	-	expression tag	UNP Q9Y3Z3
N	107	PRO	-	expression tag	UNP Q9Y3Z3
N	108	ARG	-	expression tag	UNP Q9Y3Z3
N	109	GLY	-	expression tag	UNP Q9Y3Z3
N	110	SER	-	expression tag	UNP Q9Y3Z3
N	111	HIS	-	expression tag	UNP Q9Y3Z3
N	112	MET	-	expression tag	UNP Q9Y3Z3
N	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
N	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
N	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3

*Continued on next page...*

*Continued from previous page...*

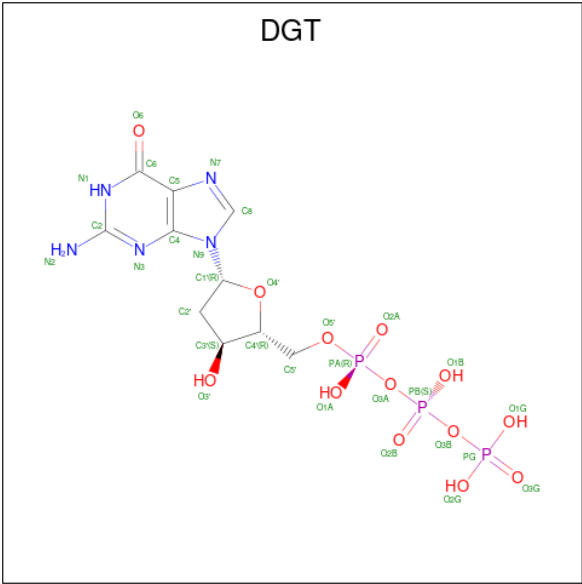
Chain	Residue	Modelled	Actual	Comment	Reference
O	92	MET	-	initiating methionine	UNP Q9Y3Z3
O	93	GLY	-	expression tag	UNP Q9Y3Z3
O	94	SER	-	expression tag	UNP Q9Y3Z3
O	95	SER	-	expression tag	UNP Q9Y3Z3
O	96	HIS	-	expression tag	UNP Q9Y3Z3
O	97	HIS	-	expression tag	UNP Q9Y3Z3
O	98	HIS	-	expression tag	UNP Q9Y3Z3
O	99	HIS	-	expression tag	UNP Q9Y3Z3
O	100	HIS	-	expression tag	UNP Q9Y3Z3
O	101	HIS	-	expression tag	UNP Q9Y3Z3
O	102	SER	-	expression tag	UNP Q9Y3Z3
O	103	SER	-	expression tag	UNP Q9Y3Z3
O	104	GLY	-	expression tag	UNP Q9Y3Z3
O	105	LEU	-	expression tag	UNP Q9Y3Z3
O	106	VAL	-	expression tag	UNP Q9Y3Z3
O	107	PRO	-	expression tag	UNP Q9Y3Z3
O	108	ARG	-	expression tag	UNP Q9Y3Z3
O	109	GLY	-	expression tag	UNP Q9Y3Z3
O	110	SER	-	expression tag	UNP Q9Y3Z3
O	111	HIS	-	expression tag	UNP Q9Y3Z3
O	112	MET	-	expression tag	UNP Q9Y3Z3
O	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
O	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
O	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3
P	92	MET	-	initiating methionine	UNP Q9Y3Z3
P	93	GLY	-	expression tag	UNP Q9Y3Z3
P	94	SER	-	expression tag	UNP Q9Y3Z3
P	95	SER	-	expression tag	UNP Q9Y3Z3
P	96	HIS	-	expression tag	UNP Q9Y3Z3
P	97	HIS	-	expression tag	UNP Q9Y3Z3
P	98	HIS	-	expression tag	UNP Q9Y3Z3
P	99	HIS	-	expression tag	UNP Q9Y3Z3
P	100	HIS	-	expression tag	UNP Q9Y3Z3
P	101	HIS	-	expression tag	UNP Q9Y3Z3
P	102	SER	-	expression tag	UNP Q9Y3Z3
P	103	SER	-	expression tag	UNP Q9Y3Z3
P	104	GLY	-	expression tag	UNP Q9Y3Z3
P	105	LEU	-	expression tag	UNP Q9Y3Z3
P	106	VAL	-	expression tag	UNP Q9Y3Z3
P	107	PRO	-	expression tag	UNP Q9Y3Z3
P	108	ARG	-	expression tag	UNP Q9Y3Z3
P	109	GLY	-	expression tag	UNP Q9Y3Z3

*Continued on next page...*

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
P	110	SER	-	expression tag	UNP Q9Y3Z3
P	111	HIS	-	expression tag	UNP Q9Y3Z3
P	112	MET	-	expression tag	UNP Q9Y3Z3
P	206	ARG	HIS	engineered mutation	UNP Q9Y3Z3
P	207	ASN	ASP	engineered mutation	UNP Q9Y3Z3
P	366	HIS	ARG	engineered mutation	UNP Q9Y3Z3

- Molecule 2 is 2'-DEOXYGUANOSINE-5'-TRIPHOSPHATE (three-letter code: DGT) (formula: C<sub>10</sub>H<sub>16</sub>N<sub>5</sub>O<sub>13</sub>P<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	A	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	B	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	C	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	D	1	Total	C	N	O	P	0	0
			31	10	5	13	3		

Continued on next page...

*Continued from previous page...*

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	E	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	E	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	E	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	F	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	F	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	G	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	G	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	H	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	I	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	I	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	J	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	J	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	J	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	K	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	K	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	L	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	M	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	M	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	N	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	O	1	Total 31	C 10	N 5	O 13	P 3	0	0
2	O	1	Total 31	C 10	N 5	O 13	P 3	0	0

*Continued on next page...*

*Continued from previous page...*

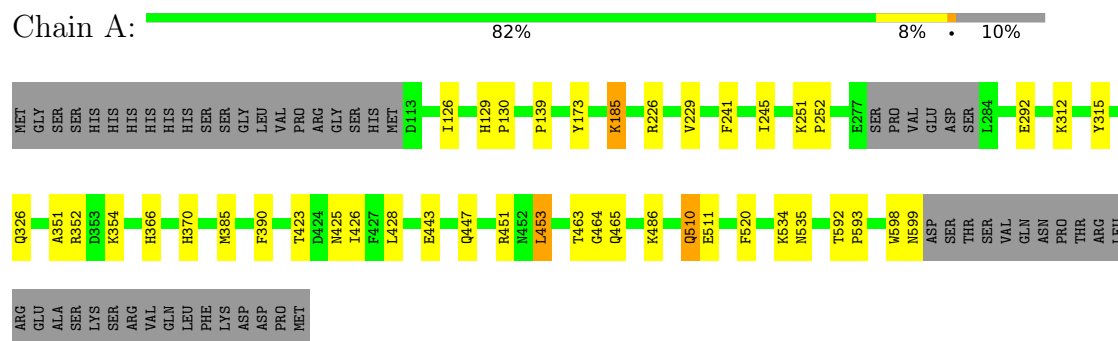
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	O	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	P	1	Total	C	N	O	P	0	0
			31	10	5	13	3		
2	P	1	Total	C	N	O	P	0	0
			31	10	5	13	3		



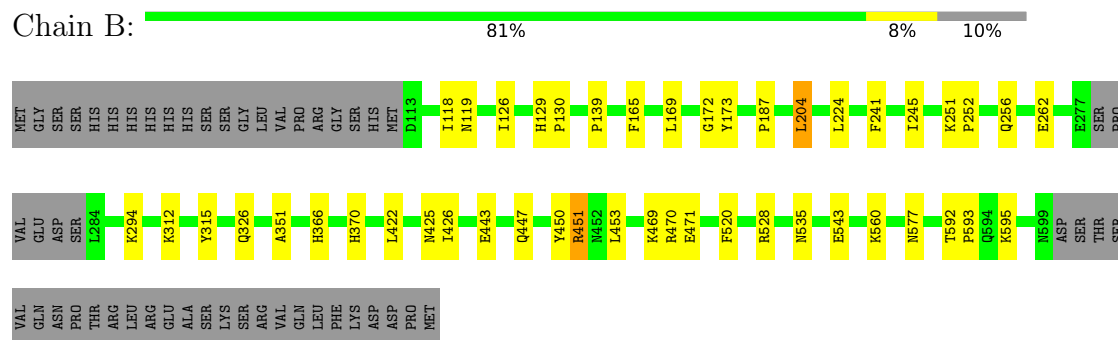
### 3 Residue-property plots

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. 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. 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.

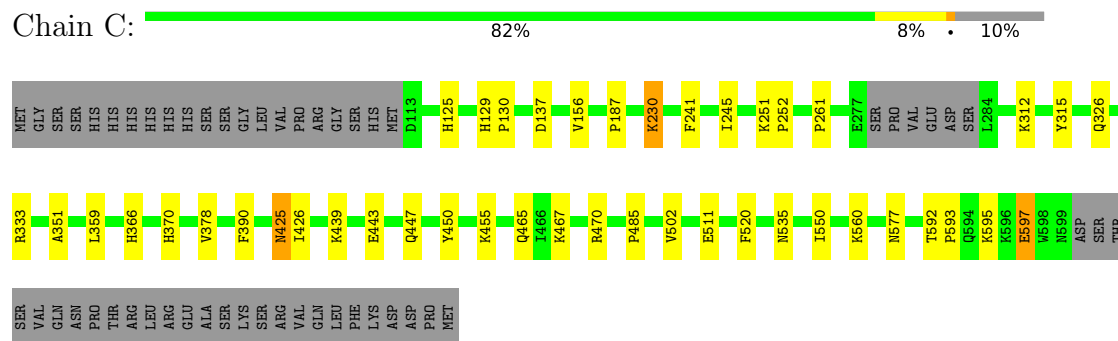
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1




- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

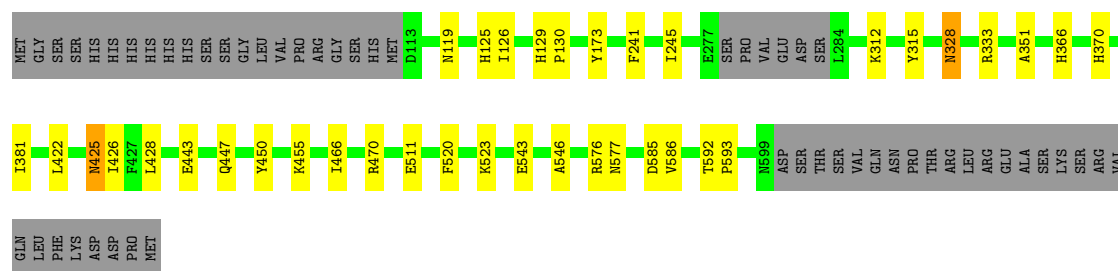


- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1




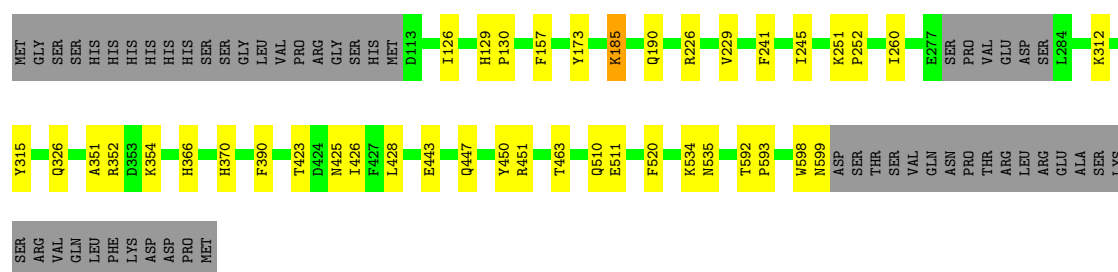
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain D:  83% 7% 10%




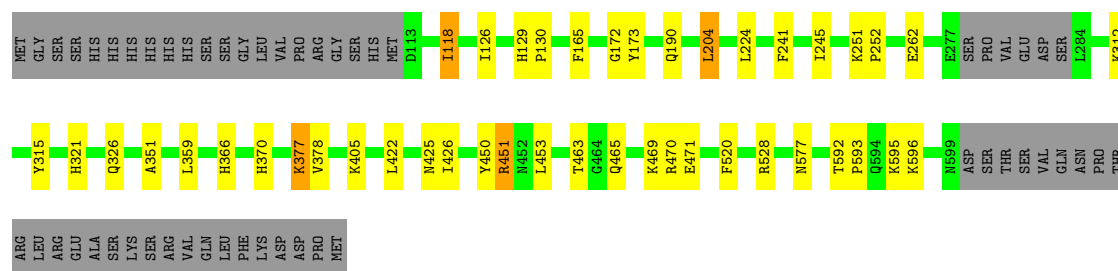
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain E:  82% 7% 10%




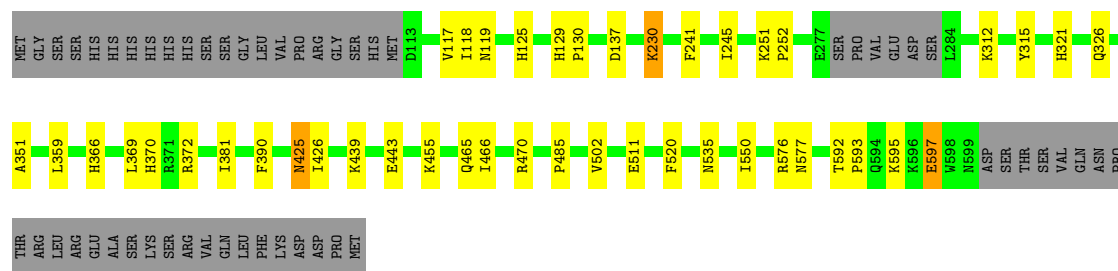
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain F:  82% 7% 10%




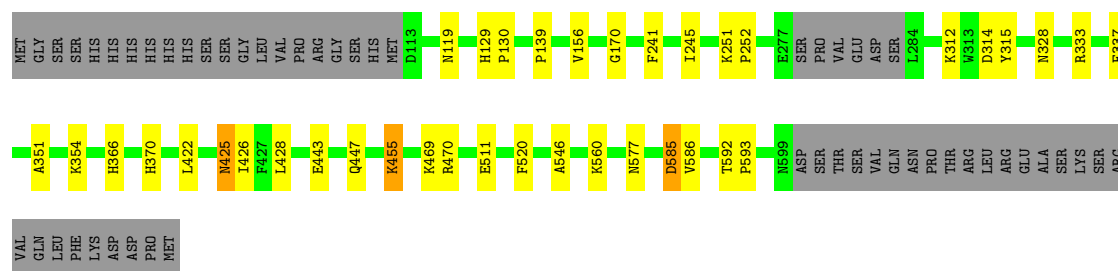
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain G:  82% 8% 10%




- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain H:  83% 7% 10%




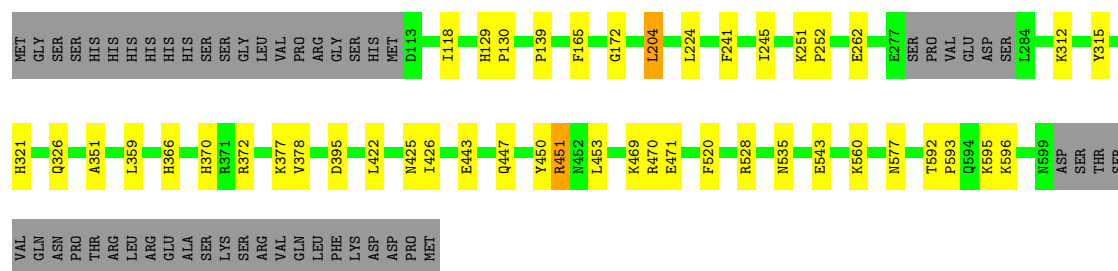
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain I:  83% 7% 10%




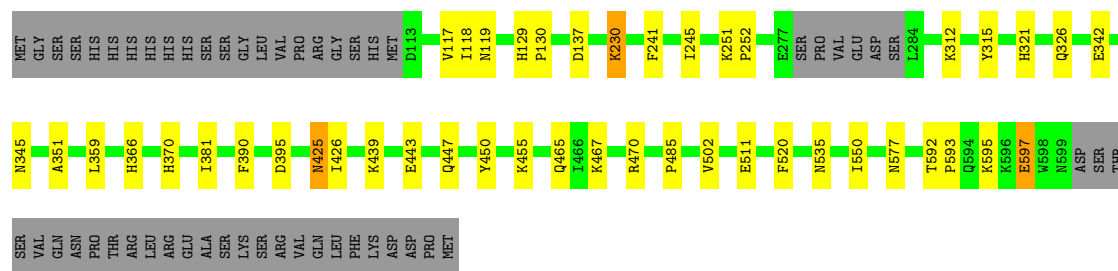
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain J:  81% 8% 10%




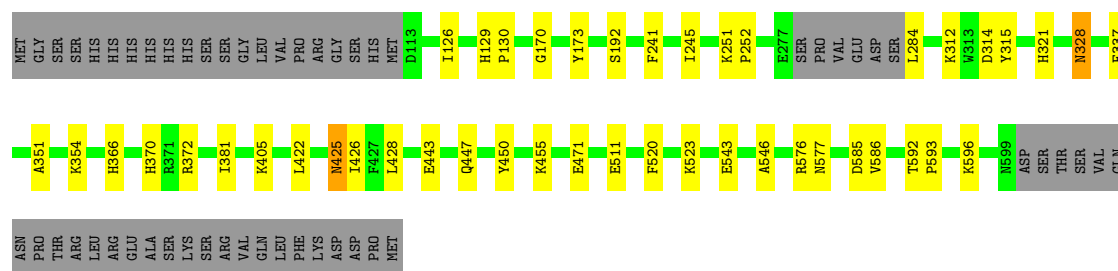
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain K:  81% 8% 10%




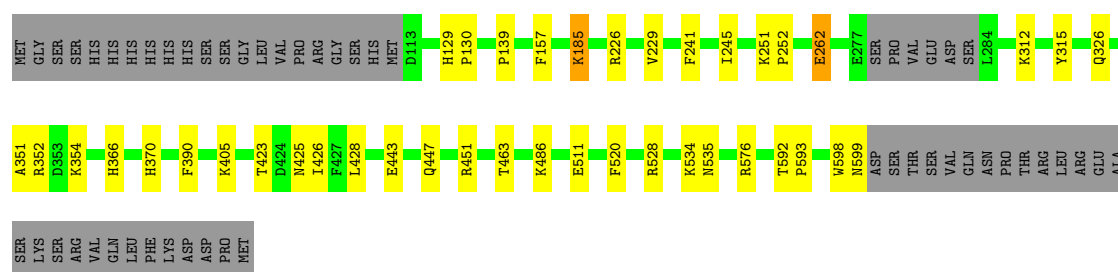
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain L: 




- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain M: 




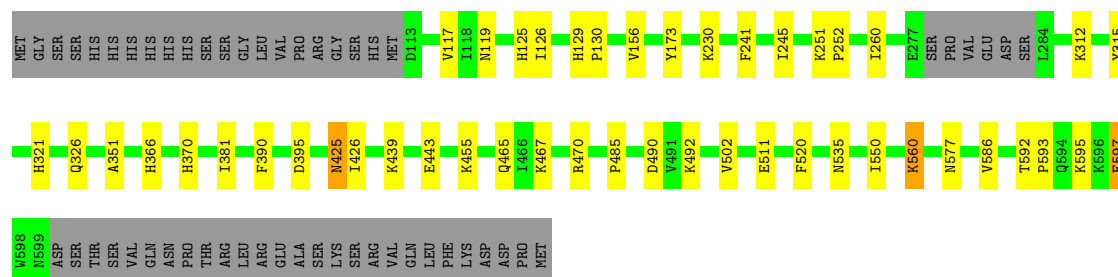
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain N: 



- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain O: 



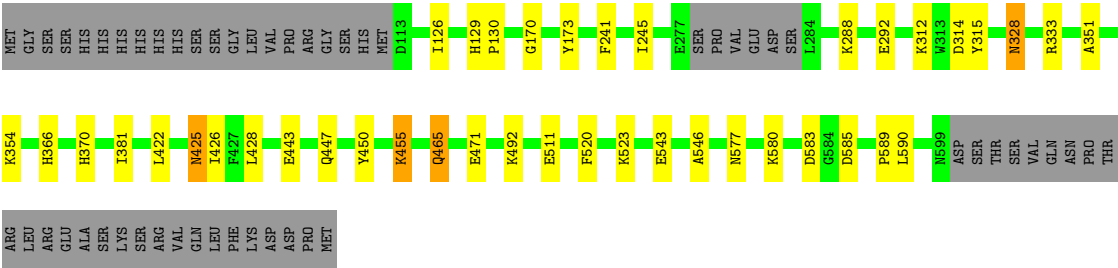
- Molecule 1: Deoxynucleoside triphosphate triphosphohydrolase SAMHD1

Chain P: 

82%

7%

10%



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	83.73Å 573.48Å 100.50Å 90.00° 114.72° 90.00°	Depositor
Resolution (Å)	49.27 – 3.57 49.27 – 3.57	Depositor EDS
% Data completeness (in resolution range)	83.9 (49.27-3.57) 91.6 (49.27-3.57)	Depositor EDS
$R_{merge}$	0.14	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.20 (at 3.57Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, $R_{free}$	0.236 , 0.274 0.297 , 0.334	Depositor DCC
$R_{free}$ test set	4552 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	159.8	Xtriage
Anisotropy	0.126	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.31 , 122.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.036 for h,-k,-h-l	Xtriage
Reported twinning fraction	0.598 for H, K, L 0.402 for -H, -K, H+L	Depositor
Outliers	0 of 92616 reflections	Xtriage
$F_o, F_c$ correlation	0.85	EDS
Total number of atoms	63936	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	188.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% 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: DGT

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	0.62	0/4025	0.73	2/5434 (0.0%)
1	B	0.63	1/4025 (0.0%)	0.77	4/5434 (0.1%)
1	C	0.62	1/4025 (0.0%)	0.72	2/5434 (0.0%)
1	D	0.60	0/4033	0.70	0/5445
1	E	0.61	0/4025	0.72	2/5434 (0.0%)
1	F	0.63	1/4025 (0.0%)	0.76	4/5434 (0.1%)
1	G	0.60	0/4025	0.71	2/5434 (0.0%)
1	H	0.60	0/4033	0.69	0/5445
1	I	0.62	0/4025	0.73	2/5434 (0.0%)
1	J	0.63	1/4025 (0.0%)	0.76	4/5434 (0.1%)
1	K	0.59	0/4025	0.71	2/5434 (0.0%)
1	L	0.60	0/4033	0.69	0/5445
1	M	0.63	1/4025 (0.0%)	0.72	2/5434 (0.0%)
1	N	0.62	0/4025	0.76	4/5434 (0.1%)
1	O	0.60	0/4025	0.71	2/5434 (0.0%)
1	P	0.61	0/4033	0.70	0/5445
All	All	0.61	5/64432 (0.0%)	0.72	32/86988 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	M	262	GLU	CD-OE2	-9.01	1.15	1.25
1	C	261	PRO	N-CD	7.37	1.58	1.47
1	J	262	GLU	CD-OE2	-5.86	1.19	1.25
1	F	262	GLU	CD-OE2	-5.78	1.19	1.25
1	B	262	GLU	CD-OE2	-5.64	1.19	1.25

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	451	ARG	NE-CZ-NH1	-11.67	114.47	120.30
1	N	451	ARG	NE-CZ-NH1	-10.78	114.91	120.30
1	J	451	ARG	NE-CZ-NH1	-10.46	115.07	120.30
1	F	451	ARG	NE-CZ-NH1	-10.21	115.19	120.30
1	B	450	TYR	CB-CG-CD2	-9.76	115.14	121.00
1	N	450	TYR	CB-CG-CD2	-9.75	115.15	121.00
1	J	450	TYR	CB-CG-CD2	-9.69	115.19	121.00
1	F	450	TYR	CB-CG-CD2	-9.43	115.34	121.00
1	B	451	ARG	NE-CZ-NH2	8.30	124.45	120.30
1	N	450	TYR	CB-CG-CD1	7.89	125.73	121.00
1	B	450	TYR	CB-CG-CD1	7.63	125.58	121.00
1	J	450	TYR	CB-CG-CD1	7.61	125.57	121.00
1	F	450	TYR	CB-CG-CD1	7.52	125.51	121.00
1	N	451	ARG	NE-CZ-NH2	7.48	124.04	120.30
1	O	470	ARG	NE-CZ-NH2	-7.28	116.66	120.30
1	F	451	ARG	NE-CZ-NH2	6.87	123.73	120.30
1	J	451	ARG	NE-CZ-NH2	6.76	123.68	120.30
1	C	485	PRO	N-CA-C	6.26	128.38	112.10
1	O	485	PRO	N-CA-C	6.21	128.23	112.10
1	G	485	PRO	N-CA-C	6.17	128.14	112.10
1	K	485	PRO	N-CA-C	6.16	128.11	112.10
1	I	534	LYS	CG-CD-CE	-5.75	94.65	111.90
1	E	534	LYS	CG-CD-CE	-5.74	94.68	111.90
1	M	534	LYS	CG-CD-CE	-5.71	94.76	111.90
1	A	534	LYS	CG-CD-CE	-5.51	95.37	111.90
1	C	470	ARG	NE-CZ-NH2	-5.27	117.66	120.30
1	G	470	ARG	NE-CZ-NH2	-5.21	117.69	120.30
1	I	185	LYS	CB-CG-CD	5.20	125.12	111.60
1	K	470	ARG	NE-CZ-NH2	-5.18	117.71	120.30
1	A	185	LYS	CB-CG-CD	5.13	124.93	111.60
1	E	185	LYS	CB-CG-CD	5.09	124.83	111.60
1	M	185	LYS	CB-CG-CD	5.07	124.78	111.60

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	3932	0	3915	30	2
1	B	3932	0	3915	42	0
1	C	3932	0	3915	35	0
1	D	3940	0	3920	26	2
1	E	3932	0	3915	31	0
1	F	3932	0	3915	35	14
1	G	3932	0	3915	37	14
1	H	3940	0	3920	30	0
1	I	3932	0	3915	29	6
1	J	3932	0	3915	33	6
1	K	3932	0	3915	39	2
1	L	3940	0	3920	31	3
1	M	3932	0	3915	28	5
1	N	3932	0	3915	31	3
1	O	3932	0	3915	38	5
1	P	3940	0	3920	26	6
2	A	93	0	36	5	0
2	B	62	0	24	2	0
2	C	62	0	24	8	0
2	D	31	0	12	1	0
2	E	93	0	36	3	0
2	F	62	0	24	6	0
2	G	62	0	24	4	0
2	H	31	0	12	7	0
2	I	62	0	24	0	0
2	J	93	0	36	11	0
2	K	62	0	24	4	0
2	L	31	0	12	2	0
2	M	62	0	24	1	0
2	N	31	0	12	0	0
2	O	93	0	36	4	0
2	P	62	0	24	6	0
All	All	63936	0	63044	448	34

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 4.

All (448) 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:J:543:GLU:HG3	1:L:543:GLU:HG3	1.72	0.70
1:N:543:GLU:HG3	1:P:543:GLU:HG3	1.73	0.70

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:377:LYS:NZ	2:P:701:DGT:O1G	2.25	0.69
2:D:701:DGT:H5'A	2:D:701:DGT:H8	1.74	0.69
1:J:118:ILE:HG12	2:K:701:DGT:H2'	1.75	0.68
1:B:543:GLU:HG3	1:D:543:GLU:HG3	1.74	0.68
1:D:511:GLU:HG2	1:D:546:ALA:HB3	1.75	0.68
1:H:511:GLU:HG2	1:H:546:ALA:HB3	1.76	0.68
2:J:701:DGT:O2A	1:K:117:VAL:HG23	1.94	0.67
1:N:118:ILE:HG12	2:O:702:DGT:H2'	1.75	0.67
1:F:172:GLY:HA3	1:F:204:LEU:HD13	1.77	0.67
1:B:118:ILE:HG12	2:C:701:DGT:H2'	1.78	0.66
1:F:165:PHE:HZ	2:F:702:DGT:O6	1.78	0.66
1:F:366:HIS:CE1	1:F:370:HIS:CD2	2.84	0.66
1:N:366:HIS:CE1	1:N:370:HIS:CD2	2.84	0.66
1:F:378:VAL:HG21	2:H:701:DGT:O3G	1.96	0.66
1:A:366:HIS:CE1	1:A:370:HIS:CD2	2.84	0.65
1:B:172:GLY:HA3	1:B:204:LEU:HD13	1.77	0.65
1:I:366:HIS:CE1	1:I:370:HIS:CD2	2.84	0.65
1:M:366:HIS:CE1	1:M:370:HIS:CD2	2.84	0.65
1:O:366:HIS:CE1	1:O:370:HIS:CD2	2.84	0.65
2:G:702:DGT:HN2A	1:H:119:ASN:ND2	1.93	0.65
1:D:366:HIS:CE1	1:D:370:HIS:CD2	2.84	0.65
1:J:172:GLY:HA3	1:J:204:LEU:HD13	1.77	0.65
1:G:366:HIS:CE1	1:G:370:HIS:CD2	2.84	0.65
1:L:366:HIS:CE1	1:L:370:HIS:CD2	2.84	0.65
1:J:366:HIS:CE1	1:J:370:HIS:CD2	2.84	0.65
1:C:366:HIS:CE1	1:C:370:HIS:CD2	2.84	0.65
1:H:366:HIS:CE1	1:H:370:HIS:CD2	2.85	0.65
1:B:366:HIS:CE1	1:B:370:HIS:CD2	2.84	0.65
1:N:172:GLY:HA3	1:N:204:LEU:HD13	1.78	0.65
1:P:366:HIS:CE1	1:P:370:HIS:CD2	2.84	0.65
1:E:366:HIS:CE1	1:E:370:HIS:CD2	2.84	0.64
1:K:366:HIS:CE1	1:K:370:HIS:CD2	2.84	0.64
2:C:702:DGT:HN2A	1:D:119:ASN:ND2	1.96	0.64
1:L:511:GLU:HG2	1:L:546:ALA:HB3	1.78	0.64
1:F:118:ILE:HG12	2:F:702:DGT:H2'	1.79	0.64
1:M:511:GLU:O	1:M:511:GLU:HG3	2.00	0.62
1:E:511:GLU:O	1:E:511:GLU:HG3	2.00	0.62
1:J:165:PHE:HZ	2:K:701:DGT:O6	1.82	0.62
1:E:326:GLN:OE1	1:G:326:GLN:HB3	2.00	0.62
1:A:326:GLN:OE1	1:C:326:GLN:HB3	2.00	0.61
1:I:390:PHE:CE2	1:I:426:ILE:HD11	2.35	0.61

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:390:PHE:CE2	1:M:426:ILE:HD11	2.35	0.61
1:I:511:GLU:HG3	1:I:511:GLU:O	2.01	0.61
1:P:511:GLU:HG2	1:P:546:ALA:HB3	1.83	0.61
1:G:117:VAL:HG21	1:H:337:PHE:CZ	2.36	0.61
1:A:511:GLU:HG3	1:A:511:GLU:O	2.00	0.60
1:E:157:PHE:CE2	2:G:702:DGT:H1'	2.36	0.60
1:E:390:PHE:CE2	1:E:426:ILE:HD11	2.35	0.60
1:K:117:VAL:HG21	1:L:337:PHE:CZ	2.36	0.60
1:A:390:PHE:CE2	1:A:426:ILE:HD11	2.36	0.60
1:K:345:ASN:CG	1:O:490:ASP:HB3	2.20	0.60
1:J:165:PHE:CZ	2:K:701:DGT:O6	2.55	0.60
1:F:165:PHE:CZ	2:F:702:DGT:O6	2.56	0.59
1:M:326:GLN:OE1	1:O:326:GLN:HB3	2.03	0.59
2:J:701:DGT:O2A	1:K:117:VAL:CG2	2.51	0.58
1:E:428:LEU:HD13	1:H:425:ASN:HB2	1.85	0.58
1:B:256:GLN:NE2	1:F:190:GLN:OE1	2.37	0.58
1:B:294:LYS:HE3	1:E:190:GLN:HE22	1.68	0.58
1:A:292:GLU:OE1	1:H:560:LYS:NZ	2.31	0.57
1:M:528:ARG:NH1	1:O:586:VAL:HG22	2.19	0.57
1:I:423:THR:O	1:I:426:ILE:HG12	2.05	0.57
1:M:428:LEU:HD13	1:P:425:ASN:HB2	1.86	0.57
1:N:425:ASN:ND2	1:O:425:ASN:OD1	2.38	0.57
1:G:117:VAL:HG23	2:H:701:DGT:O2A	2.04	0.56
1:E:423:THR:O	1:E:426:ILE:HG12	2.05	0.56
1:I:428:LEU:HD13	1:L:425:ASN:HB2	1.86	0.56
1:B:560:LYS:CE	1:P:292:GLU:OE1	2.53	0.56
1:A:423:THR:O	1:A:426:ILE:HG12	2.05	0.56
1:M:423:THR:O	1:M:426:ILE:HG12	2.05	0.56
1:A:366:HIS:CE1	1:A:370:HIS:NE2	2.75	0.55
1:J:378:VAL:HG21	2:J:701:DGT:O3G	2.05	0.55
1:I:366:HIS:CE1	1:I:370:HIS:NE2	2.75	0.55
1:A:428:LEU:HD13	1:D:425:ASN:HB2	1.88	0.55
1:E:366:HIS:CE1	1:E:370:HIS:NE2	2.74	0.55
1:L:366:HIS:CE1	1:L:370:HIS:NE2	2.75	0.55
1:P:366:HIS:CE1	1:P:370:HIS:NE2	2.75	0.55
1:F:366:HIS:CE1	1:F:370:HIS:NE2	2.75	0.55
1:M:428:LEU:CD1	1:P:425:ASN:HB2	2.36	0.55
1:D:366:HIS:CE1	1:D:370:HIS:NE2	2.75	0.55
1:I:326:GLN:OE1	1:K:326:GLN:HB3	2.07	0.55
1:M:366:HIS:CE1	1:M:370:HIS:NE2	2.75	0.55
1:N:378:VAL:HG21	2:P:701:DGT:O3G	2.06	0.55

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:366:HIS:CE1	1:G:370:HIS:NE2	2.75	0.55
1:E:428:LEU:CD1	1:H:425:ASN:HB2	2.36	0.55
1:B:366:HIS:CE1	1:B:370:HIS:NE2	2.75	0.55
1:C:187:PRO:CB	1:O:260:ILE:HD11	2.37	0.55
1:O:366:HIS:CE1	1:O:370:HIS:NE2	2.75	0.55
1:C:366:HIS:CE1	1:C:370:HIS:NE2	2.75	0.54
1:H:366:HIS:CE1	1:H:370:HIS:NE2	2.75	0.54
1:N:366:HIS:CE1	1:N:370:HIS:NE2	2.75	0.54
1:K:345:ASN:OD1	1:O:490:ASP:HB3	2.07	0.54
1:C:455:LYS:HA	1:C:455:LYS:HE2	1.90	0.54
1:J:366:HIS:CE1	1:J:370:HIS:NE2	2.75	0.54
1:K:366:HIS:CE1	1:K:370:HIS:NE2	2.75	0.54
1:B:560:LYS:NZ	1:P:292:GLU:OE1	2.39	0.54
1:I:428:LEU:CD1	1:L:425:ASN:HB2	2.37	0.54
2:J:701:DGT:O1A	1:L:354:LYS:NZ	2.39	0.54
1:K:455:LYS:HA	1:K:455:LYS:HE2	1.90	0.53
1:C:187:PRO:HB2	1:O:260:ILE:CD1	2.38	0.53
1:B:528:ARG:HH11	1:D:586:VAL:HG22	1.72	0.53
1:G:117:VAL:HG21	1:H:337:PHE:HZ	1.73	0.53
1:B:425:ASN:ND2	1:C:425:ASN:OD1	2.42	0.53
1:B:560:LYS:HE2	1:P:292:GLU:OE1	2.09	0.53
1:K:117:VAL:HG21	1:L:337:PHE:HZ	1.72	0.53
1:G:117:VAL:CG2	2:H:701:DGT:O2A	2.58	0.52
1:A:226:ARG:NH2	1:A:229:VAL:HG21	2.26	0.51
1:I:226:ARG:NH2	1:I:229:VAL:HG21	2.25	0.51
1:C:156:VAL:O	2:C:701:DGT:H1'	2.10	0.51
1:G:455:LYS:HE2	1:G:455:LYS:HA	1.92	0.51
1:I:451:ARG:HG3	2:L:701:DGT:N1	2.26	0.51
1:F:425:ASN:ND2	1:G:425:ASN:OD1	2.43	0.51
1:C:390:PHE:CZ	1:C:426:ILE:CG2	2.94	0.51
2:A:702:DGT:HN2A	1:B:119:ASN:HD21	1.59	0.51
1:H:354:LYS:NZ	2:H:701:DGT:O1A	2.35	0.51
1:D:351:ALA:O	1:D:520:PHE:HA	2.12	0.50
1:J:372:ARG:HG2	2:J:701:DGT:O6	2.10	0.50
1:K:390:PHE:CZ	1:K:426:ILE:CG2	2.95	0.50
1:P:351:ALA:O	1:P:520:PHE:HA	2.12	0.50
1:A:351:ALA:O	1:A:520:PHE:HA	2.12	0.50
1:A:425:ASN:HB2	1:D:428:LEU:HD13	1.94	0.50
1:F:351:ALA:O	1:F:520:PHE:HA	2.12	0.50
1:A:428:LEU:CD1	1:D:425:ASN:HB2	2.41	0.50
1:G:351:ALA:O	1:G:520:PHE:HA	2.12	0.50

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:M:351:ALA:O	1:M:520:PHE:HA	2.12	0.50
2:E:702:DGT:H3'	2:F:702:DGT:O3'	2.12	0.50
1:J:543:GLU:CG	1:L:543:GLU:HG3	2.41	0.50
1:L:351:ALA:O	1:L:520:PHE:HA	2.12	0.50
1:M:226:ARG:NH2	1:M:229:VAL:HG21	2.26	0.50
1:O:455:LYS:HE2	1:O:455:LYS:HA	1.94	0.50
1:G:390:PHE:CZ	1:G:426:ILE:CG2	2.94	0.50
1:I:351:ALA:O	1:I:520:PHE:HA	2.12	0.49
1:J:351:ALA:O	1:J:520:PHE:HA	2.12	0.49
1:K:351:ALA:O	1:K:520:PHE:HA	2.12	0.49
1:E:351:ALA:O	1:E:520:PHE:HA	2.12	0.49
1:I:425:ASN:HB2	1:L:428:LEU:HD13	1.95	0.49
1:N:351:ALA:O	1:N:520:PHE:HA	2.12	0.49
1:O:351:ALA:O	1:O:520:PHE:HA	2.12	0.49
1:B:351:ALA:O	1:B:520:PHE:HA	2.12	0.49
1:M:390:PHE:CE2	1:M:426:ILE:CD1	2.96	0.49
1:O:390:PHE:CZ	1:O:426:ILE:CG2	2.95	0.49
1:E:226:ARG:NH2	1:E:229:VAL:HG21	2.26	0.49
2:C:702:DGT:N2	1:D:119:ASN:ND2	2.61	0.49
1:H:351:ALA:O	1:H:520:PHE:HA	2.12	0.49
1:B:326:GLN:HB2	1:D:328:ASN:HA	1.95	0.49
2:A:703:DGT:H5'A	2:C:702:DGT:O1B	2.12	0.49
1:C:351:ALA:O	1:C:520:PHE:HA	2.12	0.49
1:M:425:ASN:HB2	1:P:428:LEU:HD13	1.94	0.49
1:I:592:THR:OG1	1:I:593:PRO:HD3	2.13	0.49
1:J:592:THR:OG1	1:J:593:PRO:HD3	2.13	0.49
1:F:451:ARG:HA	1:F:453:LEU:CD1	2.43	0.48
1:N:451:ARG:HA	1:N:453:LEU:CD1	2.43	0.48
1:K:592:THR:OG1	1:K:593:PRO:HD3	2.13	0.48
2:M:702:DGT:H3'	2:O:702:DGT:O3'	2.13	0.48
1:C:592:THR:OG1	1:C:593:PRO:HD3	2.13	0.48
1:E:352:ARG:HG3	1:E:354:LYS:HG2	1.95	0.48
1:I:139:PRO:HD3	1:L:450:TYR:CE1	2.49	0.48
1:J:451:ARG:HA	1:J:453:LEU:CD1	2.43	0.48
1:J:425:ASN:ND2	1:K:425:ASN:OD1	2.46	0.48
1:A:390:PHE:CE2	1:A:426:ILE:CD1	2.96	0.48
1:B:543:GLU:CG	1:D:543:GLU:HG3	2.43	0.48
1:E:390:PHE:CE2	1:E:426:ILE:CD1	2.96	0.48
1:G:592:THR:OG1	1:G:593:PRO:HD3	2.13	0.48
1:A:352:ARG:HG3	1:A:354:LYS:HG2	1.96	0.48
1:B:592:THR:OG1	1:B:593:PRO:HD3	2.13	0.48

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:592:THR:OG1	1:E:593:PRO:HD3	2.13	0.48
1:B:451:ARG:HA	1:B:453:LEU:CD1	2.43	0.48
1:M:592:THR:OG1	1:M:593:PRO:HD3	2.13	0.48
1:I:390:PHE:CE2	1:I:426:ILE:CD1	2.96	0.47
1:N:577:ASN:OD1	1:N:595:LYS:NZ	2.41	0.47
1:G:119:ASN:CB	2:H:701:DGT:H8	2.44	0.47
1:M:139:PRO:HD3	1:P:450:TYR:CE1	2.49	0.47
1:N:592:THR:OG1	1:N:593:PRO:HD3	2.14	0.47
1:F:592:THR:OG1	1:F:593:PRO:HD3	2.14	0.47
1:M:425:ASN:HB2	1:P:428:LEU:CD1	2.45	0.47
1:A:592:THR:OG1	1:A:593:PRO:HD3	2.13	0.47
2:A:702:DGT:O1B	1:C:378:VAL:HG21	2.15	0.47
1:F:469:LYS:HD3	1:F:471:GLU:OE2	2.15	0.47
1:G:119:ASN:HB2	2:H:701:DGT:H8	1.96	0.47
1:K:595:LYS:HE3	1:K:597:GLU:HG2	1.97	0.47
1:C:592:THR:N	1:C:593:PRO:CD	2.78	0.47
1:E:592:THR:N	1:E:593:PRO:CD	2.78	0.47
1:B:592:THR:N	1:B:593:PRO:CD	2.78	0.47
1:J:592:THR:N	1:J:593:PRO:CD	2.78	0.47
1:A:425:ASN:HB2	1:D:428:LEU:CD1	2.45	0.47
1:F:592:THR:N	1:F:593:PRO:CD	2.78	0.47
1:G:439:LYS:HE2	1:G:443:GLU:CG	2.45	0.47
1:N:592:THR:N	1:N:593:PRO:CD	2.78	0.47
1:B:294:LYS:HE3	1:E:190:GLN:NE2	2.31	0.46
1:B:451:ARG:HG3	2:B:701:DGT:N2	2.29	0.46
1:B:469:LYS:HD3	1:B:471:GLU:OE2	2.15	0.46
1:C:577:ASN:OD1	1:C:595:LYS:NZ	2.41	0.46
1:G:592:THR:N	1:G:593:PRO:CD	2.78	0.46
1:I:352:ARG:HG3	1:I:354:LYS:HG2	1.95	0.46
1:I:592:THR:N	1:I:593:PRO:CD	2.78	0.46
1:O:592:THR:OG1	1:O:593:PRO:HD3	2.13	0.46
1:A:598:TRP:O	1:A:599:ASN:HB2	2.16	0.46
2:A:702:DGT:H3'	2:C:701:DGT:O3'	2.15	0.46
1:I:598:TRP:O	1:I:599:ASN:HB2	2.15	0.46
1:K:592:THR:N	1:K:593:PRO:CD	2.78	0.46
1:N:469:LYS:HD3	1:N:471:GLU:OE2	2.15	0.46
1:O:119:ASN:CB	2:P:701:DGT:H8	2.45	0.46
1:I:157:PHE:CE2	2:K:702:DGT:H1'	2.50	0.46
1:J:359:LEU:HD23	1:J:359:LEU:HA	1.69	0.46
1:J:469:LYS:HD3	1:J:471:GLU:OE2	2.15	0.46
1:N:377:LYS:NZ	2:P:701:DGT:PG	2.87	0.46

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:592:THR:N	1:H:593:PRO:CD	2.79	0.46
1:K:439:LYS:HE2	1:K:443:GLU:CG	2.46	0.46
1:A:592:THR:N	1:A:593:PRO:CD	2.78	0.46
1:B:224:LEU:HD23	1:B:470:ARG:HH22	1.81	0.46
1:C:595:LYS:HG3	1:C:597:GLU:HG2	1.98	0.46
1:D:592:THR:N	1:D:593:PRO:CD	2.79	0.46
1:M:352:ARG:HG3	1:M:354:LYS:HG2	1.96	0.46
1:B:451:ARG:HD3	1:B:453:LEU:CD1	2.46	0.46
1:F:224:LEU:HD23	1:F:470:ARG:HH22	1.81	0.46
1:M:157:PHE:CE2	2:O:703:DGT:H1'	2.51	0.46
1:M:535:ASN:OD1	1:M:535:ASN:N	2.49	0.46
1:O:535:ASN:OD1	1:O:535:ASN:N	2.49	0.46
1:K:535:ASN:OD1	1:K:535:ASN:N	2.49	0.46
1:L:592:THR:N	1:L:593:PRO:CD	2.79	0.46
1:N:326:GLN:HB2	1:P:328:ASN:HA	1.98	0.46
1:A:510:GLN:HB3	1:A:511:GLU:H	1.67	0.46
1:L:192:SER:HB3	1:N:262:GLU:HB2	1.97	0.46
1:M:592:THR:N	1:M:593:PRO:CD	2.78	0.46
1:A:535:ASN:OD1	1:A:535:ASN:N	2.49	0.45
1:C:439:LYS:HE2	1:C:443:GLU:CG	2.46	0.45
1:I:451:ARG:HG3	2:L:701:DGT:C2	2.45	0.45
1:O:592:THR:N	1:O:593:PRO:CD	2.78	0.45
1:E:598:TRP:O	1:E:599:ASN:HB2	2.16	0.45
1:N:224:LEU:HD23	1:N:470:ARG:HH22	1.82	0.45
1:O:439:LYS:HE2	1:O:443:GLU:CG	2.45	0.45
1:G:381:ILE:HD12	1:G:381:ILE:HA	1.90	0.45
1:G:535:ASN:OD1	1:G:535:ASN:N	2.49	0.45
1:J:535:ASN:OD1	1:J:535:ASN:N	2.49	0.45
1:C:535:ASN:OD1	1:C:535:ASN:N	2.49	0.45
1:F:451:ARG:HD3	1:F:453:LEU:CD1	2.46	0.45
1:J:139:PRO:HD3	1:K:450:TYR:CE1	2.51	0.45
1:N:451:ARG:HD3	1:N:453:LEU:CD1	2.46	0.45
1:C:187:PRO:CB	1:O:260:ILE:CD1	2.94	0.45
1:D:381:ILE:HD12	1:D:381:ILE:HA	1.91	0.45
1:E:535:ASN:OD1	1:E:535:ASN:N	2.49	0.45
1:K:595:LYS:HG3	1:K:597:GLU:HG2	1.99	0.45
2:A:702:DGT:HN2A	1:B:119:ASN:ND2	2.14	0.45
1:J:451:ARG:HD3	1:J:453:LEU:CD1	2.46	0.45
1:O:595:LYS:HG3	1:O:597:GLU:HG2	1.98	0.45
1:I:425:ASN:HB2	1:L:428:LEU:CD1	2.46	0.45
1:O:502:VAL:HG22	1:O:550:ILE:HD12	1.98	0.45

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:326:GLN:HB2	1:H:328:ASN:HA	1.99	0.45
1:J:224:LEU:HD23	1:J:470:ARG:HH22	1.81	0.45
1:B:535:ASN:OD1	1:B:535:ASN:N	2.49	0.44
1:G:502:VAL:HG22	1:G:550:ILE:HD12	1.98	0.44
1:G:577:ASN:OD1	1:G:595:LYS:NZ	2.41	0.44
1:I:535:ASN:OD1	1:I:535:ASN:N	2.49	0.44
1:J:326:GLN:HB2	1:L:328:ASN:HA	1.98	0.44
1:K:502:VAL:HG22	1:K:550:ILE:HD12	1.98	0.44
1:N:535:ASN:N	1:N:535:ASN:OD1	2.49	0.44
1:C:333:ARG:HB2	1:D:125:HIS:CE1	2.52	0.44
1:E:425:ASN:HB2	1:H:428:LEU:HD13	1.98	0.44
1:K:312:LYS:HA	1:K:315:TYR:CE2	2.53	0.44
1:A:139:PRO:HD3	1:D:450:TYR:CE1	2.52	0.44
1:D:312:LYS:HA	1:D:315:TYR:CE2	2.53	0.44
1:N:312:LYS:HA	1:N:315:TYR:CE2	2.53	0.44
1:O:312:LYS:HA	1:O:315:TYR:CE2	2.53	0.44
1:O:381:ILE:HD12	1:O:381:ILE:HA	1.90	0.44
1:F:405:LYS:HB3	1:F:405:LYS:HE2	1.77	0.44
1:G:312:LYS:HA	1:G:315:TYR:CE2	2.53	0.44
1:G:595:LYS:HG3	1:G:597:GLU:HG2	1.98	0.44
1:J:312:LYS:HA	1:J:315:TYR:CE2	2.53	0.44
1:M:598:TRP:O	1:M:599:ASN:HB2	2.17	0.44
1:N:405:LYS:HE2	1:N:405:LYS:HB3	1.77	0.44
1:L:192:SER:CB	1:N:262:GLU:HB2	2.48	0.44
1:O:595:LYS:HE3	1:O:597:GLU:HG2	1.99	0.44
1:P:422:LEU:HD12	1:P:426:ILE:HG13	2.00	0.44
1:M:312:LYS:HA	1:M:315:TYR:CE2	2.53	0.44
1:C:502:VAL:HG22	1:C:550:ILE:HD12	1.99	0.44
1:L:422:LEU:HD12	1:L:426:ILE:HG13	1.99	0.44
1:C:312:LYS:HA	1:C:315:TYR:CE2	2.52	0.43
1:D:422:LEU:HD12	1:D:426:ILE:HG13	2.00	0.43
1:E:312:LYS:HA	1:E:315:TYR:CE2	2.53	0.43
1:F:377:LYS:HB2	1:F:377:LYS:HE2	1.33	0.43
1:A:312:LYS:HA	1:A:315:TYR:CE2	2.53	0.43
1:B:312:LYS:HA	1:B:315:TYR:CE2	2.53	0.43
1:F:129:HIS:CG	1:F:130:PRO:HD2	2.54	0.43
1:G:230:LYS:HB3	1:G:230:LYS:HE3	1.64	0.43
1:K:230:LYS:HE3	1:K:230:LYS:HB3	1.65	0.43
1:P:312:LYS:HA	1:P:315:TYR:CE2	2.53	0.43
1:B:451:ARG:HG3	1:C:137:ASP:OD2	2.18	0.43
1:C:230:LYS:HB3	1:C:230:LYS:HE3	1.64	0.43

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:595:LYS:HE3	1:C:597:GLU:HG2	1.99	0.43
1:F:312:LYS:HA	1:F:315:TYR:CE2	2.53	0.43
1:F:577:ASN:OD1	1:F:595:LYS:NZ	2.41	0.43
1:G:595:LYS:HE3	1:G:597:GLU:HG2	1.99	0.43
1:H:170:GLY:HA3	1:H:314:ASP:OD2	2.18	0.43
1:H:312:LYS:HA	1:H:315:TYR:CE2	2.53	0.43
1:K:359:LEU:HA	1:K:359:LEU:HD23	1.70	0.43
1:K:595:LYS:HE3	1:K:597:GLU:CG	2.48	0.43
1:P:170:GLY:HA3	1:P:314:ASP:OD2	2.18	0.43
1:A:129:HIS:CG	1:A:130:PRO:HD2	2.54	0.43
1:F:528:ARG:HH12	1:H:585:ASP:HB3	1.84	0.43
1:K:129:HIS:CG	1:K:130:PRO:HD2	2.53	0.43
1:O:156:VAL:O	2:O:702:DGT:H1'	2.18	0.43
1:B:451:ARG:HG3	2:B:701:DGT:HN2	1.83	0.43
1:I:312:LYS:HA	1:I:315:TYR:CE2	2.53	0.43
1:J:451:ARG:HG3	1:K:137:ASP:OD2	2.18	0.43
1:L:170:GLY:HA3	1:L:314:ASP:OD1	2.19	0.43
1:C:187:PRO:HB3	1:O:260:ILE:HD11	1.99	0.43
1:J:577:ASN:OD1	1:J:595:LYS:NZ	2.41	0.43
1:E:129:HIS:CG	1:E:130:PRO:HD2	2.54	0.43
1:J:528:ARG:HH11	1:L:586:VAL:HG22	1.83	0.43
1:L:312:LYS:HA	1:L:315:TYR:CE2	2.53	0.43
1:O:560:LYS:H	1:O:560:LYS:HG2	1.51	0.43
1:B:129:HIS:CG	1:B:130:PRO:HD2	2.53	0.43
1:C:125:HIS:NE2	1:D:333:ARG:HD2	2.34	0.43
1:G:118:ILE:HG12	2:G:701:DGT:H2'A	2.01	0.43
1:G:129:HIS:CG	1:G:130:PRO:HD2	2.54	0.43
1:H:129:HIS:CG	1:H:130:PRO:HD2	2.53	0.43
1:D:129:HIS:CG	1:D:130:PRO:HD2	2.54	0.43
2:F:701:DGT:H8	2:F:701:DGT:H2'A	1.89	0.43
1:H:241:PHE:O	1:H:245:ILE:HG12	2.19	0.43
1:J:129:HIS:CG	1:J:130:PRO:HD2	2.54	0.43
2:J:702:DGT:H2'A	1:K:118:ILE:HG12	2.00	0.43
1:L:381:ILE:HD12	1:L:381:ILE:HA	1.91	0.43
1:N:543:GLU:CG	1:P:543:GLU:HG3	2.43	0.43
1:C:129:HIS:CG	1:C:130:PRO:HD2	2.54	0.43
1:J:241:PHE:O	1:J:245:ILE:HG12	2.19	0.43
1:L:129:HIS:CG	1:L:130:PRO:HD2	2.53	0.43
1:N:129:HIS:CG	1:N:130:PRO:HD2	2.53	0.43
1:O:129:HIS:CG	1:O:130:PRO:HD2	2.53	0.43
1:A:241:PHE:O	1:A:245:ILE:HG12	2.19	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:422:LEU:HD12	1:B:426:ILE:HG13	2.01	0.42
1:M:129:HIS:CG	1:M:130:PRO:HD2	2.53	0.42
1:C:241:PHE:O	1:C:245:ILE:HG12	2.19	0.42
1:E:450:TYR:CE1	1:H:139:PRO:HD3	2.54	0.42
2:E:701:DGT:O1G	1:H:455:LYS:NZ	2.42	0.42
1:I:129:HIS:CG	1:I:130:PRO:HD2	2.54	0.42
1:J:422:LEU:HD12	1:J:426:ILE:HG13	2.01	0.42
1:N:241:PHE:O	1:N:245:ILE:HG12	2.19	0.42
1:N:422:LEU:HD12	1:N:426:ILE:HG13	2.01	0.42
1:F:321:HIS:CE1	1:G:321:HIS:CE1	3.06	0.42
1:J:251:LYS:HB2	1:J:252:PRO:HD3	2.02	0.42
1:P:129:HIS:CG	1:P:130:PRO:HD2	2.54	0.42
2:C:702:DGT:HN2A	1:D:119:ASN:HD21	1.63	0.42
2:J:703:DGT:O6	1:L:372:ARG:HG2	2.20	0.42
1:K:241:PHE:O	1:K:245:ILE:HG12	2.19	0.42
1:C:251:LYS:HB2	1:C:252:PRO:HD3	2.02	0.42
1:E:425:ASN:HB2	1:H:428:LEU:CD1	2.49	0.42
1:F:241:PHE:O	1:F:245:ILE:HG12	2.19	0.42
1:N:359:LEU:HD23	1:N:359:LEU:HA	1.69	0.42
1:P:241:PHE:O	1:P:245:ILE:HG12	2.19	0.42
1:E:598:TRP:O	1:E:599:ASN:CB	2.68	0.42
2:E:702:DGT:O6	1:G:372:ARG:HG2	2.19	0.42
1:F:377:LYS:H	1:F:377:LYS:HG3	1.67	0.42
2:G:702:DGT:N2	1:H:119:ASN:ND2	2.64	0.42
1:I:226:ARG:HH21	1:I:229:VAL:HG21	1.85	0.42
1:L:241:PHE:O	1:L:245:ILE:HG12	2.19	0.42
1:N:321:HIS:CE1	1:O:321:HIS:CE1	3.08	0.42
1:B:241:PHE:O	1:B:245:ILE:HG12	2.19	0.42
1:C:595:LYS:HE3	1:C:597:GLU:CG	2.49	0.42
1:F:451:ARG:HG3	1:G:137:ASP:OD2	2.20	0.42
1:H:422:LEU:HD12	1:H:426:ILE:HG13	2.00	0.42
1:M:241:PHE:O	1:M:245:ILE:HG12	2.19	0.42
1:A:226:ARG:HH21	1:A:229:VAL:HG21	1.85	0.42
1:E:241:PHE:O	1:E:245:ILE:HG12	2.19	0.42
1:F:359:LEU:HA	1:F:359:LEU:HD23	1.69	0.42
1:G:241:PHE:O	1:G:245:ILE:HG12	2.19	0.42
1:O:251:LYS:HB2	1:O:252:PRO:HD3	2.02	0.42
1:O:577:ASN:OD1	1:O:595:LYS:NZ	2.41	0.42
1:C:359:LEU:HA	1:C:359:LEU:HD23	1.70	0.42
1:D:241:PHE:O	1:D:245:ILE:HG12	2.19	0.42
1:E:425:ASN:ND2	1:H:425:ASN:OD1	2.53	0.42

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:251:LYS:HB2	1:F:252:PRO:HD3	2.02	0.42
1:F:422:LEU:HD12	1:F:426:ILE:HG13	2.01	0.42
1:I:241:PHE:O	1:I:245:ILE:HG12	2.19	0.42
1:O:119:ASN:ND2	2:P:701:DGT:H8	2.35	0.42
1:O:465:GLN:OE1	1:O:465:GLN:HA	2.20	0.42
1:A:598:TRP:O	1:A:599:ASN:CB	2.68	0.42
1:G:125:HIS:NE2	1:H:333:ARG:HD2	2.35	0.42
1:G:465:GLN:OE1	1:G:465:GLN:HA	2.20	0.42
1:G:595:LYS:HE3	1:G:597:GLU:CG	2.49	0.42
1:K:465:GLN:OE1	1:K:465:GLN:HA	2.20	0.42
1:M:251:LYS:HB2	1:M:252:PRO:HD3	2.02	0.42
1:O:595:LYS:HE3	1:O:597:GLU:CG	2.49	0.41
1:B:118:ILE:HG12	2:C:701:DGT:C2'	2.49	0.41
1:F:528:ARG:HH11	1:H:586:VAL:HG22	1.85	0.41
1:I:321:HIS:CE1	1:L:321:HIS:CE1	3.09	0.41
1:M:598:TRP:O	1:M:599:ASN:CB	2.67	0.41
1:O:241:PHE:O	1:O:245:ILE:HG12	2.19	0.41
1:B:251:LYS:HB2	1:B:252:PRO:HD3	2.02	0.41
1:B:256:GLN:NE2	1:F:190:GLN:HE22	2.18	0.41
1:B:577:ASN:OD1	1:B:595:LYS:NZ	2.41	0.41
2:J:701:DGT:H1'	1:K:119:ASN:HB2	2.02	0.41
1:L:251:LYS:HB2	1:L:252:PRO:HD3	2.03	0.41
1:L:284:LEU:HD23	1:L:284:LEU:HA	1.92	0.41
1:H:251:LYS:HB2	1:H:252:PRO:HD3	2.02	0.41
1:J:321:HIS:CE1	1:K:321:HIS:CE1	3.08	0.41
1:K:381:ILE:HD12	1:K:381:ILE:HA	1.90	0.41
1:K:465:GLN:O	1:K:467:LYS:HG3	2.21	0.41
1:D:126:ILE:HG12	1:D:173:TYR:CD1	2.55	0.41
1:F:126:ILE:HG12	1:F:173:TYR:CD1	2.56	0.41
1:I:598:TRP:O	1:I:599:ASN:CB	2.68	0.41
1:O:465:GLN:O	1:O:467:LYS:HG3	2.21	0.41
1:A:385:MET:SD	1:A:453:LEU:HB3	2.61	0.41
1:B:256:GLN:NE2	1:F:190:GLN:NE2	2.68	0.41
1:K:577:ASN:OD1	1:K:595:LYS:NZ	2.41	0.41
2:J:701:DGT:H8	1:K:119:ASN:HB2	2.03	0.41
1:A:251:LYS:HB2	1:A:252:PRO:HD3	2.03	0.41
1:C:465:GLN:OE1	1:C:465:GLN:HA	2.20	0.41
1:E:443:GLU:O	1:E:447:GLN:HG2	2.21	0.41
1:F:465:GLN:O	1:F:465:GLN:HG3	2.19	0.41
1:K:443:GLU:O	1:K:447:GLN:HG2	2.21	0.41
1:N:126:ILE:HG12	1:N:173:TYR:CD1	2.56	0.41

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:126:ILE:HG12	1:P:173:TYR:CD1	2.56	0.41
1:B:139:PRO:HD3	1:C:450:TYR:CE1	2.56	0.41
1:C:465:GLN:O	1:C:467:LYS:HG3	2.21	0.41
1:E:126:ILE:HG12	1:E:173:TYR:CD1	2.56	0.41
1:G:251:LYS:HB2	1:G:252:PRO:HD3	2.02	0.41
1:O:125:HIS:NE2	1:P:333:ARG:HD2	2.36	0.41
1:P:354:LYS:NZ	2:P:701:DGT:O1A	2.33	0.41
1:P:443:GLU:O	1:P:447:GLN:HG2	2.21	0.41
1:E:251:LYS:HB2	1:E:252:PRO:HD3	2.03	0.41
1:G:119:ASN:HB2	2:H:701:DGT:H1'	2.03	0.41
1:G:359:LEU:HA	1:G:359:LEU:HD23	1.70	0.41
1:P:455:LYS:HE3	1:P:455:LYS:HA	2.03	0.41
1:B:187:PRO:CB	1:E:260:ILE:HD11	2.51	0.40
1:C:443:GLU:O	1:C:447:GLN:HG2	2.21	0.40
2:F:701:DGT:H2'	1:H:156:VAL:HG12	2.02	0.40
1:H:443:GLU:O	1:H:447:GLN:HG2	2.21	0.40
1:I:381:ILE:HD12	1:I:381:ILE:HA	1.91	0.40
2:J:701:DGT:H8	1:K:119:ASN:CB	2.51	0.40
1:L:126:ILE:HG12	1:L:173:TYR:CD1	2.56	0.40
1:L:443:GLU:O	1:L:447:GLN:HG2	2.21	0.40
1:M:443:GLU:O	1:M:447:GLN:HG2	2.21	0.40
1:B:126:ILE:HG12	1:B:173:TYR:CD1	2.56	0.40
1:O:492:LYS:HE3	1:O:492:LYS:HB3	1.94	0.40
1:P:381:ILE:HD12	1:P:381:ILE:HA	1.91	0.40
1:A:443:GLU:O	1:A:447:GLN:HG2	2.21	0.40
1:D:443:GLU:O	1:D:447:GLN:HG2	2.21	0.40
1:G:369:LEU:HD23	1:G:369:LEU:HA	1.96	0.40
1:J:451:ARG:HG3	2:J:702:DGT:N2	2.37	0.40
1:K:251:LYS:HB2	1:K:252:PRO:HD3	2.02	0.40
1:A:126:ILE:HG12	1:A:173:TYR:CD1	2.57	0.40
1:J:443:GLU:O	1:J:447:GLN:HG2	2.22	0.40
1:M:226:ARG:HH21	1:M:229:VAL:HG21	1.87	0.40
1:N:251:LYS:HB2	1:N:252:PRO:HD3	2.02	0.40
1:O:126:ILE:HG12	1:O:173:TYR:CD1	2.56	0.40
1:B:165:PHE:CZ	1:B:169:LEU:HD11	2.56	0.40
1:B:443:GLU:O	1:B:447:GLN:HG2	2.22	0.40

All (34) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:492:LYS:CE	1:J:395:ASP:OD1[1_454]	1.17	1.03
1:F:465:GLN:OE1	1:G:576:ARG:O[1_455]	1.19	1.01
1:F:465:GLN:OE1	1:G:576:ARG:C[1_455]	1.22	0.98
1:F:465:GLN:CD	1:G:576:ARG:C[1_455]	1.31	0.89
1:F:465:GLN:CD	1:G:576:ARG:O[1_455]	1.34	0.86
1:F:465:GLN:OE1	1:G:577:ASN:N[1_455]	1.41	0.79
1:M:405:LYS:NZ	1:N:396:TYR:OH[1_454]	1.43	0.77
1:I:492:LYS:CE	1:J:395:ASP:CG[1_454]	1.48	0.72
1:O:395:ASP:OD1	1:P:492:LYS:CE[1_556]	1.57	0.63
1:F:465:GLN:CG	1:G:576:ARG:O[1_455]	1.67	0.53
1:F:465:GLN:OE1	1:G:577:ASN:CA[1_455]	1.68	0.52
1:F:463:THR:O	1:G:466:ILE:CD1[1_455]	1.74	0.46
1:J:560:LYS:NZ	1:M:262:GLU:OE1[1_656]	1.77	0.43
1:F:465:GLN:CD	1:G:577:ASN:N[1_455]	1.78	0.42
1:I:492:LYS:NZ	1:J:395:ASP:OD1[1_454]	1.82	0.38
1:A:464:GLY:CA	1:D:466:ILE:CD1[1_455]	1.88	0.32
1:A:465:GLN:OE1	1:D:576:ARG:O[1_455]	1.88	0.32
1:I:492:LYS:CE	1:J:395:ASP:CB[1_454]	1.88	0.32
1:K:597:GLU:O	1:L:596:LYS:NZ[1_455]	1.89	0.31
1:O:395:ASP:OD1	1:P:492:LYS:NZ[1_556]	1.89	0.31
1:I:492:LYS:CD	1:J:395:ASP:CB[1_454]	1.90	0.30
1:O:395:ASP:CG	1:P:492:LYS:CE[1_556]	1.93	0.27
1:O:395:ASP:CG	1:P:492:LYS:CD[1_556]	1.98	0.22
1:F:465:GLN:NE2	1:G:577:ASN:N[1_455]	1.99	0.21
1:F:465:GLN:OE1	1:G:577:ASN:C[1_455]	2.01	0.19
1:M:405:LYS:CD	1:N:395:ASP:OD2[1_454]	2.01	0.19
1:M:576:ARG:NH1	1:P:465:GLN:OE1[1_455]	2.02	0.18
1:M:405:LYS:CE	1:N:395:ASP:OD2[1_454]	2.03	0.17
1:F:465:GLN:NE2	1:G:576:ARG:C[1_455]	2.05	0.15
1:K:395:ASP:OD2	1:L:405:LYS:NZ[1_556]	2.10	0.10
1:F:465:GLN:NE2	1:G:577:ASN:CG[1_455]	2.12	0.08
1:F:463:THR:C	1:G:466:ILE:CD1[1_455]	2.14	0.06
1:I:465:GLN:OE1	1:L:576:ARG:CZ[1_455]	2.15	0.05
1:O:395:ASP:OD2	1:P:492:LYS:CE[1_556]	2.18	0.02

## 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	477/535 (89%)	469 (98%)	8 (2%)	0	100	100
1	B	477/535 (89%)	469 (98%)	8 (2%)	0	100	100
1	C	477/535 (89%)	470 (98%)	7 (2%)	0	100	100
1	D	478/535 (89%)	471 (98%)	7 (2%)	0	100	100
1	E	477/535 (89%)	468 (98%)	9 (2%)	0	100	100
1	F	477/535 (89%)	469 (98%)	8 (2%)	0	100	100
1	G	477/535 (89%)	470 (98%)	7 (2%)	0	100	100
1	H	478/535 (89%)	470 (98%)	8 (2%)	0	100	100
1	I	477/535 (89%)	466 (98%)	11 (2%)	0	100	100
1	J	477/535 (89%)	469 (98%)	8 (2%)	0	100	100
1	K	477/535 (89%)	470 (98%)	7 (2%)	0	100	100
1	L	478/535 (89%)	471 (98%)	7 (2%)	0	100	100
1	M	477/535 (89%)	468 (98%)	9 (2%)	0	100	100
1	N	477/535 (89%)	468 (98%)	9 (2%)	0	100	100
1	O	477/535 (89%)	470 (98%)	7 (2%)	0	100	100
1	P	478/535 (89%)	470 (98%)	7 (2%)	1 (0%)	47	80
All	All	7636/8560 (89%)	7508 (98%)	127 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	P	589	PRO

### 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	427/477 (90%)	421 (99%)	6 (1%)	67	85

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	427/477 (90%)	426 (100%)	1 (0%)	93	98
1	C	427/477 (90%)	422 (99%)	5 (1%)	71	87
1	D	428/477 (90%)	421 (98%)	7 (2%)	62	83
1	E	427/477 (90%)	423 (99%)	4 (1%)	78	90
1	F	427/477 (90%)	423 (99%)	4 (1%)	78	90
1	G	427/477 (90%)	423 (99%)	4 (1%)	78	90
1	H	428/477 (90%)	422 (99%)	6 (1%)	67	85
1	I	427/477 (90%)	423 (99%)	4 (1%)	78	90
1	J	427/477 (90%)	424 (99%)	3 (1%)	84	93
1	K	427/477 (90%)	422 (99%)	5 (1%)	71	87
1	L	428/477 (90%)	421 (98%)	7 (2%)	62	83
1	M	427/477 (90%)	423 (99%)	4 (1%)	78	90
1	N	427/477 (90%)	424 (99%)	3 (1%)	84	93
1	O	427/477 (90%)	421 (99%)	6 (1%)	67	85
1	P	428/477 (90%)	416 (97%)	12 (3%)	43	73
All	All	6836/7632 (90%)	6755 (99%)	81 (1%)	71	87

All (81) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	185	LYS
1	A	451	ARG
1	A	453	LEU
1	A	463	THR
1	A	486	LYS
1	A	510	GLN
1	B	204	LEU
1	C	230	LYS
1	C	425	ASN
1	C	511	GLU
1	C	560	LYS
1	C	597	GLU
1	D	328	ASN
1	D	425	ASN
1	D	455	LYS
1	D	470	ARG
1	D	523	LYS

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	D	577	ASN
1	D	585	ASP
1	E	185	LYS
1	E	451	ARG
1	E	463	THR
1	E	510	GLN
1	F	118	ILE
1	F	204	LEU
1	F	377	LYS
1	F	596	LYS
1	G	230	LYS
1	G	425	ASN
1	G	511	GLU
1	G	597	GLU
1	H	425	ASN
1	H	455	LYS
1	H	469	LYS
1	H	470	ARG
1	H	577	ASN
1	H	585	ASP
1	I	185	LYS
1	I	451	ARG
1	I	453	LEU
1	I	463	THR
1	J	204	LEU
1	J	377	LYS
1	J	596	LYS
1	K	230	LYS
1	K	342	GLU
1	K	425	ASN
1	K	511	GLU
1	K	597	GLU
1	L	328	ASN
1	L	425	ASN
1	L	455	LYS
1	L	471	GLU
1	L	523	LYS
1	L	577	ASN
1	L	585	ASP
1	M	185	LYS
1	M	451	ARG
1	M	463	THR

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type
1	M	486	LYS
1	N	204	LEU
1	N	325	ILE
1	N	596	LYS
1	O	117	VAL
1	O	230	LYS
1	O	425	ASN
1	O	511	GLU
1	O	560	LYS
1	O	597	GLU
1	P	288	LYS
1	P	328	ASN
1	P	425	ASN
1	P	455	LYS
1	P	465	GLN
1	P	471	GLU
1	P	523	LYS
1	P	577	ASN
1	P	580	LYS
1	P	583	ASP
1	P	585	ASP
1	P	590	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (113) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	180	HIS
1	A	190	GLN
1	A	207	ASN
1	A	210	HIS
1	A	366	HIS
1	A	425	ASN
1	A	510	GLN
1	B	119	ASN
1	B	207	ASN
1	B	210	HIS
1	B	235	GLN
1	B	256	GLN
1	B	322	HIS
1	B	366	HIS
1	B	425	ASN
1	B	599	ASN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	C	190	GLN
1	C	207	ASN
1	C	210	HIS
1	C	235	GLN
1	C	322	HIS
1	C	366	HIS
1	C	599	ASN
1	D	119	ASN
1	D	207	ASN
1	D	210	HIS
1	D	322	HIS
1	D	326	GLN
1	D	366	HIS
1	D	539	GLN
1	D	594	GLN
1	D	599	ASN
1	E	180	HIS
1	E	190	GLN
1	E	366	HIS
1	E	425	ASN
1	F	190	GLN
1	F	207	ASN
1	F	210	HIS
1	F	235	GLN
1	F	322	HIS
1	F	366	HIS
1	F	425	ASN
1	F	582	GLN
1	F	599	ASN
1	G	207	ASN
1	G	210	HIS
1	G	235	GLN
1	G	322	HIS
1	G	366	HIS
1	G	599	ASN
1	H	119	ASN
1	H	207	ASN
1	H	210	HIS
1	H	322	HIS
1	H	326	GLN
1	H	366	HIS
1	H	539	GLN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	H	594	GLN
1	H	599	ASN
1	I	180	HIS
1	I	190	GLN
1	I	366	HIS
1	I	425	ASN
1	J	119	ASN
1	J	207	ASN
1	J	210	HIS
1	J	235	GLN
1	J	322	HIS
1	J	366	HIS
1	J	425	ASN
1	J	599	ASN
1	K	207	ASN
1	K	210	HIS
1	K	235	GLN
1	K	322	HIS
1	K	366	HIS
1	K	599	ASN
1	L	207	ASN
1	L	210	HIS
1	L	322	HIS
1	L	326	GLN
1	L	366	HIS
1	L	539	GLN
1	L	594	GLN
1	L	599	ASN
1	M	180	HIS
1	M	190	GLN
1	M	207	ASN
1	M	210	HIS
1	M	366	HIS
1	M	425	ASN
1	N	207	ASN
1	N	210	HIS
1	N	235	GLN
1	N	322	HIS
1	N	366	HIS
1	N	425	ASN
1	N	599	ASN
1	O	207	ASN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	O	210	HIS
1	O	235	GLN
1	O	322	HIS
1	O	366	HIS
1	O	599	ASN
1	P	207	ASN
1	P	210	HIS
1	P	322	HIS
1	P	326	GLN
1	P	366	HIS
1	P	539	GLN
1	P	594	GLN
1	P	599	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

32 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	DGT	K	702	-	26,33,33	1.04	1 (3%)	32,52,52	2.10	4 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	DGT	J	702	-	26,33,33	1.01	1 (3%)	32,52,52	2.14	4 (12%)
2	DGT	I	702	-	26,33,33	1.05	1 (3%)	32,52,52	2.09	4 (12%)
2	DGT	K	701	-	26,33,33	1.03	1 (3%)	32,52,52	2.11	4 (12%)
2	DGT	A	701	-	26,33,33	1.02	1 (3%)	32,52,52	2.23	4 (12%)
2	DGT	P	702	-	26,33,33	1.06	1 (3%)	32,52,52	2.07	4 (12%)
2	DGT	J	701	-	26,33,33	1.05	1 (3%)	32,52,52	2.15	4 (12%)
2	DGT	O	702	-	26,33,33	1.01	1 (3%)	32,52,52	2.11	4 (12%)
2	DGT	G	702	-	26,33,33	1.06	1 (3%)	32,52,52	2.10	4 (12%)
2	DGT	H	701	-	26,33,33	1.09	1 (3%)	32,52,52	2.18	4 (12%)
2	DGT	F	702	-	26,33,33	1.04	1 (3%)	32,52,52	2.13	4 (12%)
2	DGT	B	701	-	26,33,33	1.03	1 (3%)	32,52,52	2.10	4 (12%)
2	DGT	M	701	-	26,33,33	1.05	1 (3%)	32,52,52	2.20	4 (12%)
2	DGT	A	703	-	26,33,33	1.05	1 (3%)	32,52,52	2.07	4 (12%)
2	DGT	C	701	-	26,33,33	0.99	1 (3%)	32,52,52	2.10	4 (12%)
2	DGT	I	701	-	26,33,33	1.02	1 (3%)	32,52,52	2.20	5 (15%)
2	DGT	G	701	-	26,33,33	1.00	1 (3%)	32,52,52	2.11	4 (12%)
2	DGT	D	701	-	26,33,33	1.08	1 (3%)	32,52,52	2.13	4 (12%)
2	DGT	E	702	-	26,33,33	1.04	1 (3%)	32,52,52	2.08	4 (12%)
2	DGT	J	703	-	26,33,33	1.07	1 (3%)	32,52,52	2.10	4 (12%)
2	DGT	E	703	-	26,33,33	1.05	1 (3%)	32,52,52	2.09	4 (12%)
2	DGT	O	703	-	26,33,33	1.07	2 (7%)	32,52,52	2.13	4 (12%)
2	DGT	C	702	-	26,33,33	1.01	1 (3%)	32,52,52	2.14	4 (12%)
2	DGT	E	701	-	26,33,33	1.02	1 (3%)	32,52,52	2.21	4 (12%)
2	DGT	N	701	-	26,33,33	1.12	1 (3%)	32,52,52	2.14	4 (12%)
2	DGT	M	702	-	26,33,33	1.04	1 (3%)	32,52,52	2.06	4 (12%)
2	DGT	O	701	-	26,33,33	1.01	1 (3%)	32,52,52	2.10	4 (12%)
2	DGT	P	701	-	26,33,33	1.06	1 (3%)	32,52,52	2.13	4 (12%)
2	DGT	A	702	-	26,33,33	1.10	2 (7%)	32,52,52	2.07	4 (12%)
2	DGT	B	702	-	26,33,33	1.06	1 (3%)	32,52,52	2.10	4 (12%)
2	DGT	F	701	-	26,33,33	1.08	1 (3%)	32,52,52	2.07	4 (12%)
2	DGT	L	701	-	26,33,33	1.04	1 (3%)	32,52,52	2.08	4 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	DGT	K	702	-	-	4/18/34/34	0/3/3/3
2	DGT	J	702	-	-	6/18/34/34	0/3/3/3
2	DGT	I	702	-	-	5/18/34/34	0/3/3/3
2	DGT	K	701	-	-	3/18/34/34	0/3/3/3
2	DGT	A	701	-	-	5/18/34/34	0/3/3/3
2	DGT	P	702	-	-	4/18/34/34	0/3/3/3
2	DGT	J	701	-	-	6/18/34/34	0/3/3/3
2	DGT	O	702	-	-	3/18/34/34	0/3/3/3
2	DGT	G	702	-	-	3/18/34/34	0/3/3/3
2	DGT	H	701	-	-	6/18/34/34	0/3/3/3
2	DGT	F	702	-	-	3/18/34/34	0/3/3/3
2	DGT	B	701	-	-	6/18/34/34	0/3/3/3
2	DGT	M	701	-	-	5/18/34/34	0/3/3/3
2	DGT	A	703	-	-	3/18/34/34	0/3/3/3
2	DGT	C	701	-	-	3/18/34/34	0/3/3/3
2	DGT	I	701	-	-	5/18/34/34	0/3/3/3
2	DGT	G	701	-	-	6/18/34/34	0/3/3/3
2	DGT	D	701	-	-	6/18/34/34	0/3/3/3
2	DGT	E	702	-	-	5/18/34/34	0/3/3/3
2	DGT	J	703	-	-	5/18/34/34	0/3/3/3
2	DGT	E	703	-	-	5/18/34/34	0/3/3/3
2	DGT	O	703	-	-	4/18/34/34	0/3/3/3
2	DGT	C	702	-	-	4/18/34/34	0/3/3/3
2	DGT	E	701	-	-	5/18/34/34	0/3/3/3
2	DGT	N	701	-	-	5/18/34/34	0/3/3/3
2	DGT	M	702	-	-	5/18/34/34	0/3/3/3
2	DGT	O	701	-	-	6/18/34/34	0/3/3/3
2	DGT	P	701	-	-	6/18/34/34	0/3/3/3
2	DGT	A	702	-	-	4/18/34/34	0/3/3/3
2	DGT	B	702	-	-	5/18/34/34	0/3/3/3
2	DGT	F	701	-	-	5/18/34/34	0/3/3/3
2	DGT	L	701	-	-	5/18/34/34	0/3/3/3

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	701	DGT	C6-N1	4.24	1.40	1.33
2	F	701	DGT	C6-N1	4.19	1.40	1.33
2	N	701	DGT	C6-N1	4.18	1.40	1.33
2	A	702	DGT	C6-N1	4.13	1.40	1.33
2	J	703	DGT	C6-N1	4.10	1.40	1.33
2	F	702	DGT	C6-N1	4.10	1.40	1.33
2	J	701	DGT	C6-N1	4.07	1.40	1.33
2	I	702	DGT	C6-N1	4.00	1.40	1.33
2	E	703	DGT	C6-N1	3.98	1.40	1.33
2	B	702	DGT	C6-N1	3.97	1.40	1.33
2	E	702	DGT	C6-N1	3.97	1.39	1.33
2	D	701	DGT	C6-N1	3.97	1.39	1.33
2	B	701	DGT	C6-N1	3.95	1.39	1.33
2	K	701	DGT	C6-N1	3.95	1.39	1.33
2	P	702	DGT	C6-N1	3.94	1.39	1.33
2	A	703	DGT	C6-N1	3.93	1.39	1.33
2	M	701	DGT	C6-N1	3.93	1.39	1.33
2	P	701	DGT	C6-N1	3.93	1.39	1.33
2	I	701	DGT	C6-N1	3.91	1.39	1.33
2	M	702	DGT	C6-N1	3.90	1.39	1.33
2	G	702	DGT	C6-N1	3.90	1.39	1.33
2	O	703	DGT	C6-N1	3.89	1.39	1.33
2	O	702	DGT	C6-N1	3.86	1.39	1.33
2	O	701	DGT	C6-N1	3.85	1.39	1.33
2	G	701	DGT	C6-N1	3.85	1.39	1.33
2	L	701	DGT	C6-N1	3.83	1.39	1.33
2	K	702	DGT	C6-N1	3.82	1.39	1.33
2	J	702	DGT	C6-N1	3.82	1.39	1.33
2	E	701	DGT	C6-N1	3.77	1.39	1.33
2	A	701	DGT	C6-N1	3.75	1.39	1.33
2	C	702	DGT	C6-N1	3.72	1.39	1.33
2	C	701	DGT	C6-N1	3.70	1.39	1.33
2	O	703	DGT	C8-N7	-2.08	1.31	1.34
2	A	702	DGT	C8-N7	-2.01	1.31	1.34

All (129) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	701	DGT	C5-C6-N1	-9.01	111.11	123.43
2	E	701	DGT	C5-C6-N1	-8.97	111.16	123.43
2	N	701	DGT	C5-C6-N1	-8.96	111.18	123.43
2	J	702	DGT	C5-C6-N1	-8.93	111.21	123.43
2	H	701	DGT	C5-C6-N1	-8.93	111.22	123.43

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	701	DGT	C5-C6-N1	-8.91	111.25	123.43
2	I	701	DGT	C5-C6-N1	-8.88	111.28	123.43
2	D	701	DGT	C5-C6-N1	-8.87	111.30	123.43
2	J	701	DGT	C5-C6-N1	-8.87	111.30	123.43
2	B	702	DGT	C5-C6-N1	-8.81	111.38	123.43
2	E	703	DGT	C5-C6-N1	-8.76	111.44	123.43
2	J	703	DGT	C5-C6-N1	-8.76	111.45	123.43
2	G	701	DGT	C5-C6-N1	-8.76	111.45	123.43
2	A	702	DGT	C5-C6-N1	-8.76	111.45	123.43
2	F	702	DGT	C5-C6-N1	-8.73	111.50	123.43
2	K	701	DGT	C5-C6-N1	-8.71	111.51	123.43
2	K	702	DGT	C5-C6-N1	-8.71	111.52	123.43
2	I	702	DGT	C5-C6-N1	-8.71	111.52	123.43
2	P	702	DGT	C5-C6-N1	-8.69	111.54	123.43
2	A	703	DGT	C5-C6-N1	-8.69	111.54	123.43
2	F	701	DGT	C5-C6-N1	-8.67	111.57	123.43
2	B	701	DGT	C5-C6-N1	-8.66	111.58	123.43
2	O	701	DGT	C5-C6-N1	-8.65	111.60	123.43
2	P	701	DGT	C5-C6-N1	-8.64	111.61	123.43
2	L	701	DGT	C5-C6-N1	-8.64	111.61	123.43
2	G	702	DGT	C5-C6-N1	-8.60	111.67	123.43
2	O	703	DGT	C5-C6-N1	-8.60	111.67	123.43
2	O	702	DGT	C5-C6-N1	-8.60	111.67	123.43
2	M	702	DGT	C5-C6-N1	-8.54	111.75	123.43
2	E	702	DGT	C5-C6-N1	-8.54	111.76	123.43
2	C	702	DGT	C5-C6-N1	-8.50	111.81	123.43
2	C	701	DGT	C5-C6-N1	-8.45	111.87	123.43
2	A	701	DGT	C6-N1-C2	6.26	125.87	115.93
2	E	701	DGT	C6-N1-C2	6.17	125.73	115.93
2	M	701	DGT	C6-N1-C2	6.08	125.59	115.93
2	I	701	DGT	C6-N1-C2	6.08	125.58	115.93
2	G	701	DGT	C6-N1-C2	5.94	125.37	115.93
2	J	702	DGT	C6-N1-C2	5.89	125.29	115.93
2	B	701	DGT	C6-N1-C2	5.89	125.28	115.93
2	H	701	DGT	C6-N1-C2	5.87	125.25	115.93
2	N	701	DGT	C6-N1-C2	5.86	125.25	115.93
2	J	703	DGT	C6-N1-C2	5.86	125.24	115.93
2	D	701	DGT	C6-N1-C2	5.85	125.23	115.93
2	O	701	DGT	C6-N1-C2	5.83	125.20	115.93
2	K	701	DGT	C6-N1-C2	5.83	125.19	115.93
2	J	701	DGT	C6-N1-C2	5.81	125.16	115.93
2	P	701	DGT	C6-N1-C2	5.81	125.16	115.93

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	702	DGT	C6-N1-C2	5.80	125.14	115.93
2	I	702	DGT	C6-N1-C2	5.79	125.13	115.93
2	C	702	DGT	C6-N1-C2	5.79	125.13	115.93
2	O	703	DGT	C6-N1-C2	5.78	125.11	115.93
2	F	702	DGT	C6-N1-C2	5.78	125.11	115.93
2	E	702	DGT	C6-N1-C2	5.76	125.08	115.93
2	C	701	DGT	C6-N1-C2	5.74	125.05	115.93
2	L	701	DGT	C6-N1-C2	5.74	125.04	115.93
2	F	701	DGT	C6-N1-C2	5.73	125.03	115.93
2	O	702	DGT	C6-N1-C2	5.72	125.02	115.93
2	P	702	DGT	C6-N1-C2	5.71	125.01	115.93
2	A	703	DGT	C6-N1-C2	5.71	125.01	115.93
2	K	702	DGT	C6-N1-C2	5.69	124.97	115.93
2	E	703	DGT	C6-N1-C2	5.69	124.97	115.93
2	M	702	DGT	C6-N1-C2	5.68	124.96	115.93
2	A	702	DGT	C6-N1-C2	5.58	124.80	115.93
2	G	702	DGT	C6-N1-C2	5.55	124.74	115.93
2	J	701	DGT	C2-N3-C4	-3.14	111.78	115.36
2	D	701	DGT	C2-N3-C4	-3.08	111.84	115.36
2	A	702	DGT	C2-N3-C4	-3.07	111.85	115.36
2	H	701	DGT	C2-N3-C4	-3.05	111.87	115.36
2	C	702	DGT	C2-N3-C4	-3.04	111.88	115.36
2	P	701	DGT	N3-C2-N1	-2.99	123.23	127.22
2	N	701	DGT	C2-N3-C4	-2.97	111.97	115.36
2	A	701	DGT	N3-C2-N1	-2.94	123.30	127.22
2	M	701	DGT	N3-C2-N1	-2.94	123.30	127.22
2	B	701	DGT	N3-C2-N1	-2.93	123.31	127.22
2	C	701	DGT	N3-C2-N1	-2.91	123.34	127.22
2	E	701	DGT	N3-C2-N1	-2.90	123.36	127.22
2	G	701	DGT	N3-C2-N1	-2.89	123.36	127.22
2	K	701	DGT	N3-C2-N1	-2.86	123.40	127.22
2	J	702	DGT	C2-N3-C4	-2.85	112.10	115.36
2	O	702	DGT	N3-C2-N1	-2.82	123.46	127.22
2	F	702	DGT	N3-C2-N1	-2.81	123.47	127.22
2	O	701	DGT	N3-C2-N1	-2.81	123.48	127.22
2	B	702	DGT	C2-N3-C4	-2.81	112.15	115.36
2	A	701	DGT	C2-N3-C4	-2.80	112.16	115.36
2	I	701	DGT	N3-C2-N1	-2.79	123.50	127.22
2	J	703	DGT	C2-N3-C4	-2.79	112.17	115.36
2	O	703	DGT	C2-N3-C4	-2.76	112.21	115.36
2	O	703	DGT	N3-C2-N1	-2.76	123.55	127.22
2	I	701	DGT	C2-N3-C4	-2.75	112.22	115.36

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	703	DGT	N3-C2-N1	-2.75	123.56	127.22
2	E	701	DGT	C2-N3-C4	-2.74	112.23	115.36
2	E	702	DGT	N3-C2-N1	-2.73	123.58	127.22
2	I	702	DGT	C2-N3-C4	-2.73	112.24	115.36
2	M	701	DGT	C2-N3-C4	-2.73	112.24	115.36
2	N	701	DGT	N3-C2-N1	-2.73	123.58	127.22
2	E	702	DGT	C2-N3-C4	-2.72	112.25	115.36
2	A	703	DGT	N3-C2-N1	-2.71	123.61	127.22
2	F	701	DGT	N3-C2-N1	-2.70	123.62	127.22
2	P	702	DGT	N3-C2-N1	-2.69	123.64	127.22
2	L	701	DGT	N3-C2-N1	-2.69	123.64	127.22
2	I	702	DGT	N3-C2-N1	-2.68	123.64	127.22
2	M	702	DGT	N3-C2-N1	-2.68	123.64	127.22
2	E	703	DGT	N3-C2-N1	-2.68	123.65	127.22
2	B	701	DGT	C2-N3-C4	-2.67	112.31	115.36
2	H	701	DGT	N3-C2-N1	-2.67	123.67	127.22
2	M	702	DGT	C2-N3-C4	-2.66	112.32	115.36
2	D	701	DGT	N3-C2-N1	-2.65	123.68	127.22
2	G	702	DGT	C2-N3-C4	-2.64	112.34	115.36
2	C	702	DGT	N3-C2-N1	-2.63	123.71	127.22
2	P	702	DGT	C2-N3-C4	-2.63	112.35	115.36
2	K	702	DGT	C2-N3-C4	-2.62	112.37	115.36
2	K	702	DGT	N3-C2-N1	-2.61	123.74	127.22
2	J	702	DGT	N3-C2-N1	-2.61	123.74	127.22
2	F	701	DGT	C2-N3-C4	-2.61	112.38	115.36
2	B	702	DGT	N3-C2-N1	-2.60	123.75	127.22
2	L	701	DGT	C2-N3-C4	-2.60	112.39	115.36
2	P	701	DGT	C2-N3-C4	-2.54	112.45	115.36
2	O	701	DGT	C2-N3-C4	-2.53	112.46	115.36
2	K	701	DGT	C2-N3-C4	-2.51	112.49	115.36
2	F	702	DGT	C2-N3-C4	-2.50	112.50	115.36
2	J	701	DGT	N3-C2-N1	-2.50	123.89	127.22
2	G	701	DGT	C2-N3-C4	-2.48	112.52	115.36
2	C	701	DGT	C2-N3-C4	-2.48	112.52	115.36
2	A	702	DGT	N3-C2-N1	-2.45	123.95	127.22
2	G	702	DGT	N3-C2-N1	-2.45	123.96	127.22
2	O	702	DGT	C2-N3-C4	-2.42	112.60	115.36
2	E	703	DGT	C2-N3-C4	-2.40	112.61	115.36
2	A	703	DGT	C2-N3-C4	-2.34	112.68	115.36
2	I	701	DGT	PB-O3B-PG	-2.04	125.81	132.83

There are no chirality outliers.

All (151) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	701	DGT	C5'-O5'-PA-O3A
2	A	701	DGT	C5'-O5'-PA-O2A
2	A	702	DGT	PB-O3B-PG-O1G
2	B	701	DGT	C5'-O5'-PA-O3A
2	B	701	DGT	C5'-O5'-PA-O1A
2	B	701	DGT	C5'-O5'-PA-O2A
2	B	702	DGT	C5'-O5'-PA-O1A
2	B	702	DGT	C5'-O5'-PA-O2A
2	B	702	DGT	O4'-C4'-C5'-O5'
2	C	701	DGT	C3'-C4'-C5'-O5'
2	C	702	DGT	O4'-C4'-C5'-O5'
2	D	701	DGT	C5'-O5'-PA-O3A
2	D	701	DGT	C5'-O5'-PA-O2A
2	D	701	DGT	C4'-C5'-O5'-PA
2	D	701	DGT	C3'-C4'-C5'-O5'
2	E	701	DGT	C5'-O5'-PA-O3A
2	E	701	DGT	C5'-O5'-PA-O2A
2	E	702	DGT	PB-O3B-PG-O1G
2	F	701	DGT	C5'-O5'-PA-O1A
2	F	701	DGT	C5'-O5'-PA-O2A
2	F	701	DGT	O4'-C4'-C5'-O5'
2	F	702	DGT	C3'-C4'-C5'-O5'
2	G	701	DGT	C5'-O5'-PA-O3A
2	G	701	DGT	C5'-O5'-PA-O1A
2	G	701	DGT	C5'-O5'-PA-O2A
2	G	702	DGT	O4'-C4'-C5'-O5'
2	H	701	DGT	C5'-O5'-PA-O2A
2	H	701	DGT	C3'-C4'-C5'-O5'
2	I	701	DGT	C5'-O5'-PA-O3A
2	I	701	DGT	C5'-O5'-PA-O2A
2	I	702	DGT	PB-O3B-PG-O1G
2	J	701	DGT	PB-O3A-PA-O5'
2	J	701	DGT	C5'-O5'-PA-O2A
2	J	701	DGT	C4'-C5'-O5'-PA
2	J	701	DGT	C3'-C4'-C5'-O5'
2	J	702	DGT	C5'-O5'-PA-O1A
2	J	702	DGT	C5'-O5'-PA-O2A
2	J	703	DGT	C5'-O5'-PA-O1A
2	J	703	DGT	C5'-O5'-PA-O2A
2	J	703	DGT	O4'-C4'-C5'-O5'
2	K	701	DGT	C3'-C4'-C5'-O5'
2	K	702	DGT	O4'-C4'-C5'-O5'

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms
2	L	701	DGT	O4'-C4'-C5'-O5'
2	M	701	DGT	C5'-O5'-PA-O3A
2	M	701	DGT	C5'-O5'-PA-O2A
2	M	702	DGT	PB-O3B-PG-O1G
2	N	701	DGT	C5'-O5'-PA-O1A
2	N	701	DGT	C5'-O5'-PA-O2A
2	N	701	DGT	O4'-C4'-C5'-O5'
2	O	701	DGT	C5'-O5'-PA-O3A
2	O	701	DGT	C5'-O5'-PA-O1A
2	O	701	DGT	C5'-O5'-PA-O2A
2	O	702	DGT	C3'-C4'-C5'-O5'
2	O	703	DGT	O4'-C4'-C5'-O5'
2	P	701	DGT	PB-O3A-PA-O5'
2	P	701	DGT	C5'-O5'-PA-O2A
2	P	701	DGT	C4'-C5'-O5'-PA
2	P	701	DGT	C3'-C4'-C5'-O5'
2	A	702	DGT	O4'-C4'-C5'-O5'
2	E	702	DGT	O4'-C4'-C5'-O5'
2	E	703	DGT	O4'-C4'-C5'-O5'
2	I	702	DGT	O4'-C4'-C5'-O5'
2	M	702	DGT	O4'-C4'-C5'-O5'
2	P	702	DGT	O4'-C4'-C5'-O5'
2	H	701	DGT	C4'-C5'-O5'-PA
2	B	702	DGT	C3'-C4'-C5'-O5'
2	C	701	DGT	O4'-C4'-C5'-O5'
2	C	702	DGT	C3'-C4'-C5'-O5'
2	D	701	DGT	O4'-C4'-C5'-O5'
2	F	701	DGT	C3'-C4'-C5'-O5'
2	F	702	DGT	O4'-C4'-C5'-O5'
2	G	702	DGT	C3'-C4'-C5'-O5'
2	H	701	DGT	O4'-C4'-C5'-O5'
2	J	701	DGT	O4'-C4'-C5'-O5'
2	J	703	DGT	C3'-C4'-C5'-O5'
2	K	701	DGT	O4'-C4'-C5'-O5'
2	K	702	DGT	C3'-C4'-C5'-O5'
2	N	701	DGT	C3'-C4'-C5'-O5'
2	O	702	DGT	O4'-C4'-C5'-O5'
2	O	703	DGT	C3'-C4'-C5'-O5'
2	P	701	DGT	O4'-C4'-C5'-O5'
2	A	702	DGT	C3'-C4'-C5'-O5'
2	A	701	DGT	O4'-C4'-C5'-O5'
2	E	701	DGT	O4'-C4'-C5'-O5'

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms
2	M	701	DGT	O4'-C4'-C5'-O5'
2	I	701	DGT	O4'-C4'-C5'-O5'
2	I	702	DGT	C3'-C4'-C5'-O5'
2	C	702	DGT	PB-O3B-PG-O3G
2	K	702	DGT	PB-O3B-PG-O3G
2	E	703	DGT	PG-O3B-PB-O2B
2	P	702	DGT	PG-O3B-PB-O2B
2	E	702	DGT	C3'-C4'-C5'-O5'
2	E	703	DGT	C3'-C4'-C5'-O5'
2	M	702	DGT	C3'-C4'-C5'-O5'
2	D	701	DGT	PB-O3A-PA-O5'
2	H	701	DGT	PB-O3A-PA-O5'
2	O	703	DGT	PB-O3B-PG-O3G
2	H	701	DGT	C5'-O5'-PA-O3A
2	J	701	DGT	C5'-O5'-PA-O3A
2	J	702	DGT	C5'-O5'-PA-O3A
2	P	701	DGT	C5'-O5'-PA-O3A
2	A	701	DGT	C5'-O5'-PA-O1A
2	E	701	DGT	C5'-O5'-PA-O1A
2	I	701	DGT	C5'-O5'-PA-O1A
2	M	701	DGT	C5'-O5'-PA-O1A
2	P	702	DGT	C3'-C4'-C5'-O5'
2	B	701	DGT	PG-O3B-PB-O1B
2	G	701	DGT	PG-O3B-PB-O1B
2	J	702	DGT	PG-O3B-PB-O1B
2	L	701	DGT	PG-O3B-PB-O2B
2	O	701	DGT	PG-O3B-PB-O1B
2	P	702	DGT	PG-O3B-PB-O1B
2	L	701	DGT	C3'-C4'-C5'-O5'
2	G	702	DGT	PB-O3B-PG-O3G
2	E	703	DGT	PA-O3A-PB-O2B
2	L	701	DGT	PA-O3A-PB-O2B
2	J	702	DGT	PA-O3A-PB-O3B
2	O	701	DGT	PA-O3A-PB-O3B
2	A	701	DGT	C3'-C4'-C5'-O5'
2	E	701	DGT	C3'-C4'-C5'-O5'
2	I	701	DGT	C3'-C4'-C5'-O5'
2	M	701	DGT	C3'-C4'-C5'-O5'
2	A	702	DGT	PB-O3B-PG-O3G
2	E	702	DGT	PB-O3B-PG-O3G
2	I	702	DGT	PB-O3B-PG-O3G
2	M	702	DGT	PB-O3B-PG-O3G

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms
2	B	701	DGT	PA-O3A-PB-O3B
2	G	701	DGT	PA-O3A-PB-O3B
2	B	702	DGT	C5'-O5'-PA-O3A
2	C	701	DGT	C5'-O5'-PA-O3A
2	F	701	DGT	C5'-O5'-PA-O3A
2	F	702	DGT	C5'-O5'-PA-O3A
2	J	703	DGT	C5'-O5'-PA-O3A
2	K	701	DGT	C5'-O5'-PA-O3A
2	N	701	DGT	C5'-O5'-PA-O3A
2	O	702	DGT	C5'-O5'-PA-O3A
2	A	703	DGT	O4'-C4'-C5'-O5'
2	A	703	DGT	PA-O3A-PB-O2B
2	B	701	DGT	PA-O3A-PB-O2B
2	E	703	DGT	PG-O3B-PB-O1B
2	G	701	DGT	PA-O3A-PB-O2B
2	J	702	DGT	PA-O3A-PB-O2B
2	L	701	DGT	PG-O3B-PB-O1B
2	O	701	DGT	PA-O3A-PB-O2B
2	A	703	DGT	C5'-O5'-PA-O2A
2	C	702	DGT	C5'-O5'-PA-O2A
2	E	702	DGT	C5'-O5'-PA-O2A
2	I	702	DGT	C5'-O5'-PA-O2A
2	K	702	DGT	C5'-O5'-PA-O2A
2	M	702	DGT	C5'-O5'-PA-O2A
2	O	703	DGT	C5'-O5'-PA-O2A

There are no ring outliers.

23 monomers are involved in 60 short contacts:

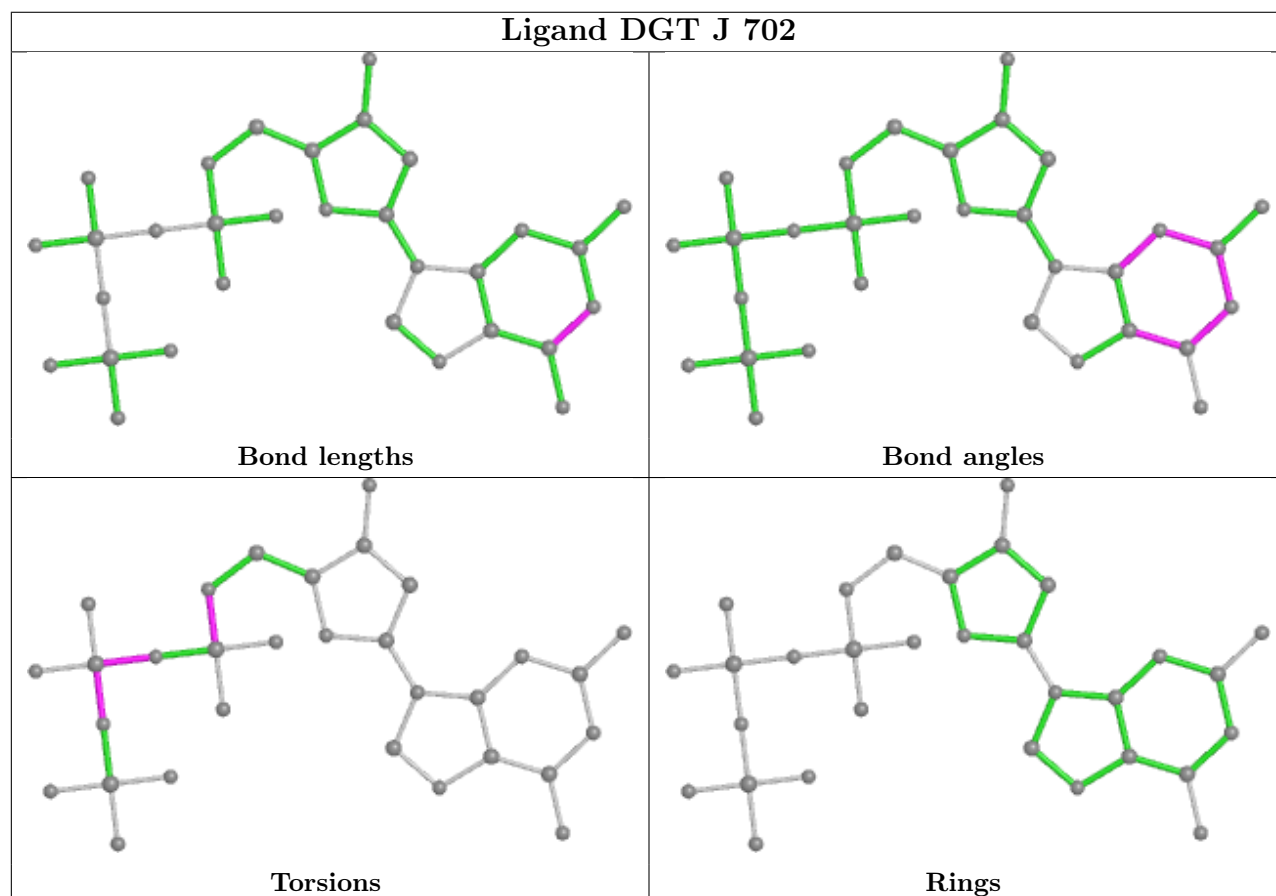
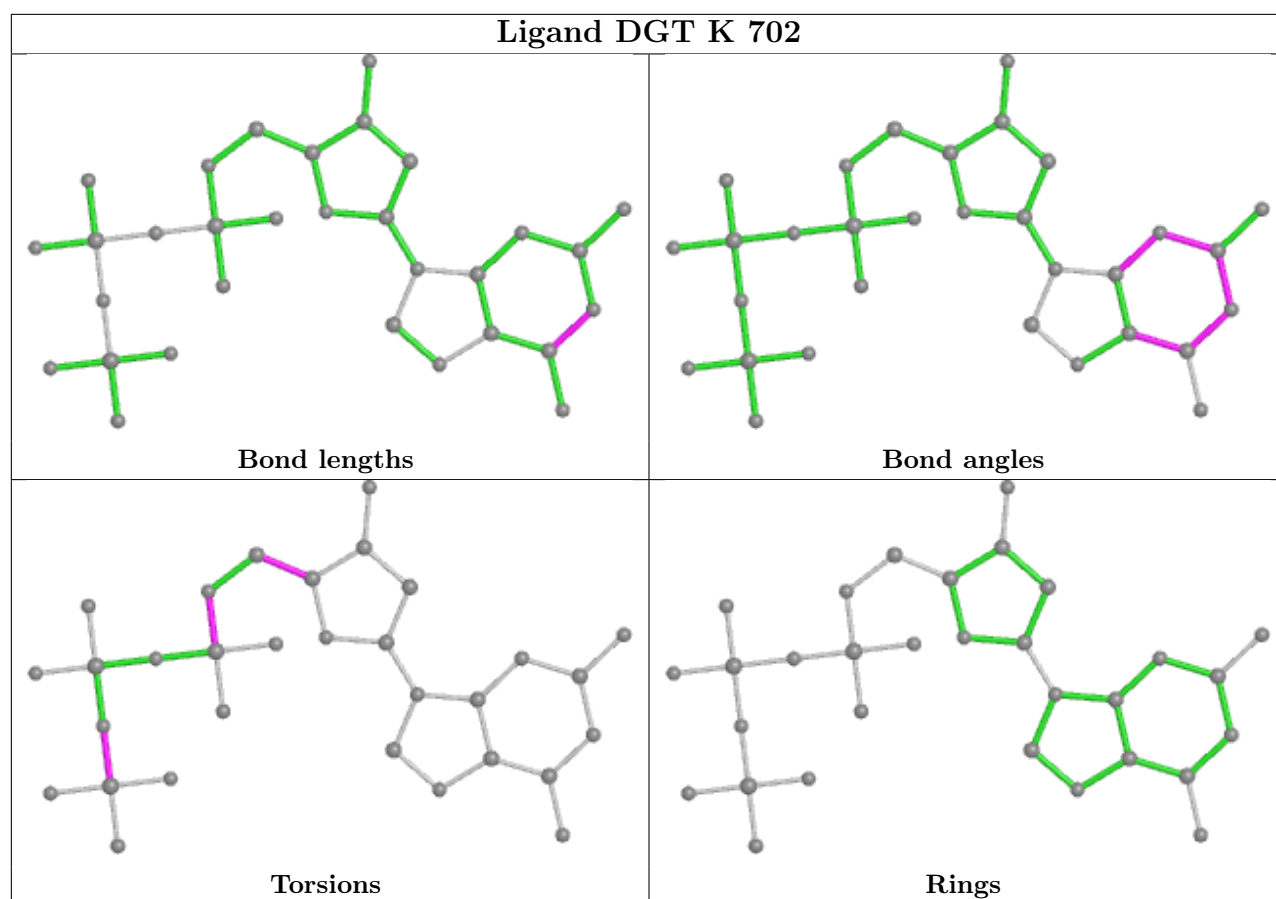
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	K	702	DGT	1	0
2	J	702	DGT	2	0
2	K	701	DGT	3	0
2	J	701	DGT	8	0
2	O	702	DGT	3	0
2	G	702	DGT	3	0
2	H	701	DGT	7	0
2	F	702	DGT	4	0
2	B	701	DGT	2	0
2	A	703	DGT	1	0
2	C	701	DGT	4	0
2	G	701	DGT	1	0

*Continued on next page...*

*Continued from previous page...*

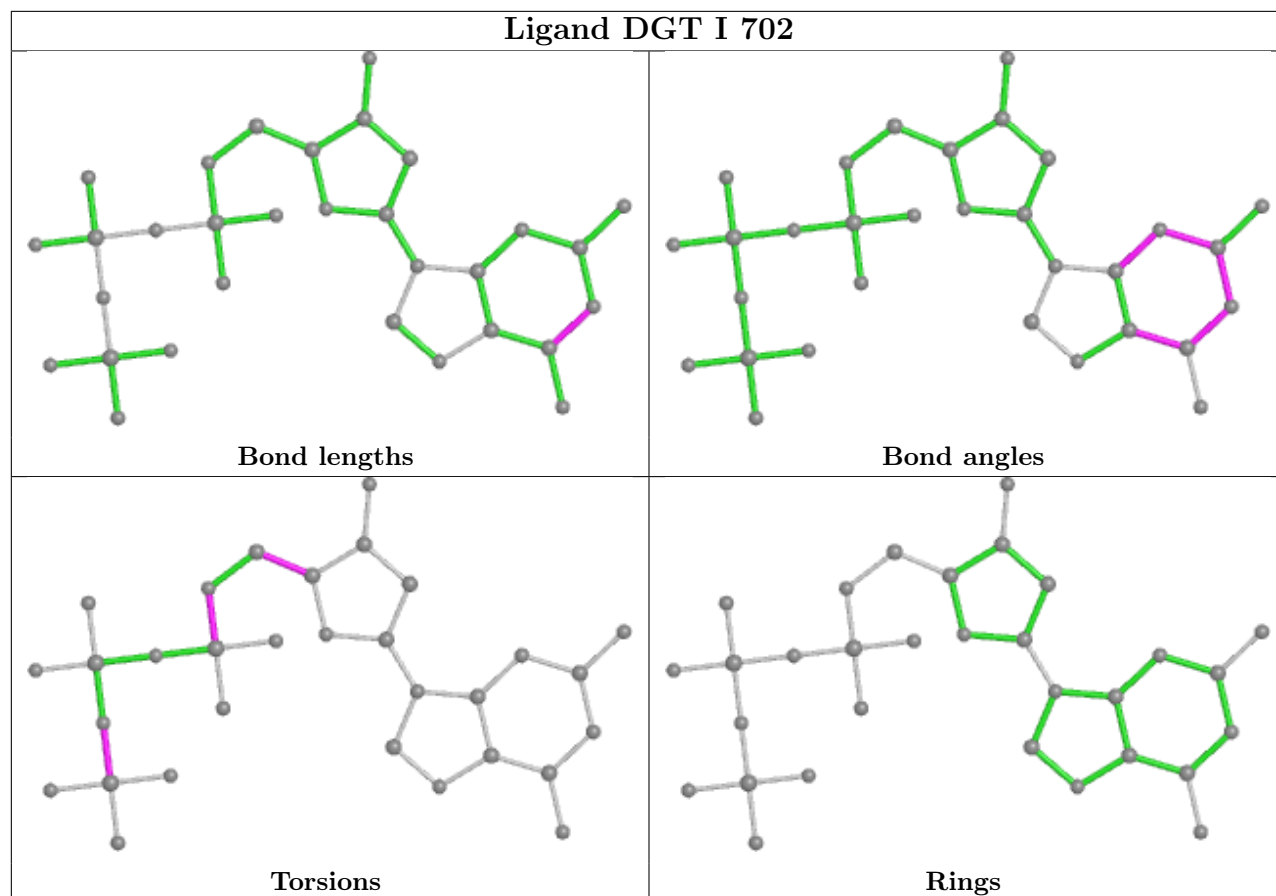
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	701	DGT	1	0
2	E	702	DGT	2	0
2	J	703	DGT	1	0
2	O	703	DGT	1	0
2	C	702	DGT	4	0
2	E	701	DGT	1	0
2	M	702	DGT	1	0
2	P	701	DGT	6	0
2	A	702	DGT	4	0
2	F	701	DGT	2	0
2	L	701	DGT	2	0

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.

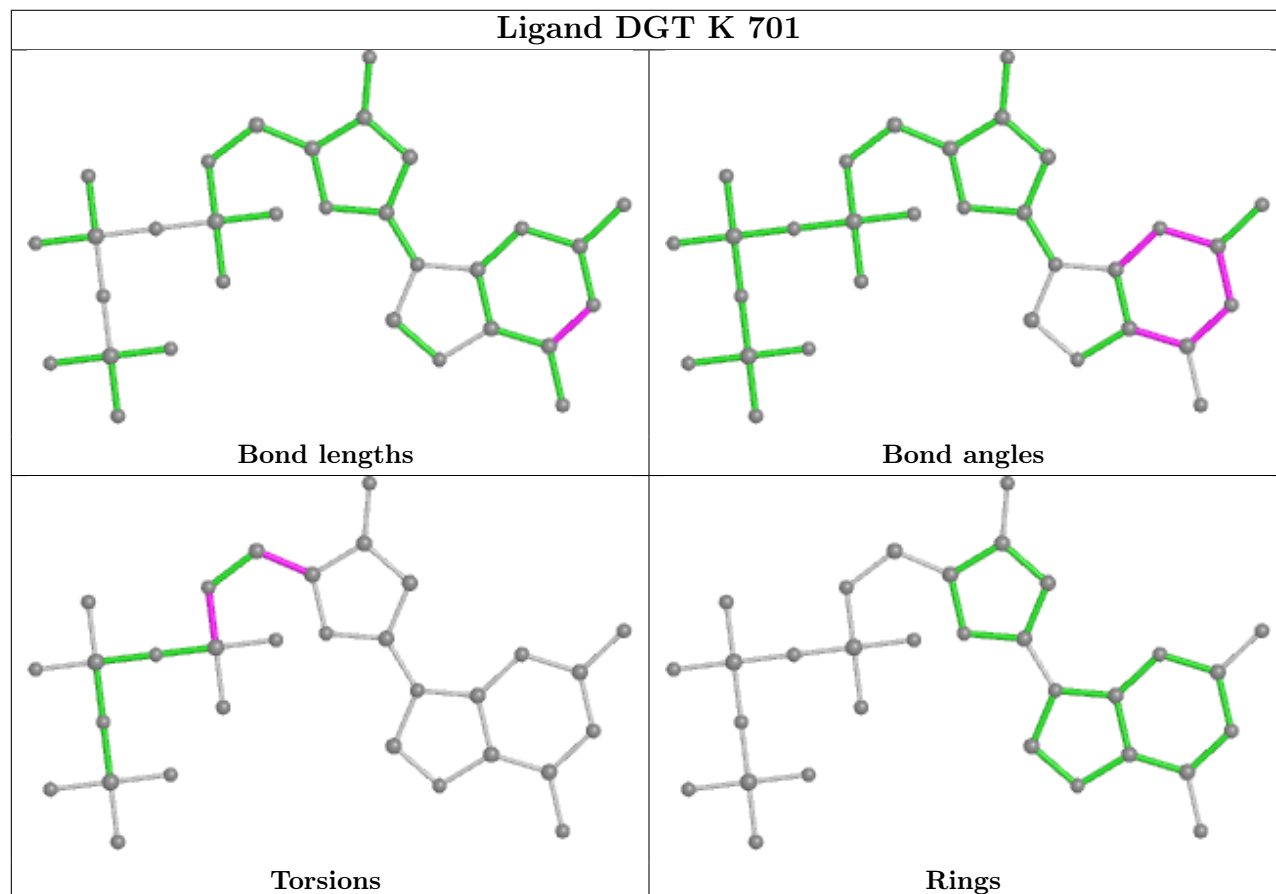


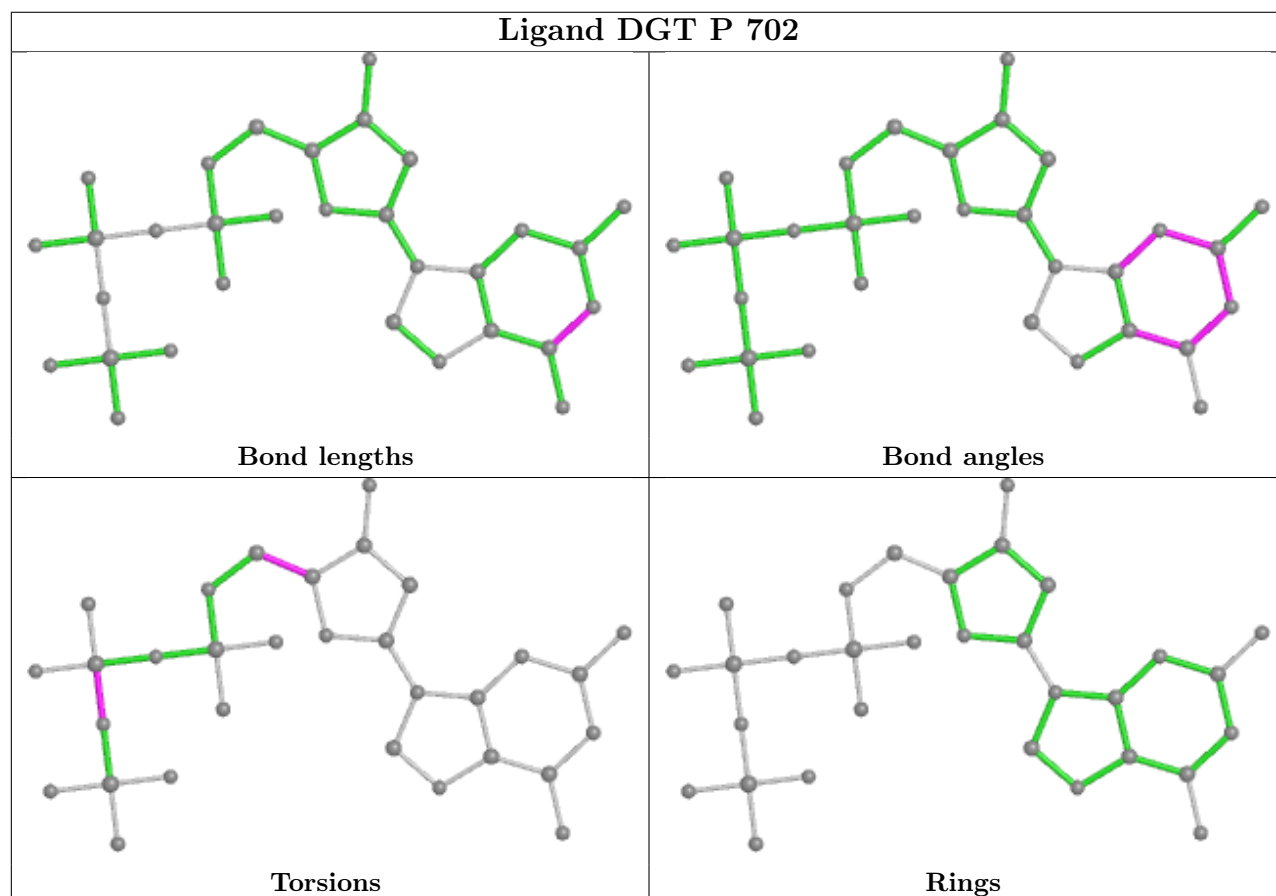
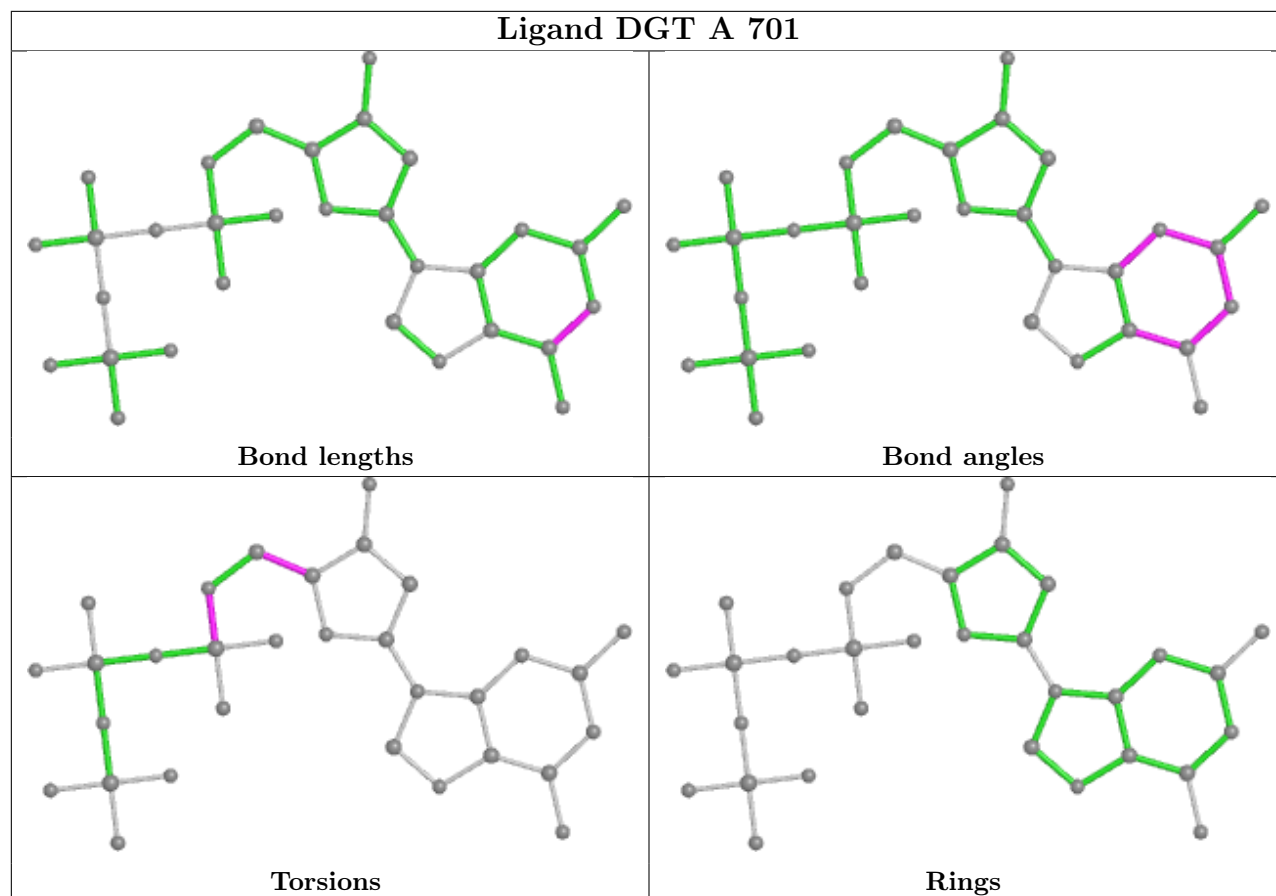


## Ligand DGT I 702

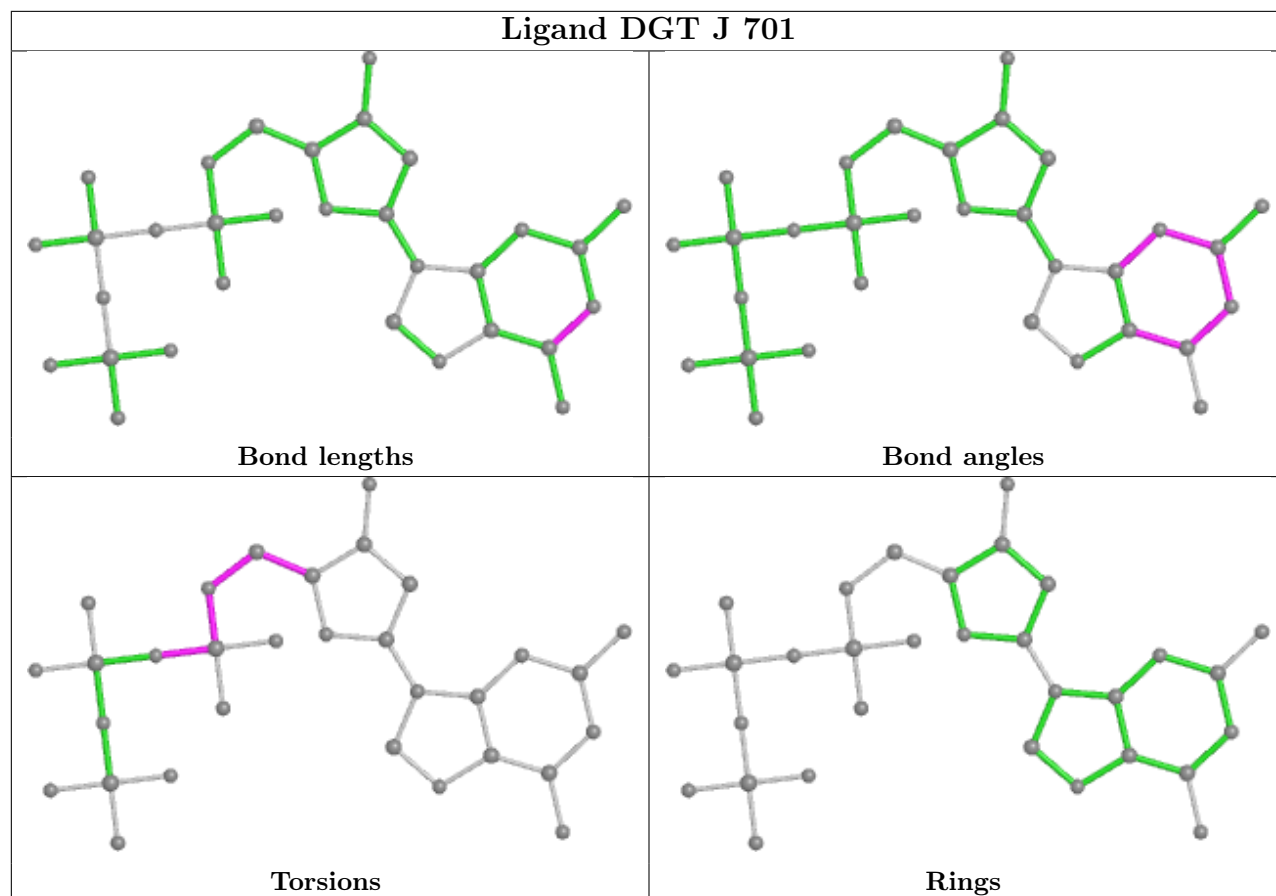


## Ligand DGT K 701

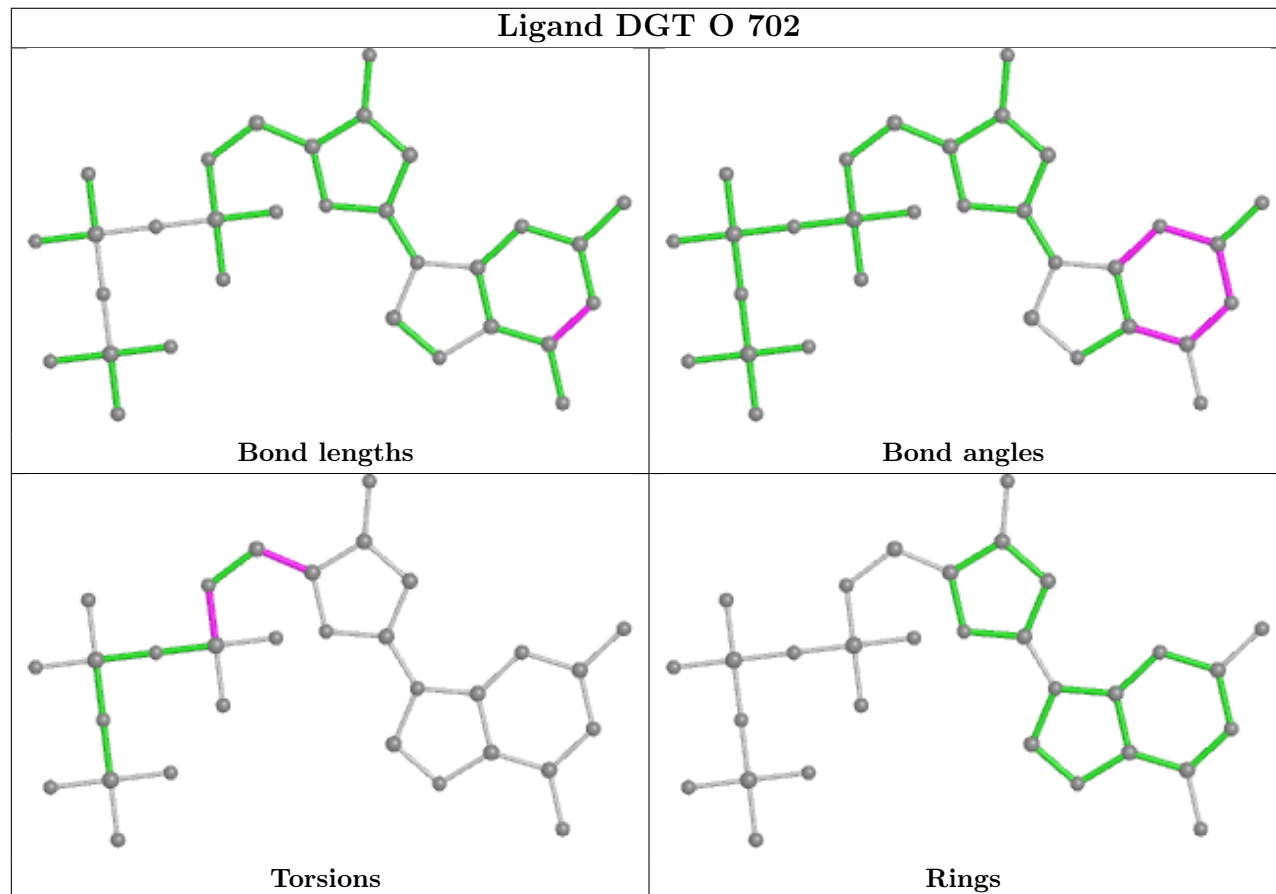


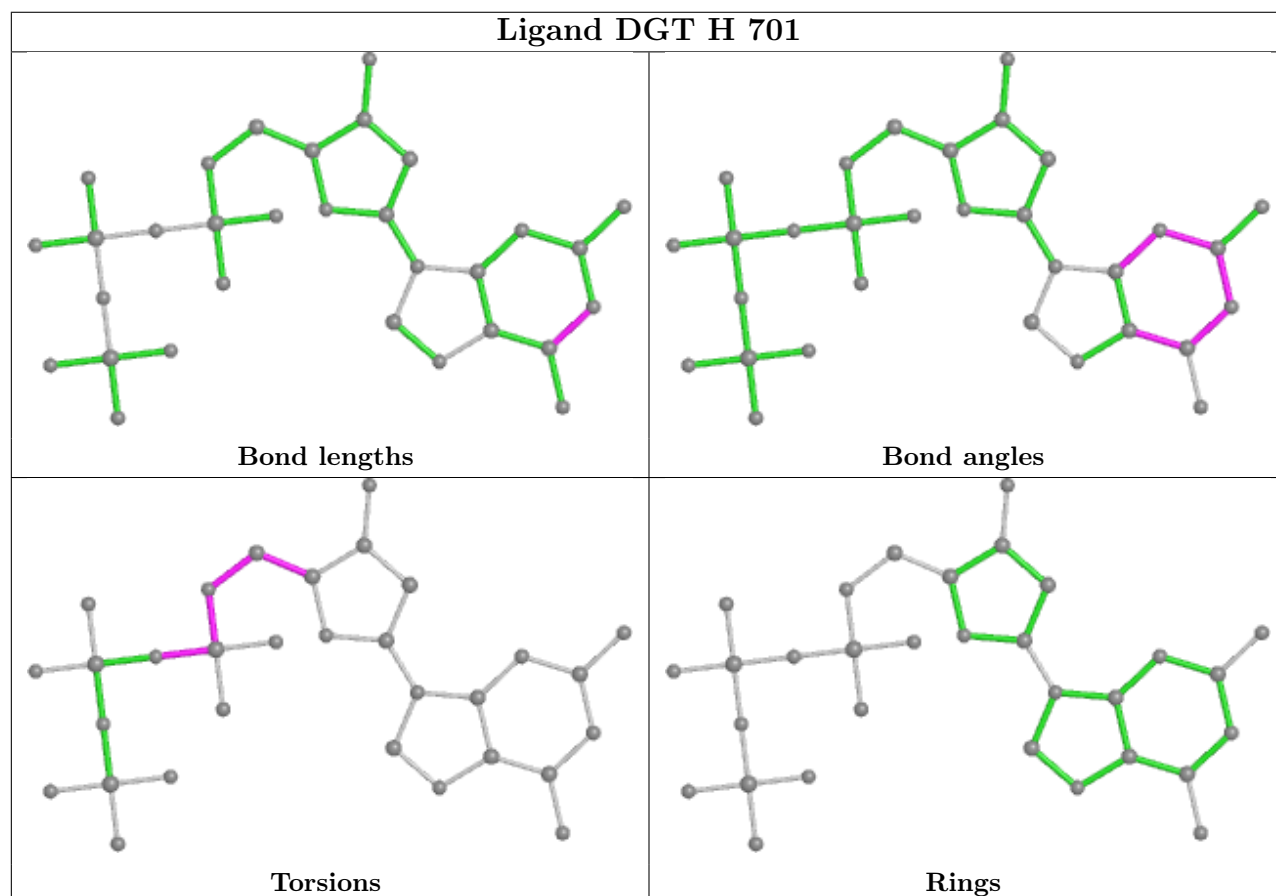
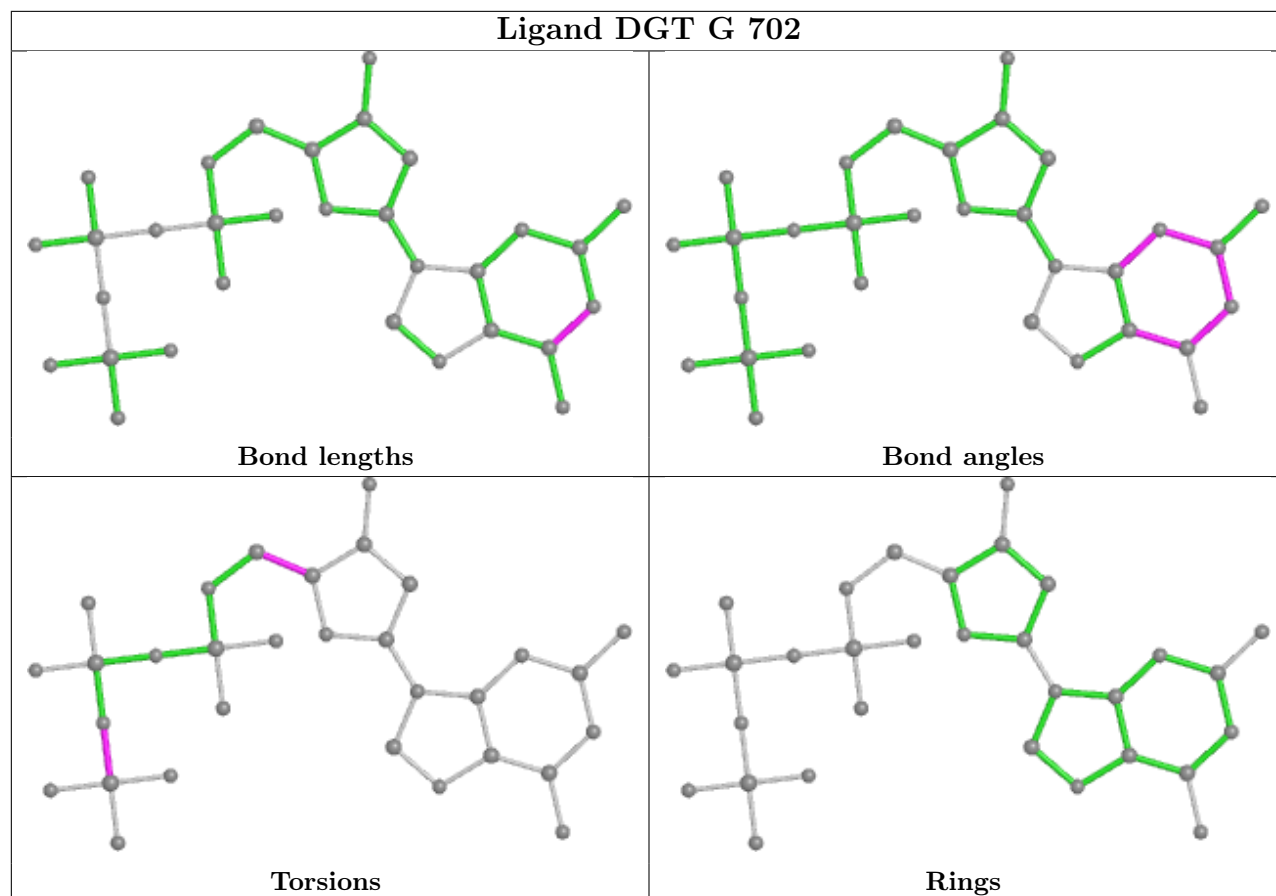


## Ligand DGT J 701

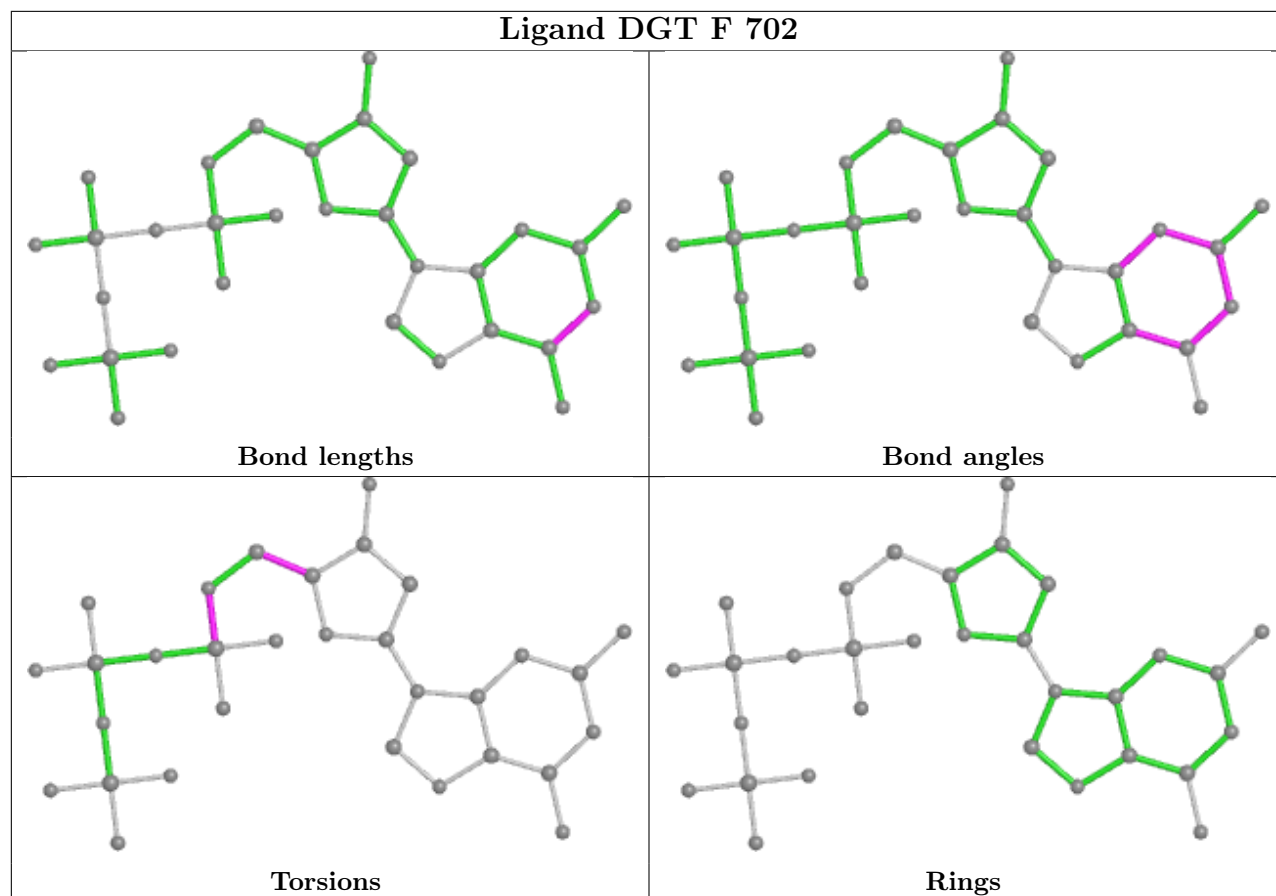


## Ligand DGT O 702

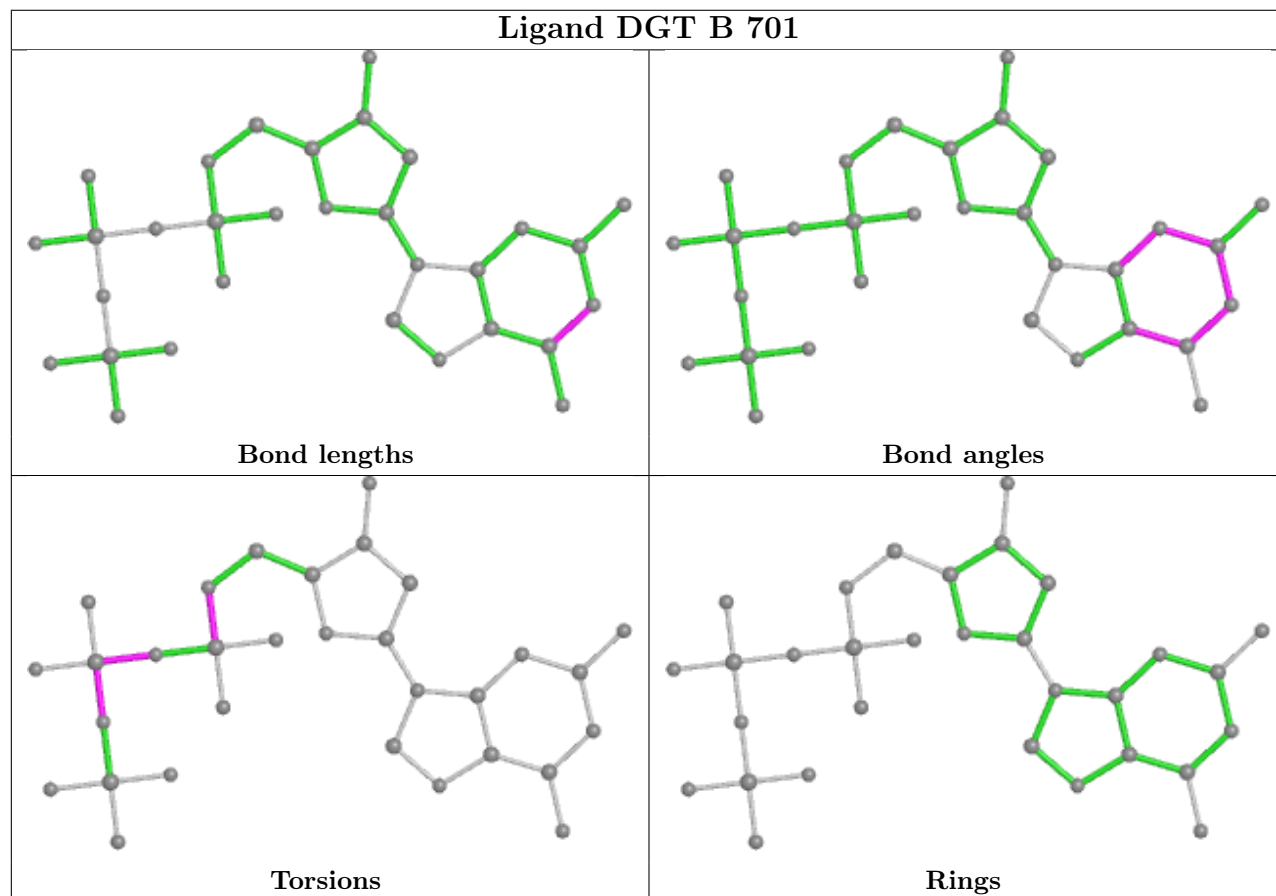




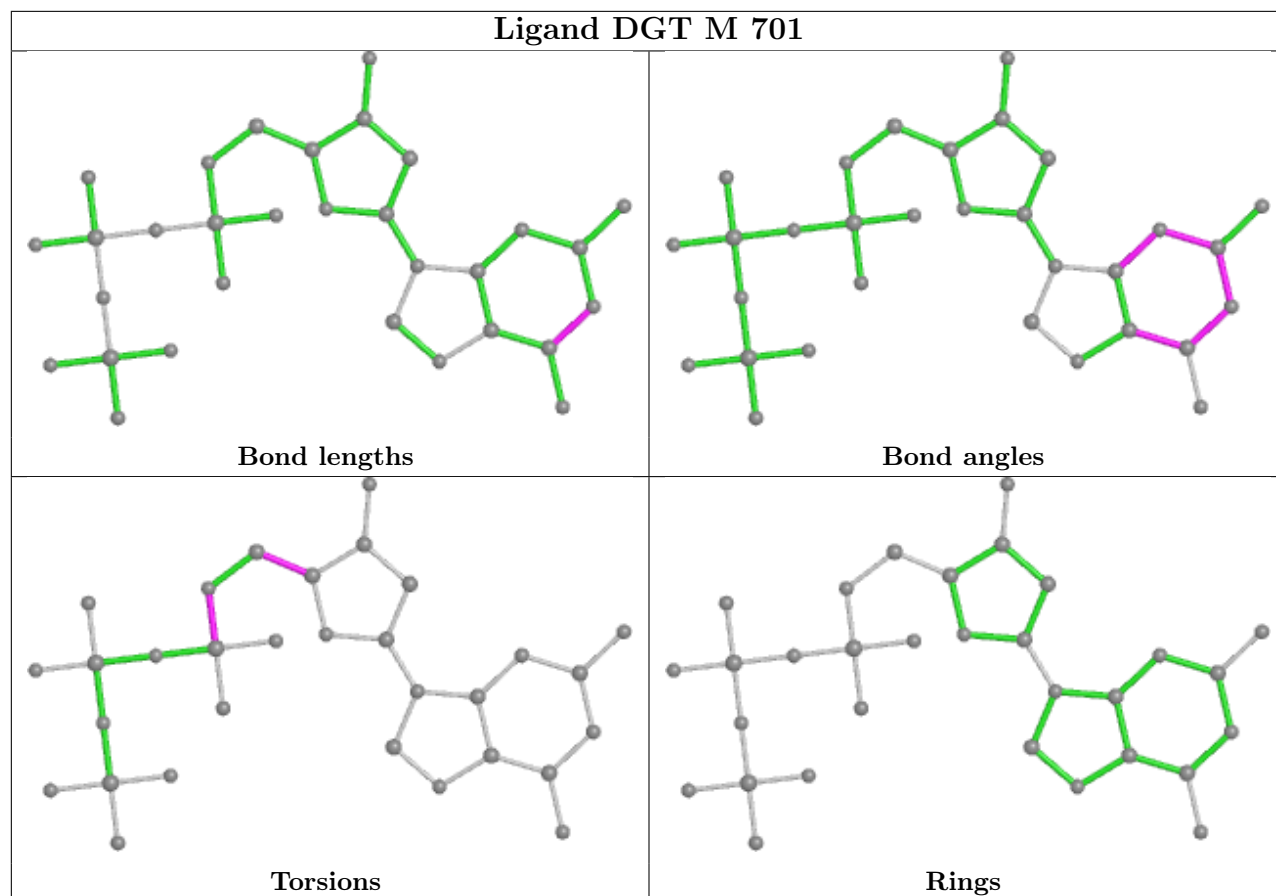
## Ligand DGT F 702



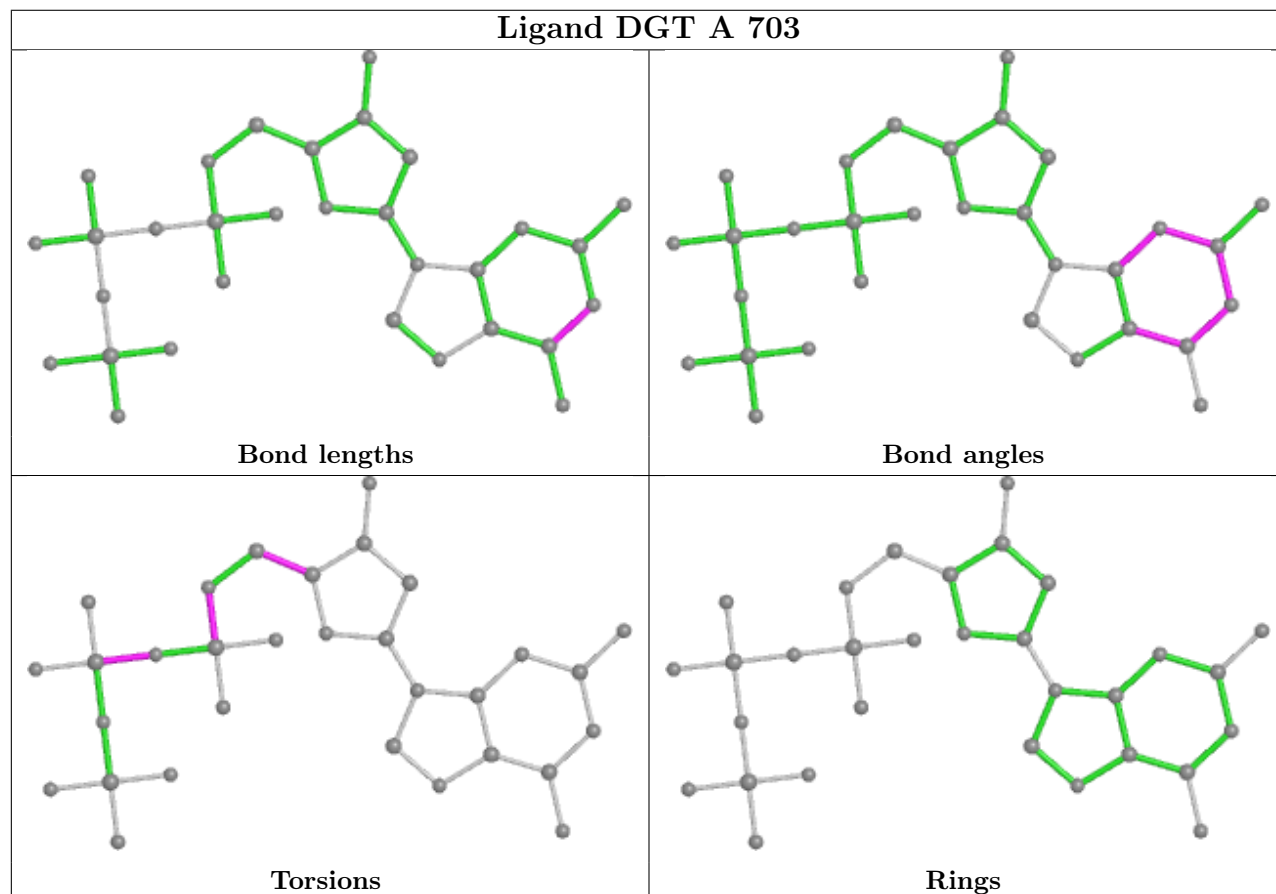
## Ligand DGT B 701

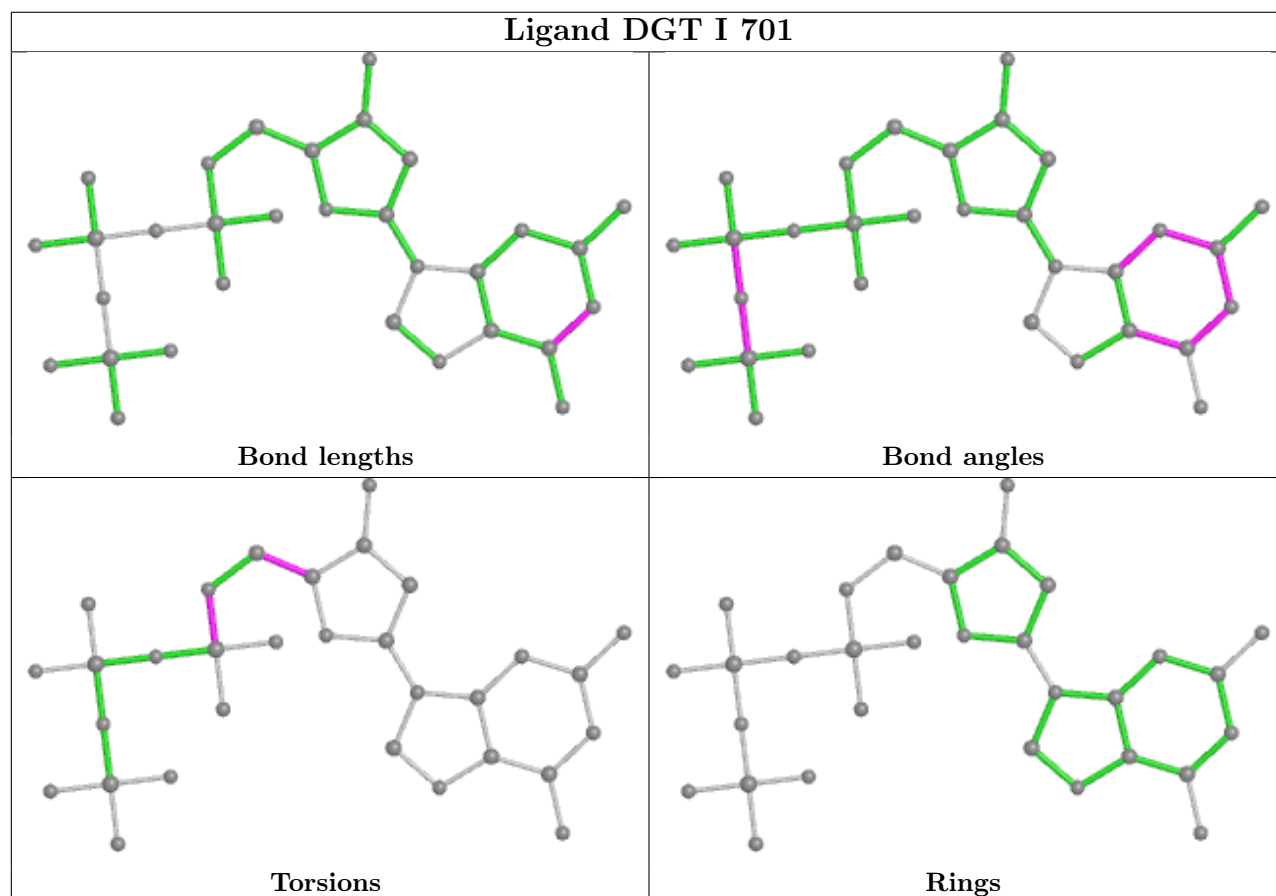
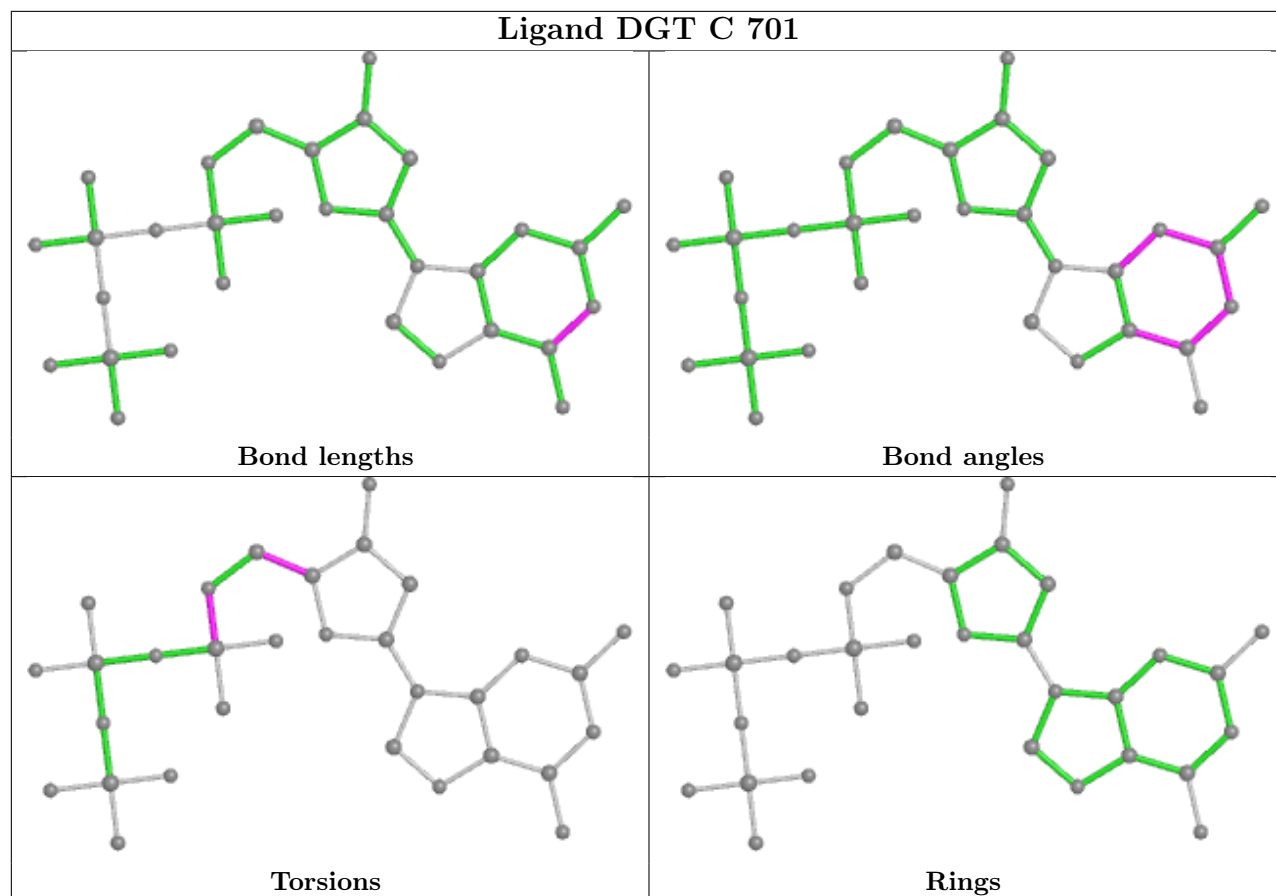


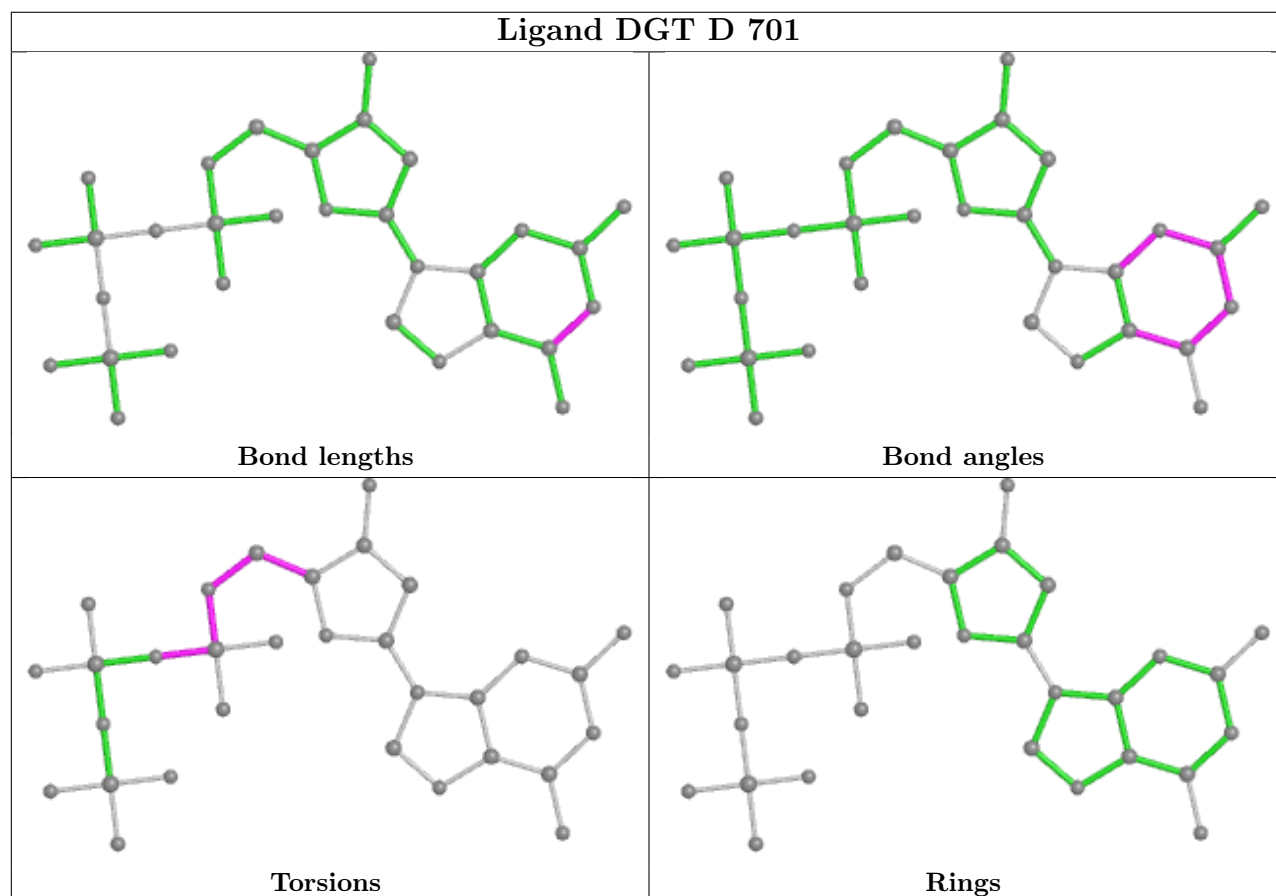
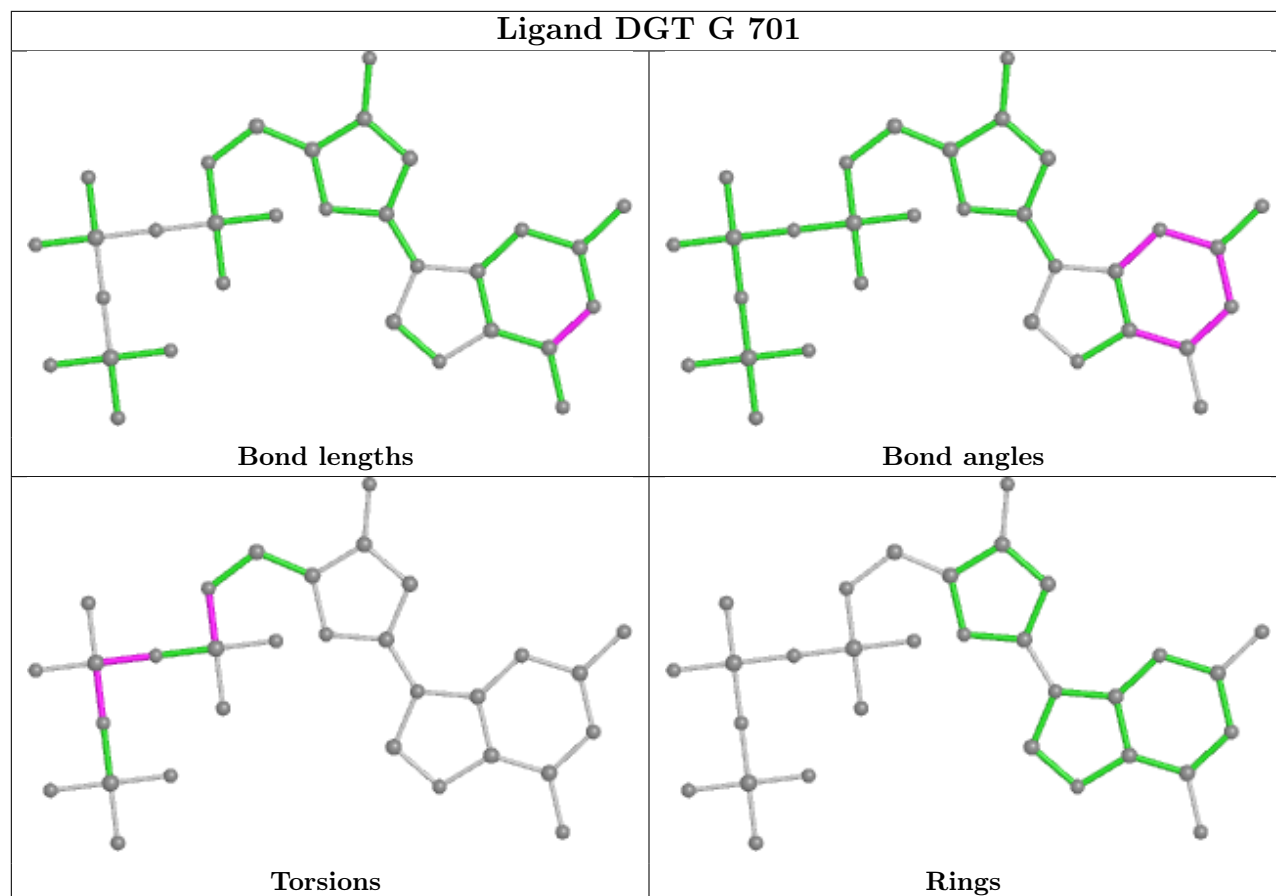
## Ligand DGT M 701



## Ligand DGT A 703

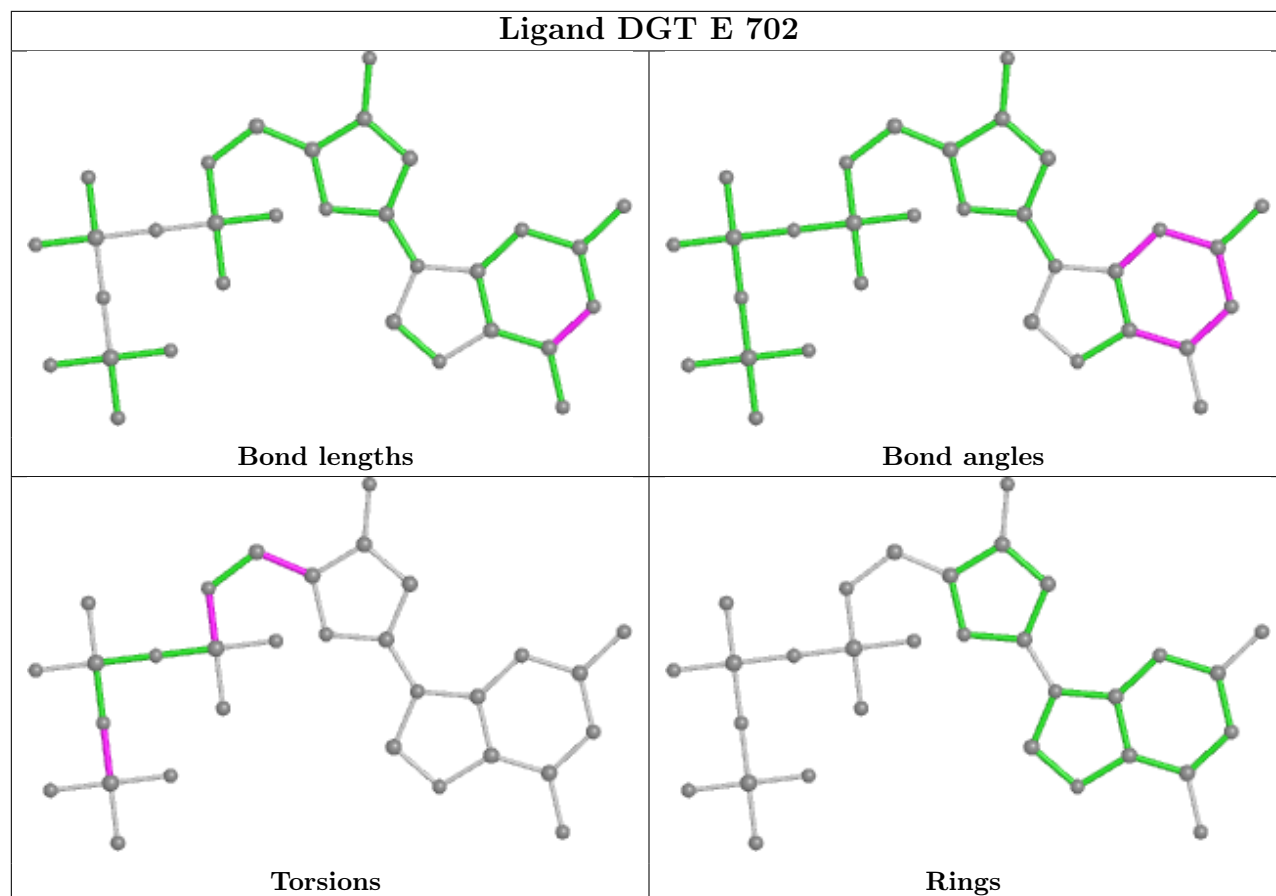




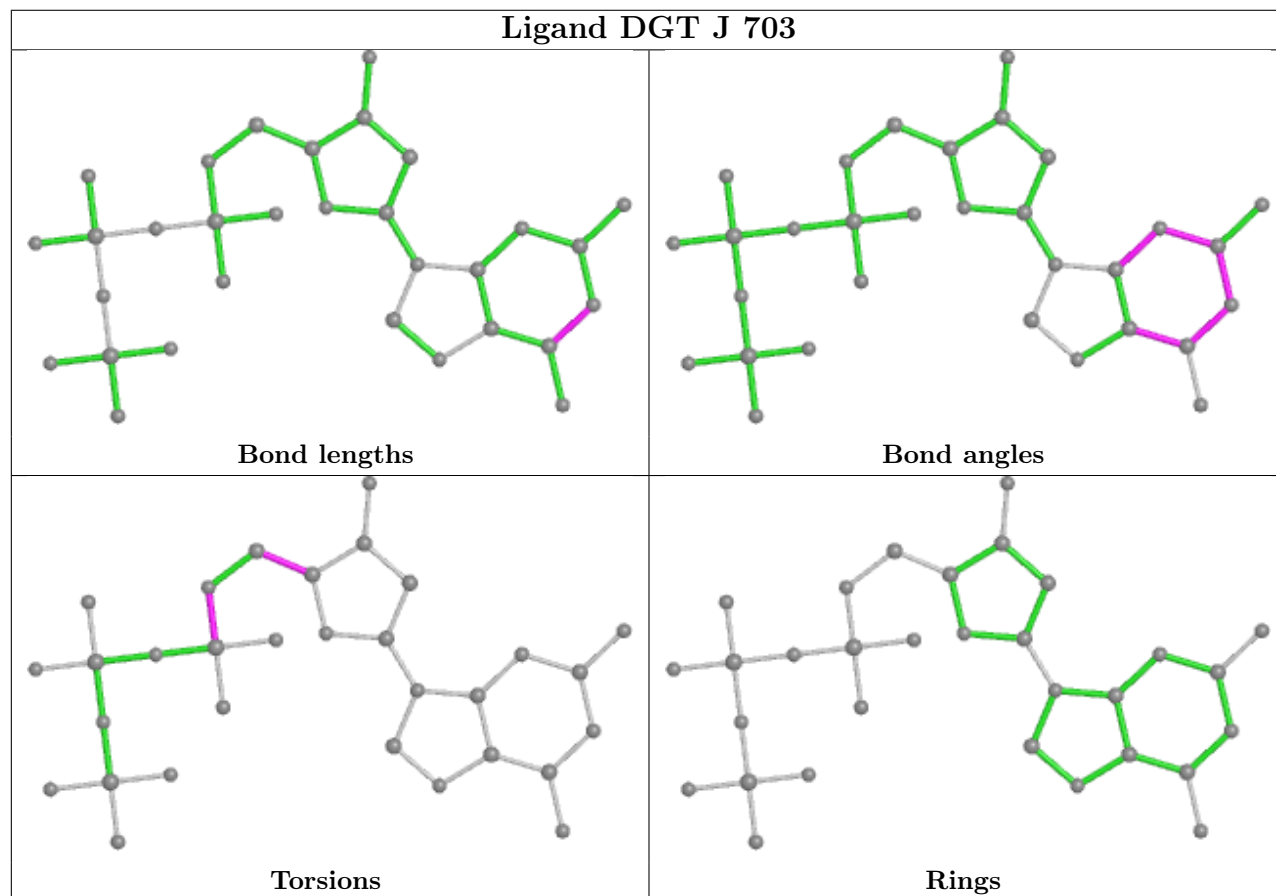




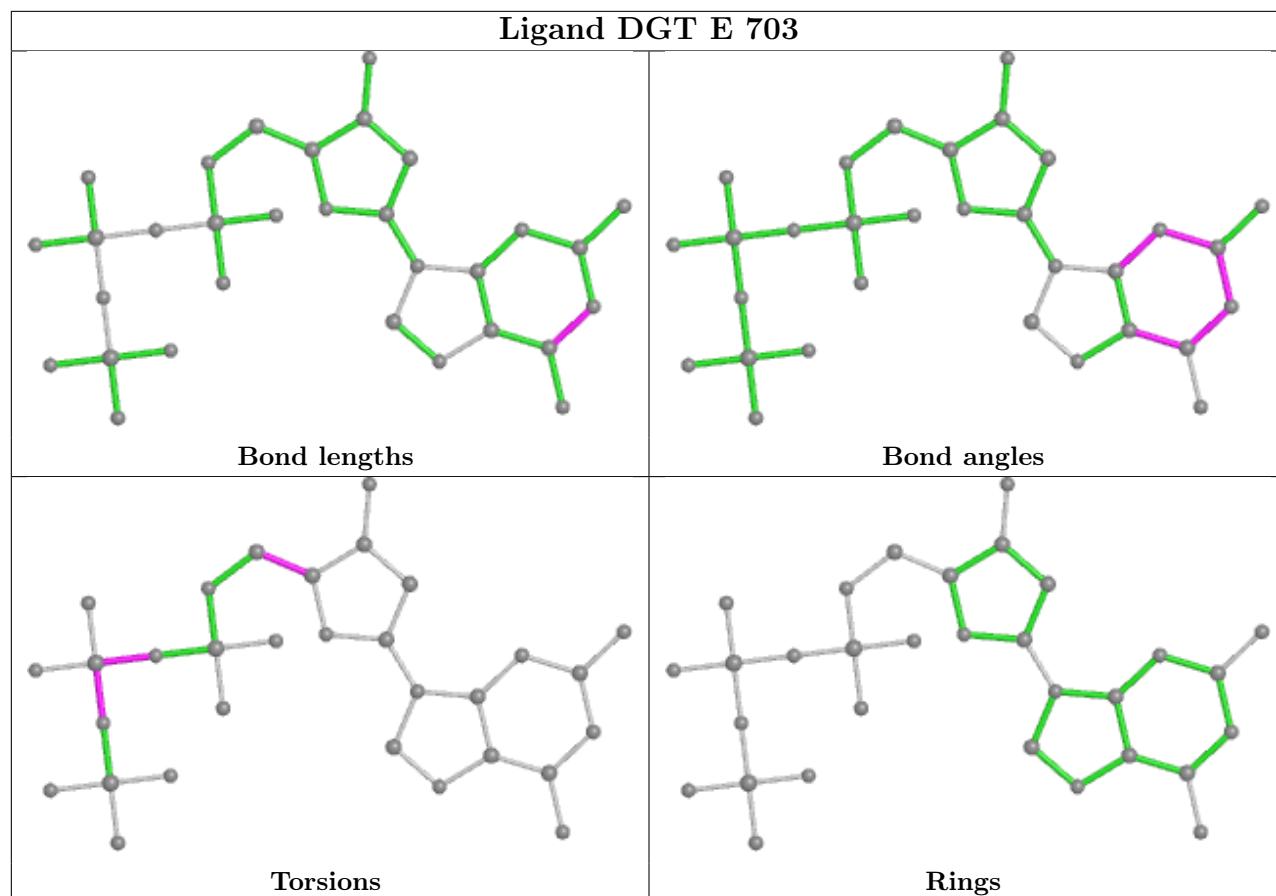
## Ligand DGT E 702



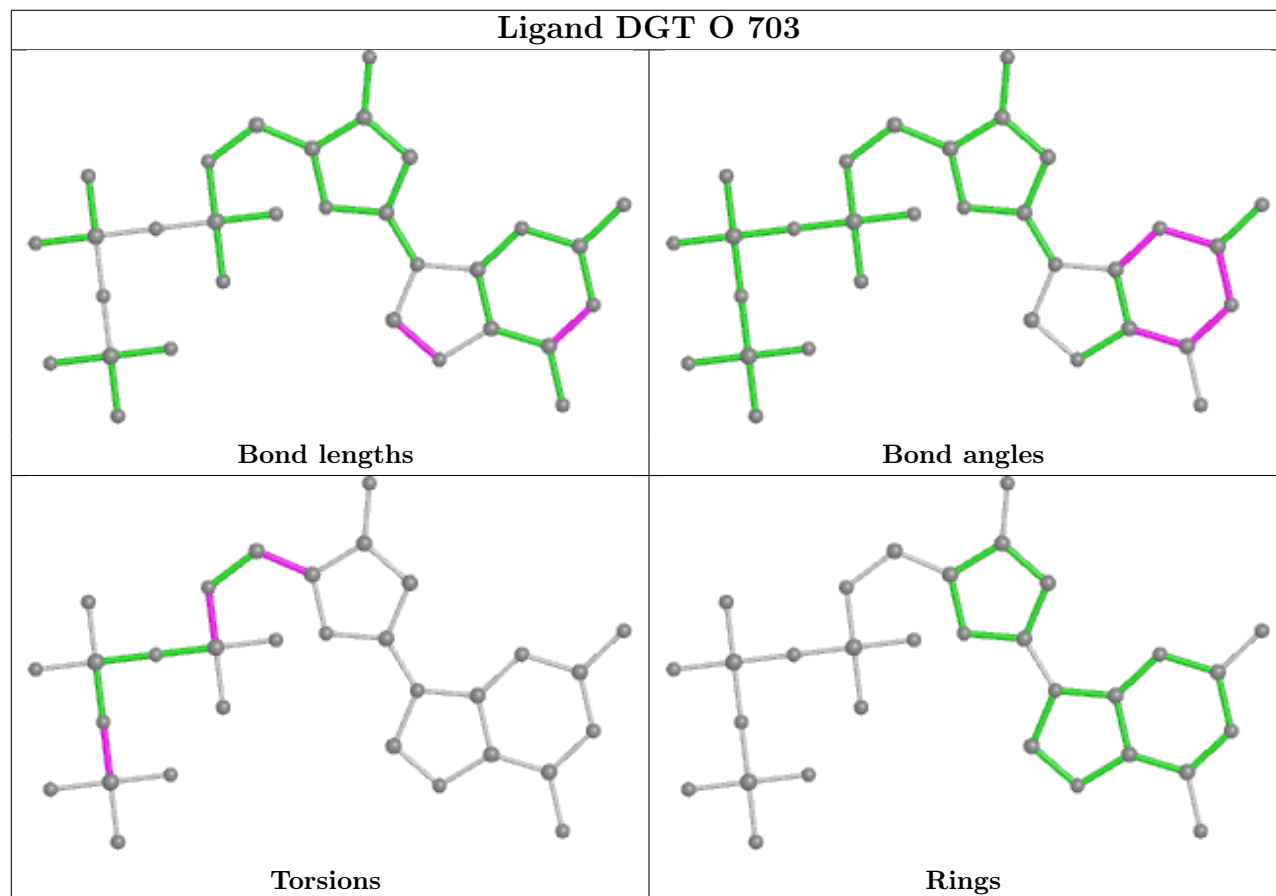
## Ligand DGT J 703

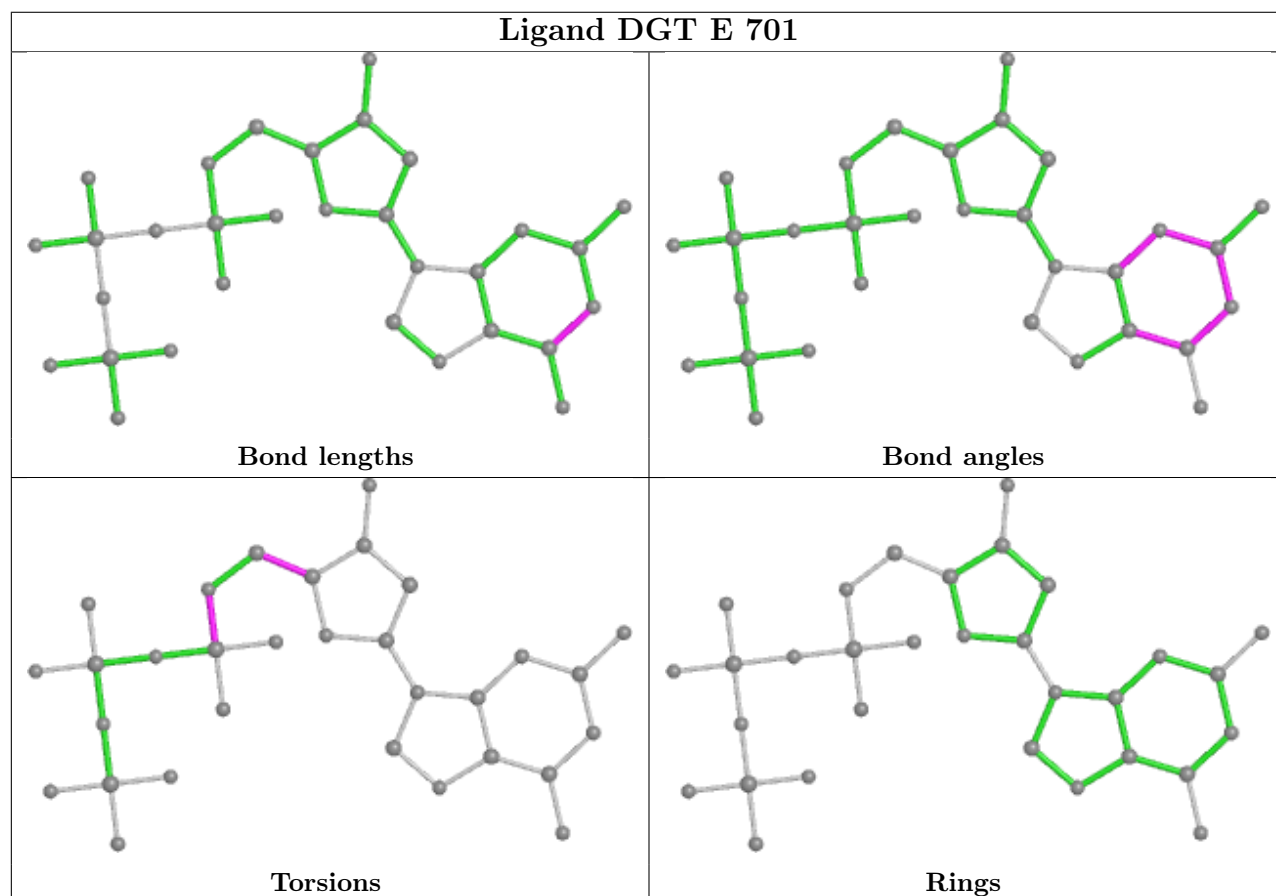
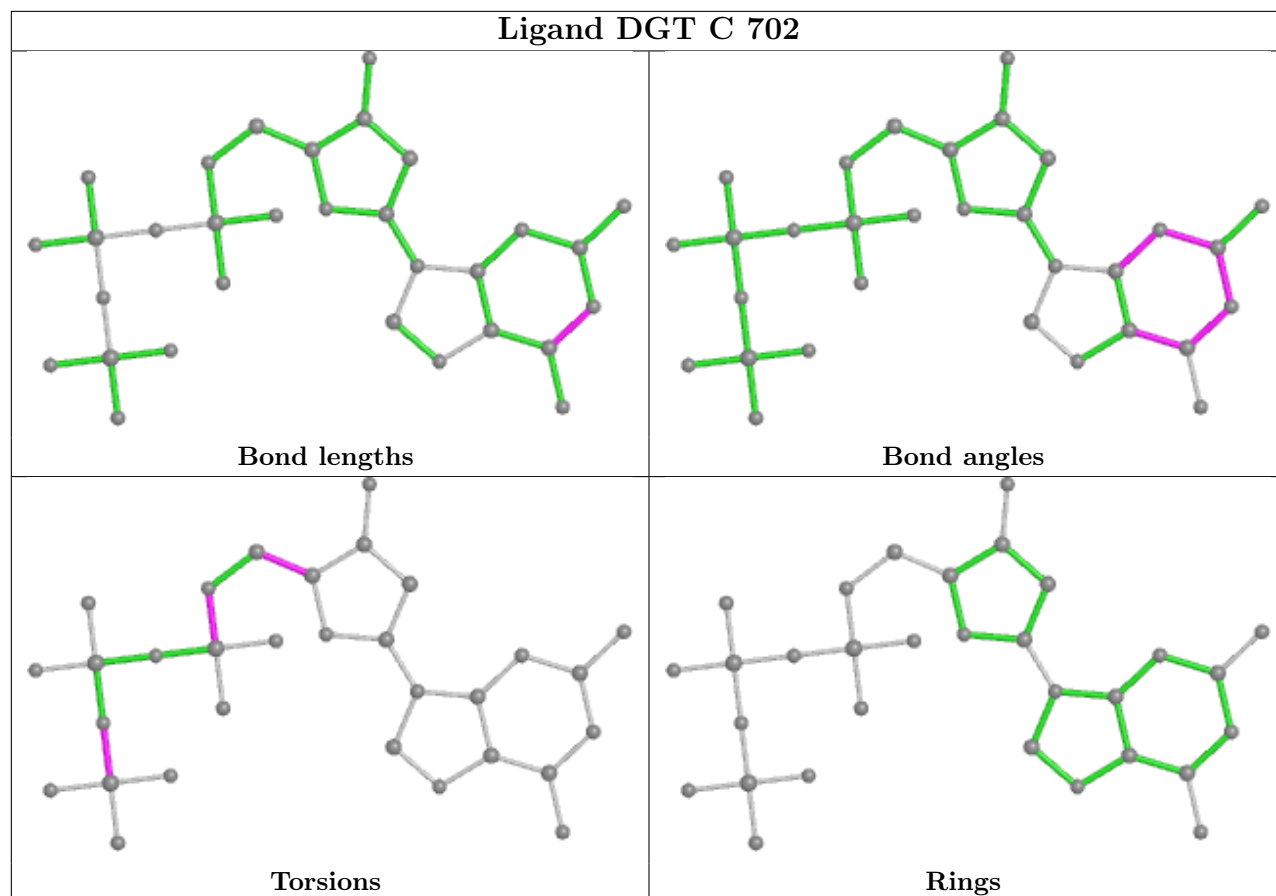


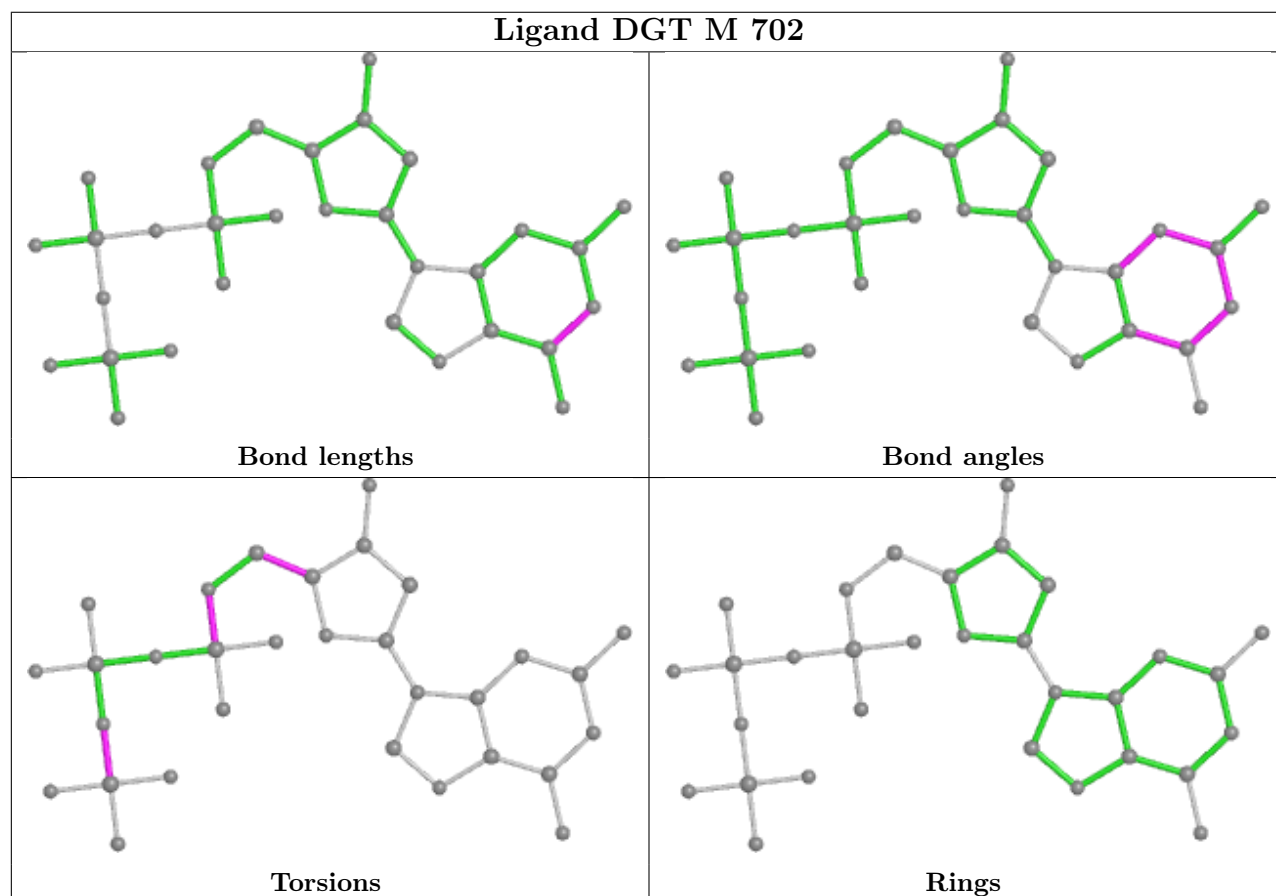
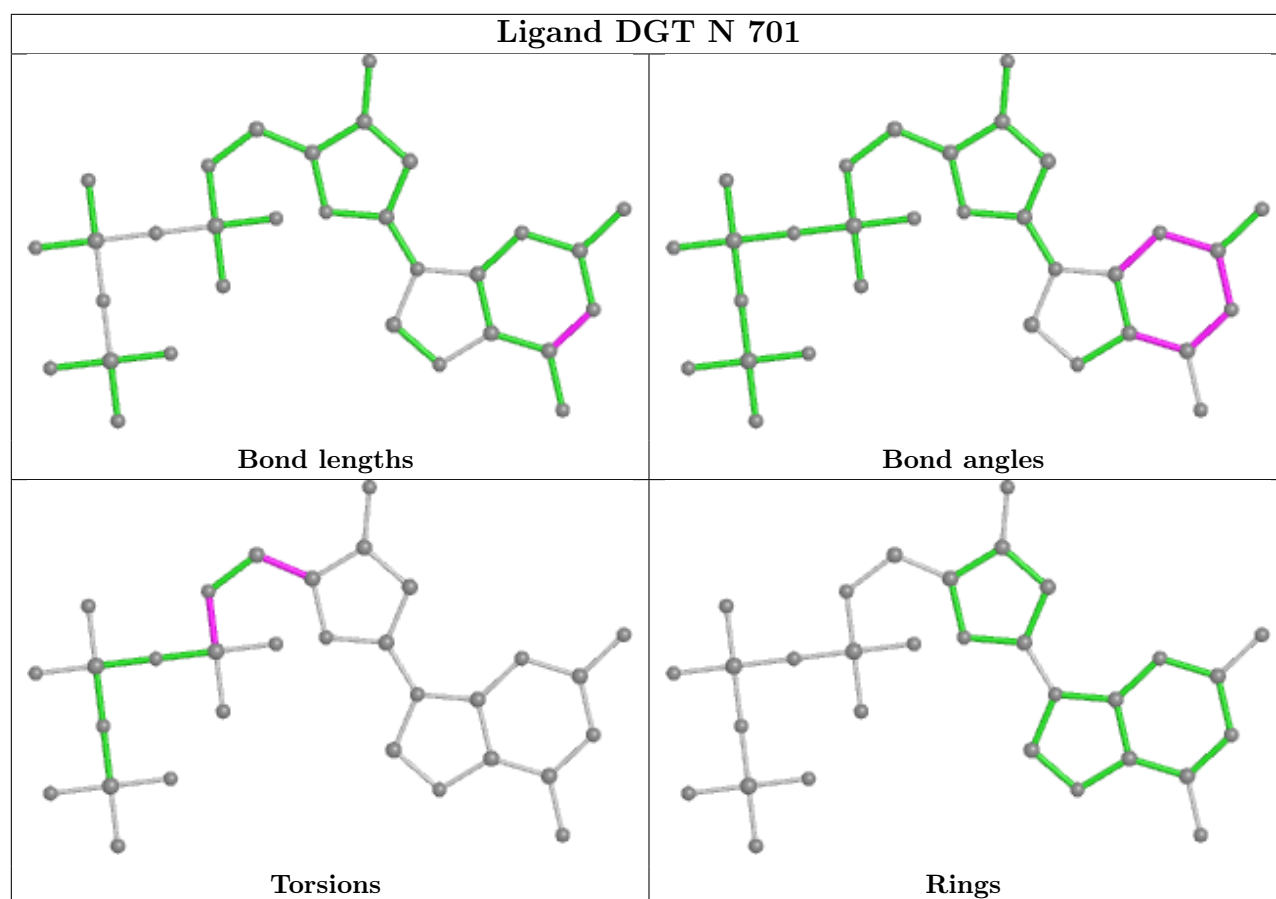
## Ligand DGT E 703

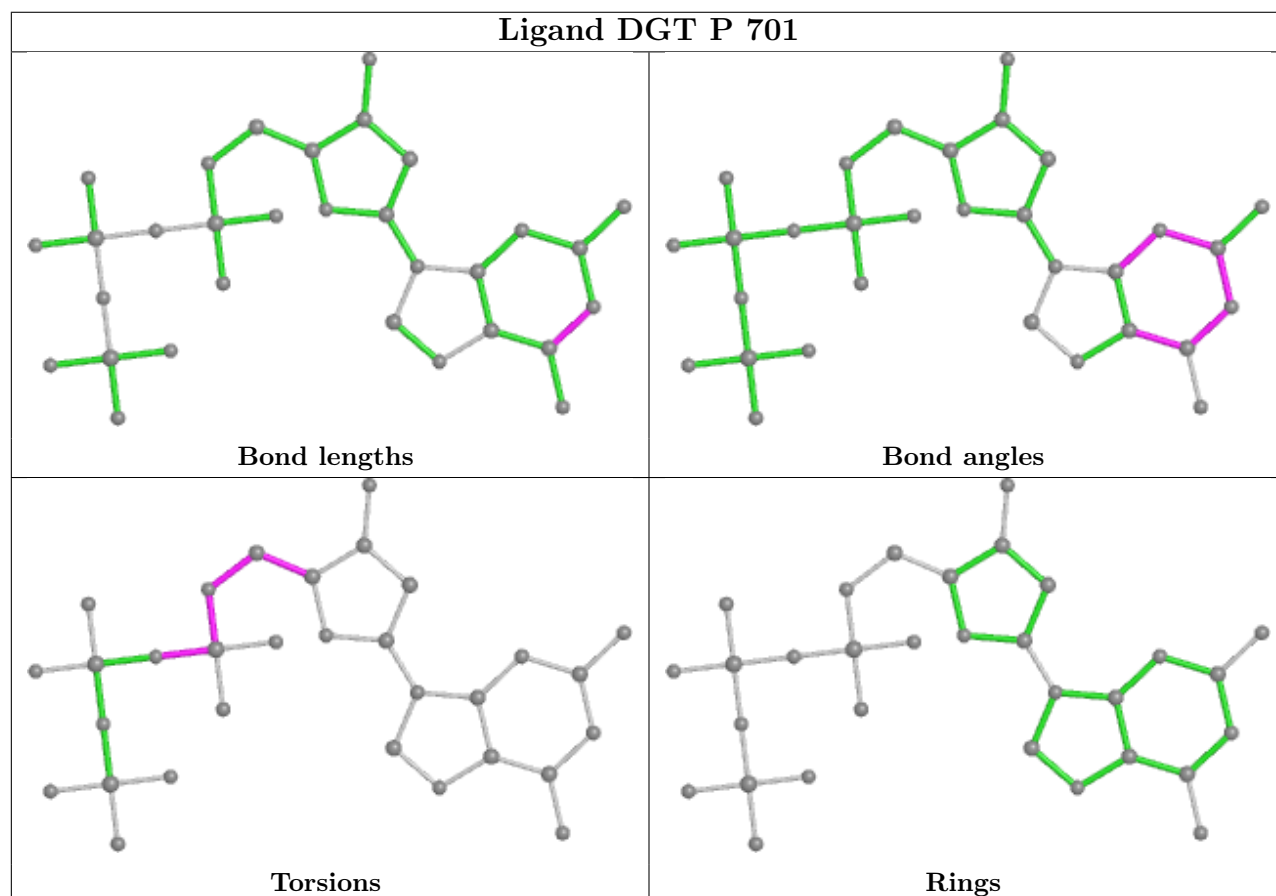
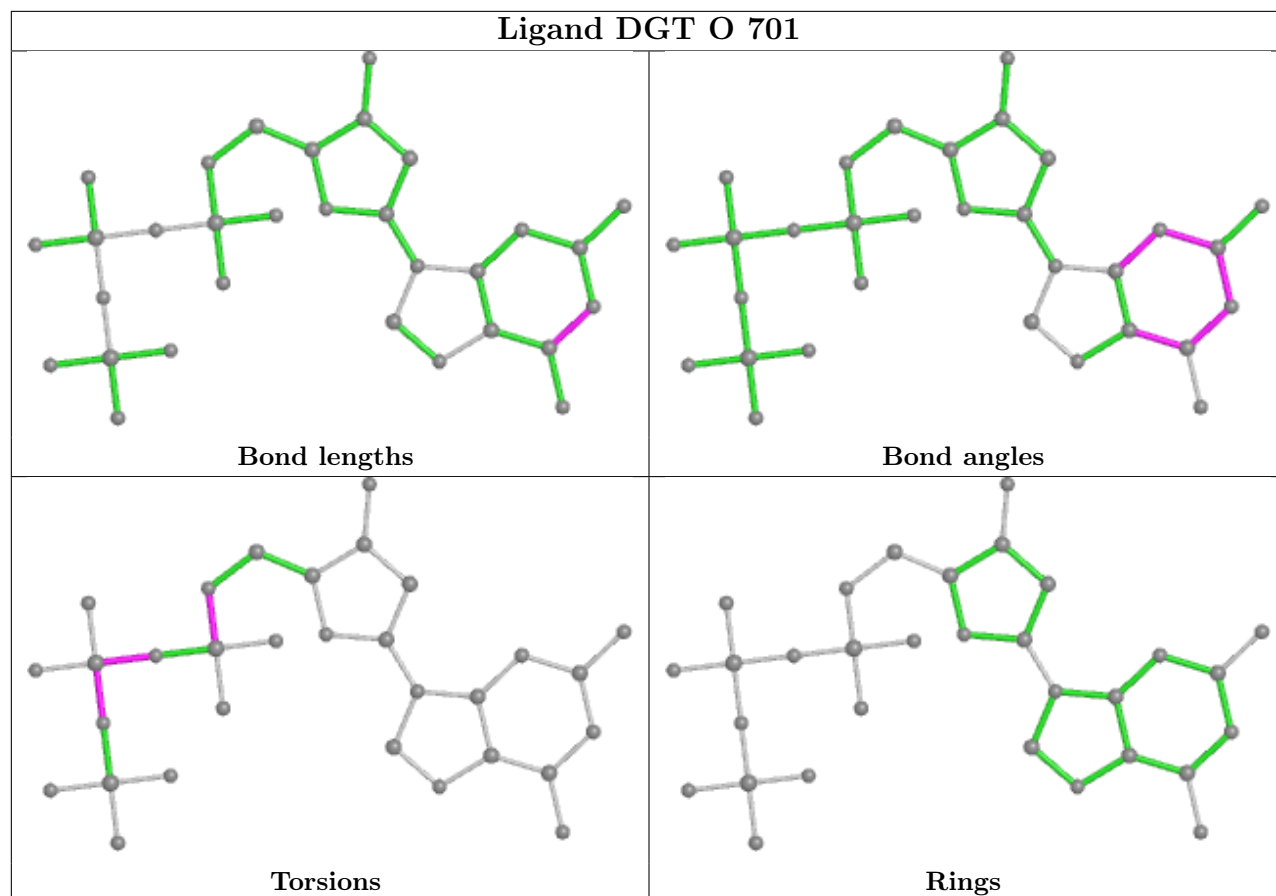


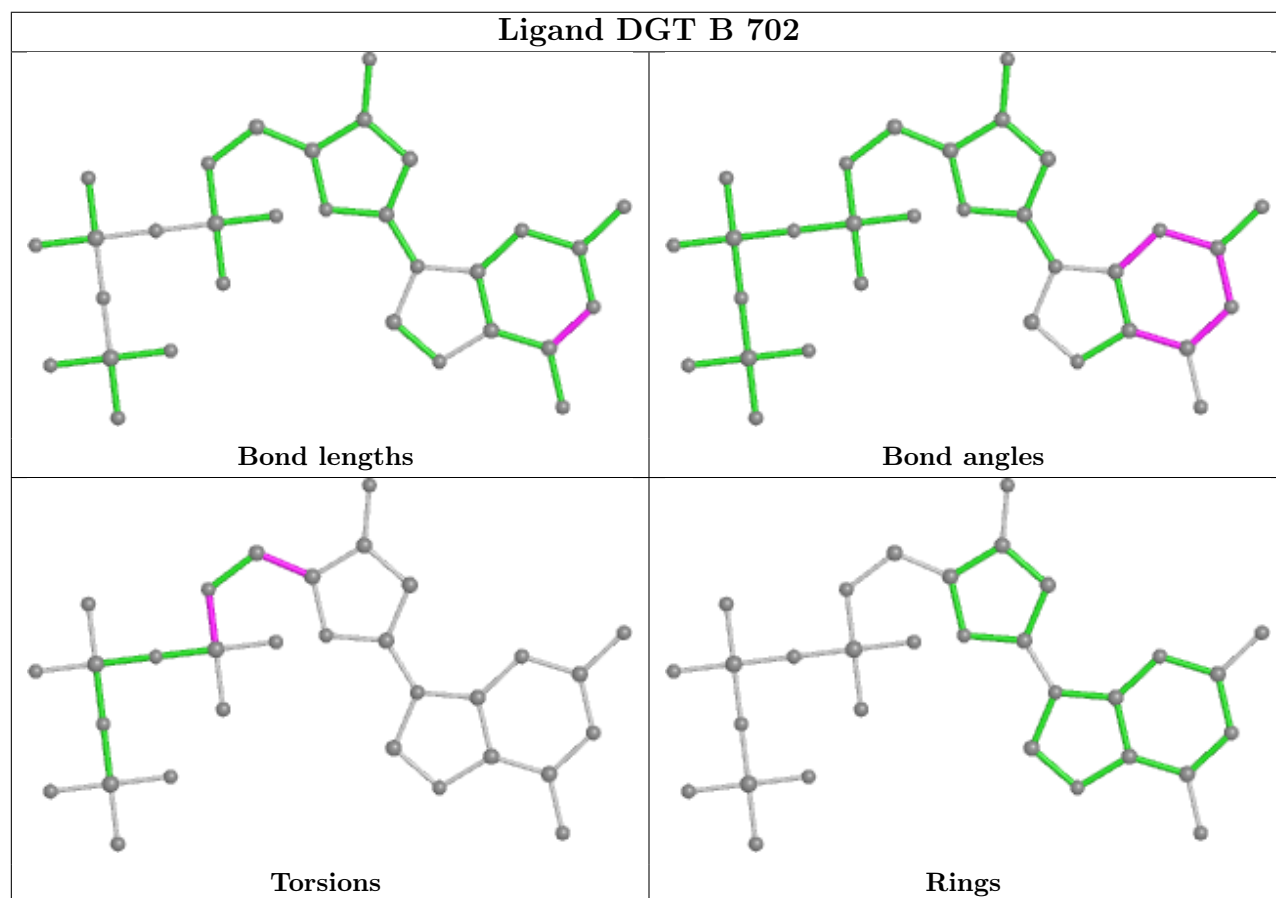
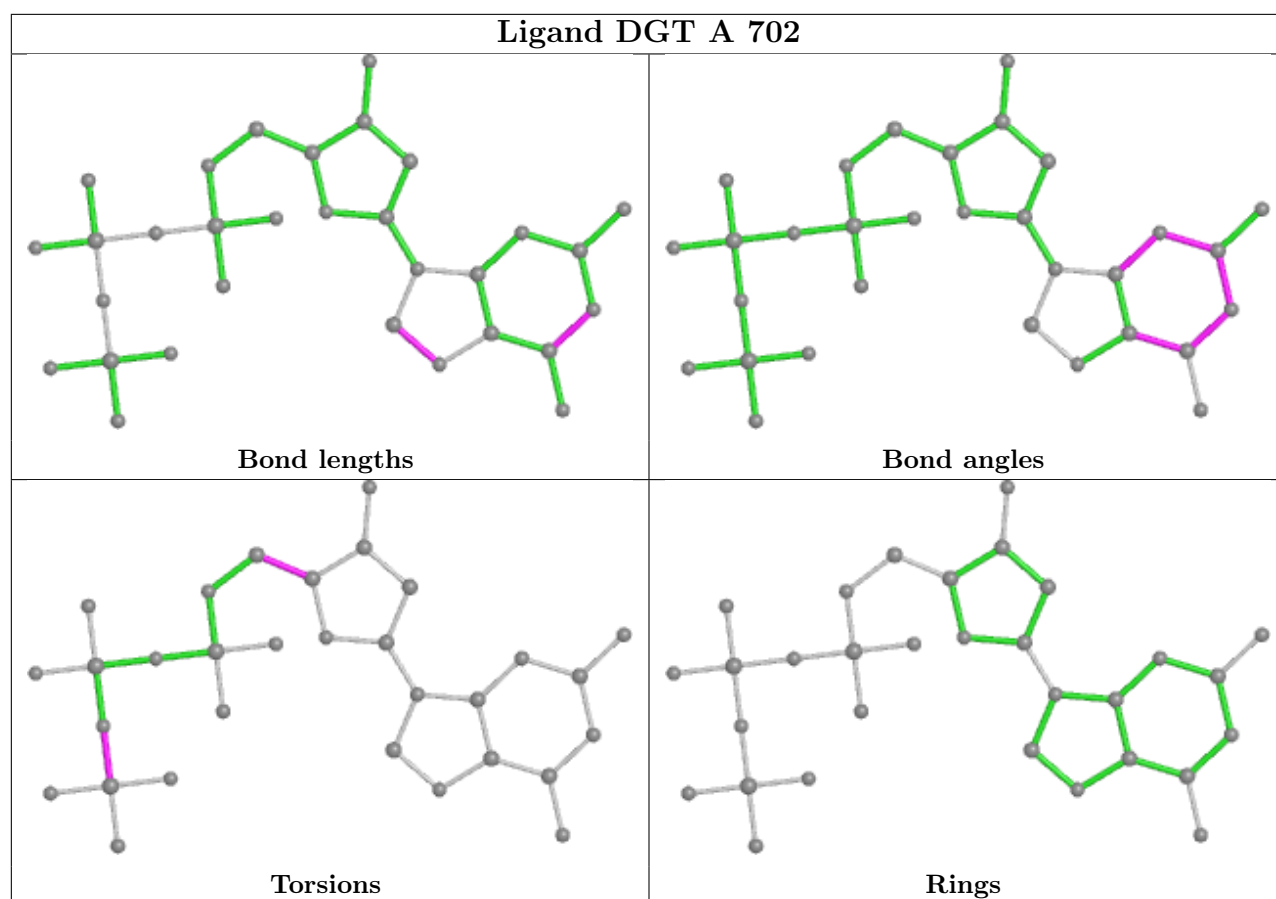
## Ligand DGT O 703



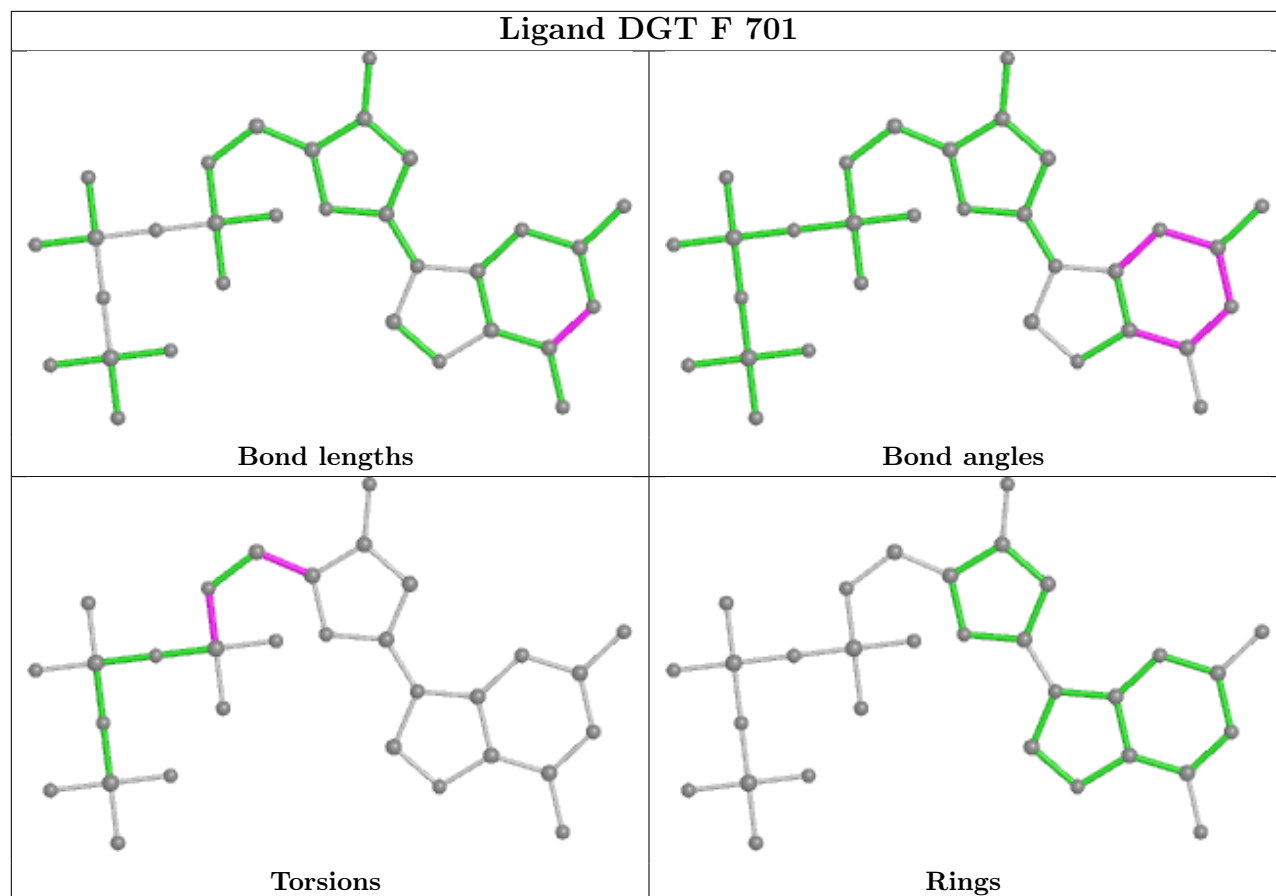




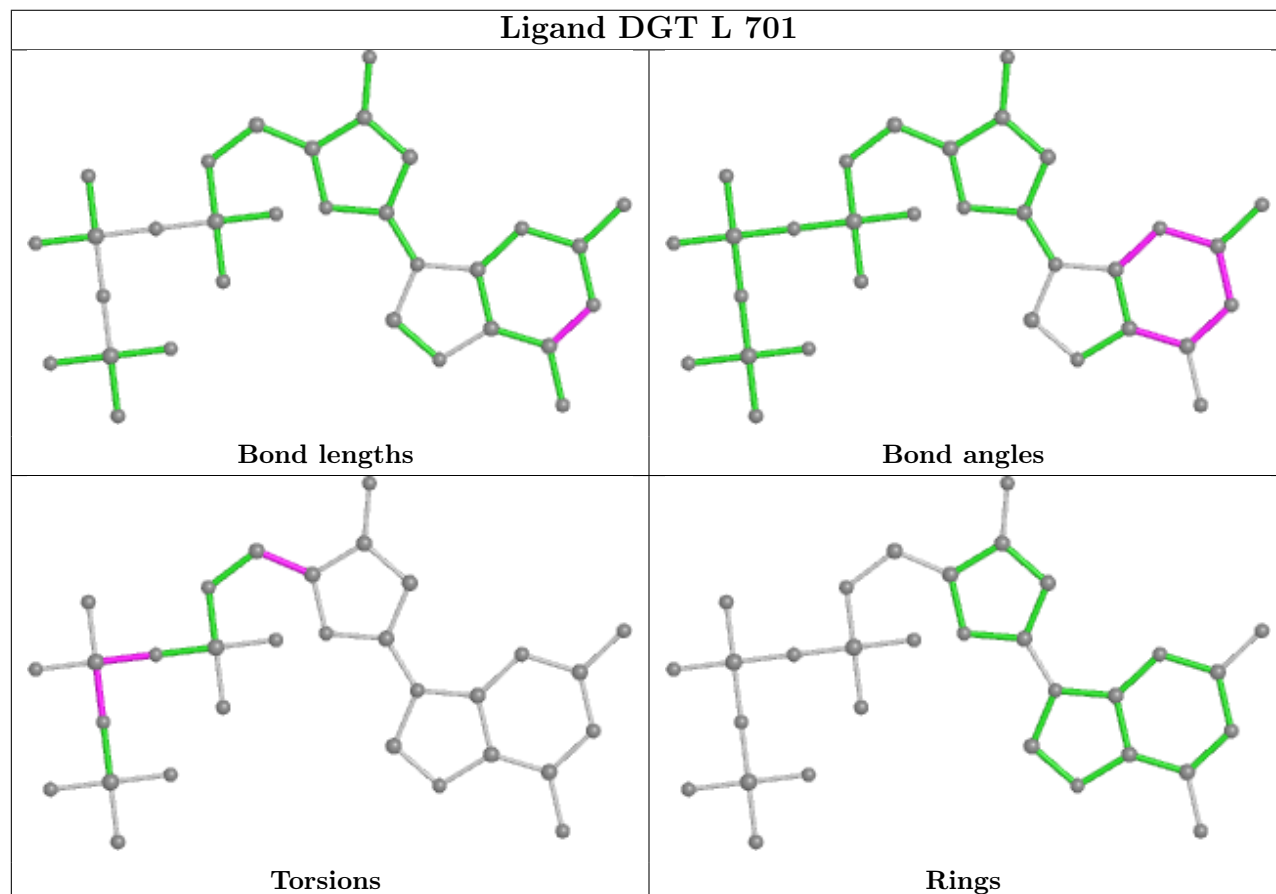




## Ligand DGT F 701



## Ligand DGT L 701



## 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](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

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

Unable to reproduce the depositors R factor - this section is therefore empty.

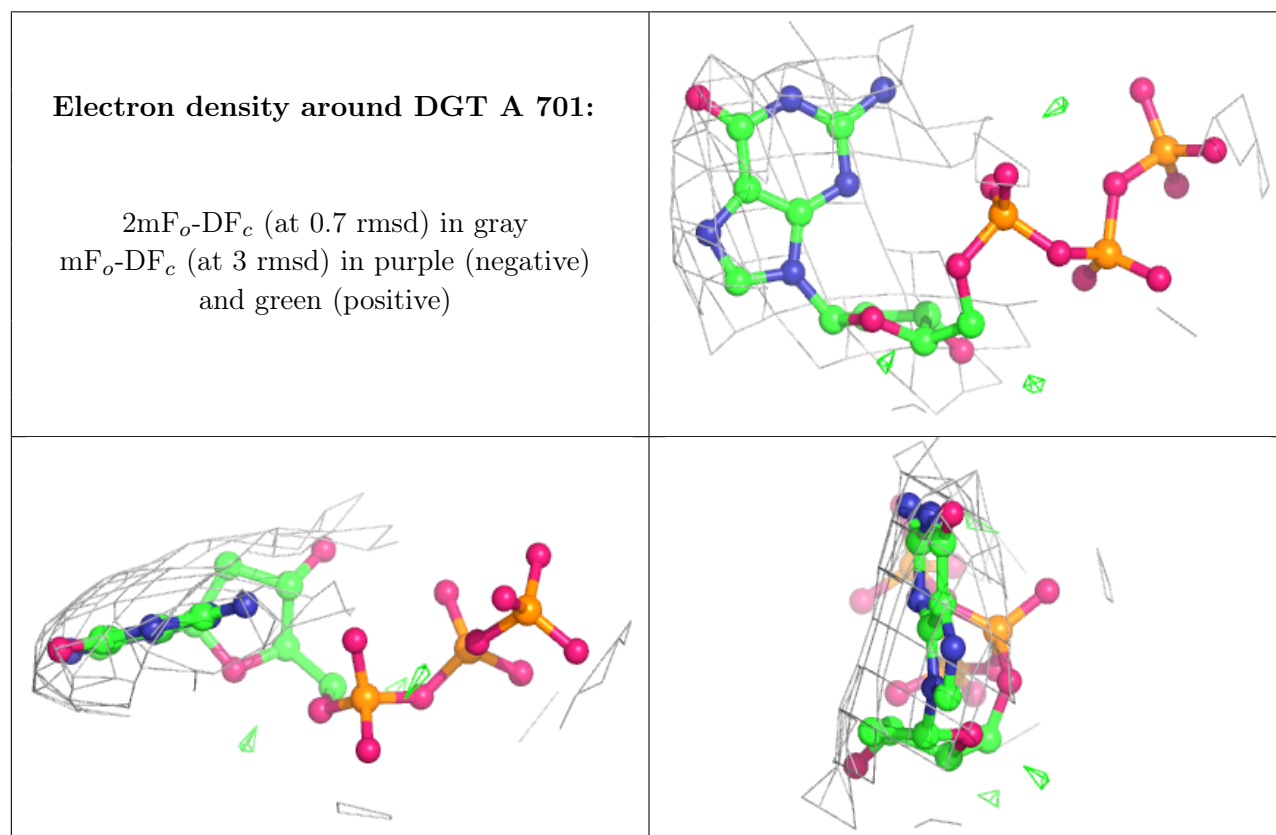
### 6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.4 Ligands [i](#)

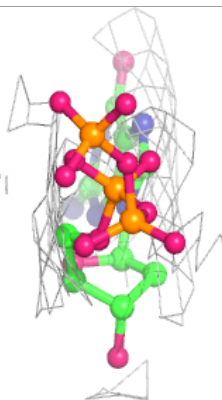
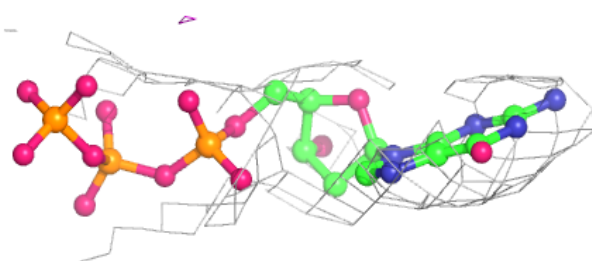
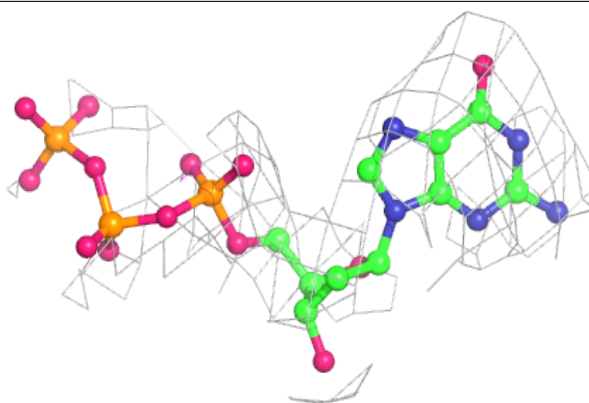
Unable to reproduce the depositors R factor - this section is therefore empty.

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.

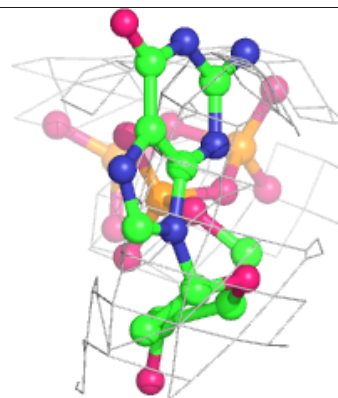
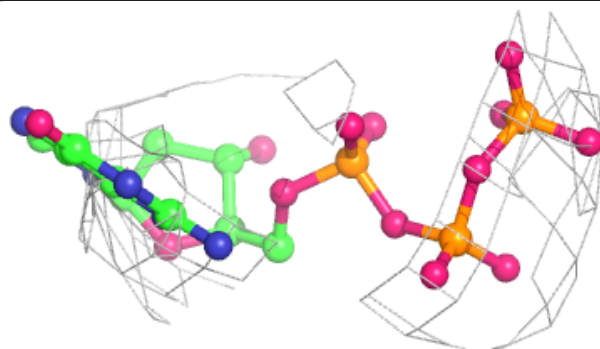
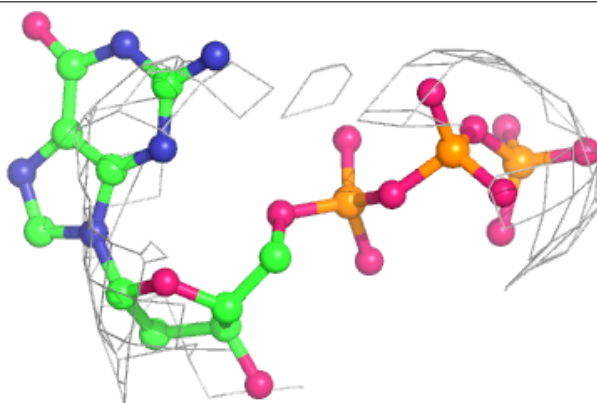


**Electron density around DGT A 702:**

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

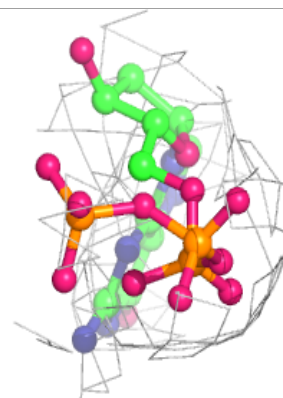
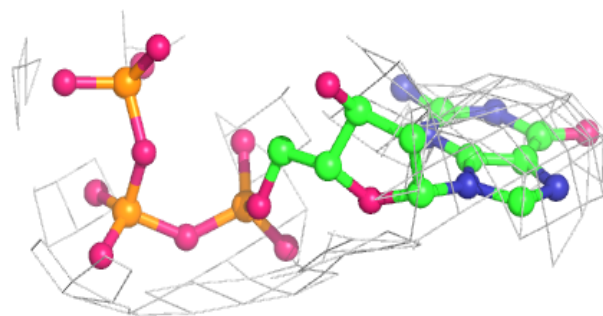
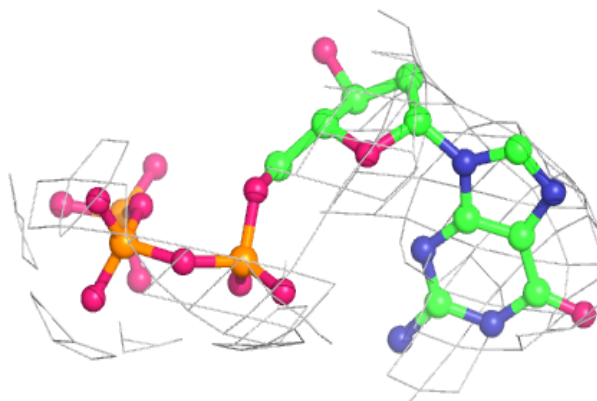
**Electron density around DGT A 703:**

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

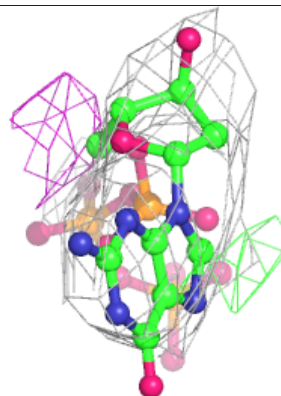
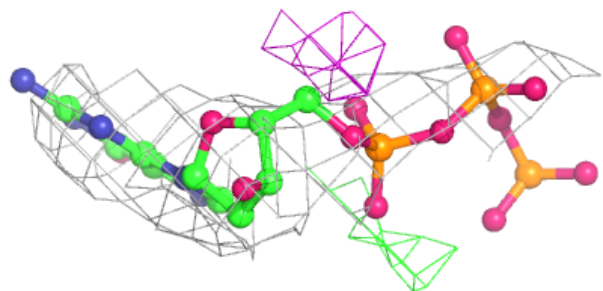
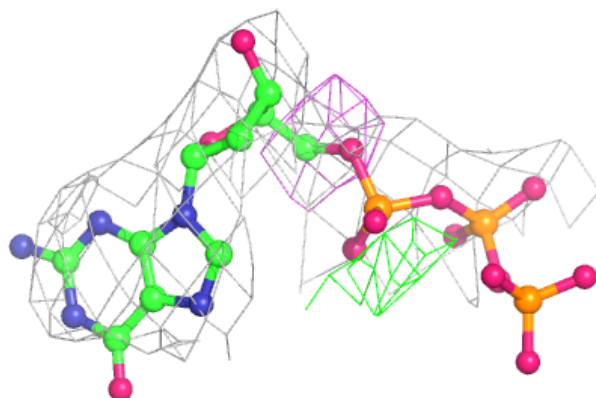


**Electron density around DGT B 701:**

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

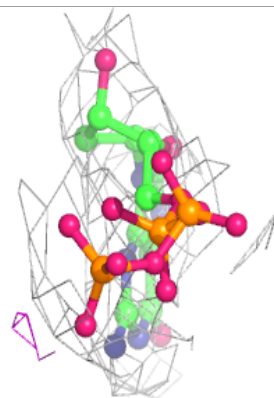
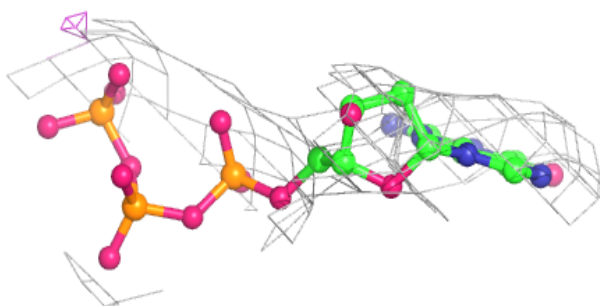
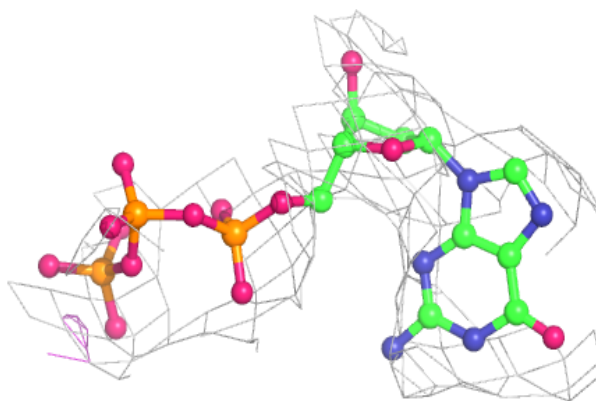
**Electron density around DGT B 702:**

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

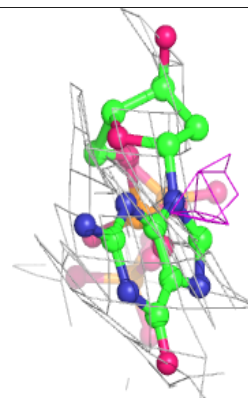
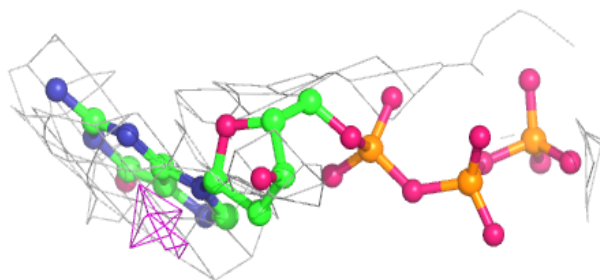
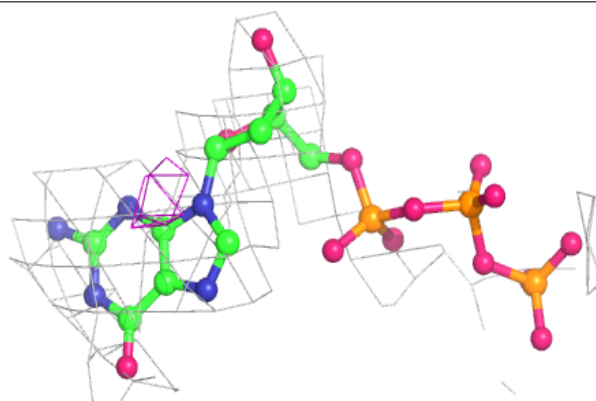


**Electron density around DGT C 701:**

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

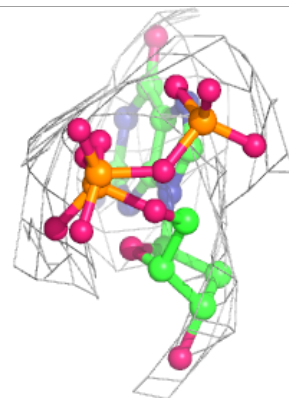
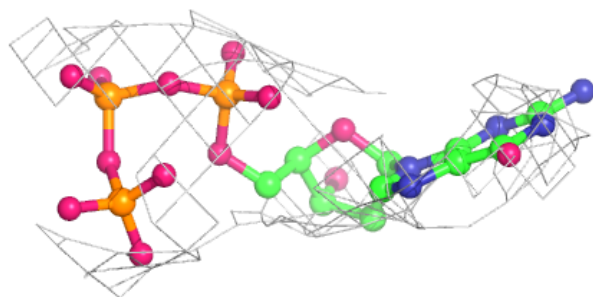
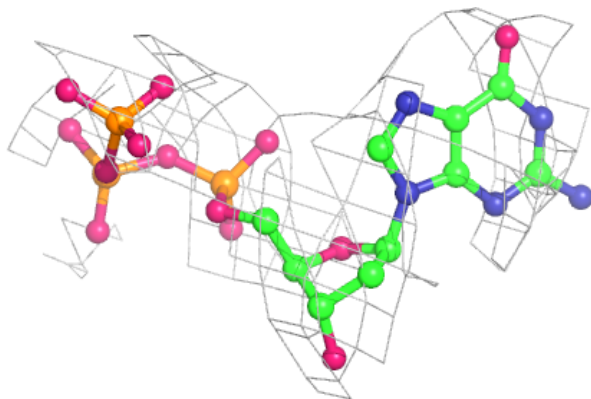
**Electron density around DGT C 702:**

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

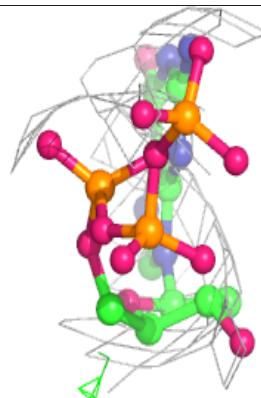
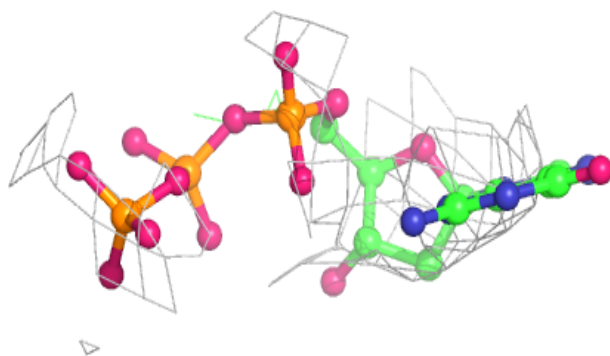
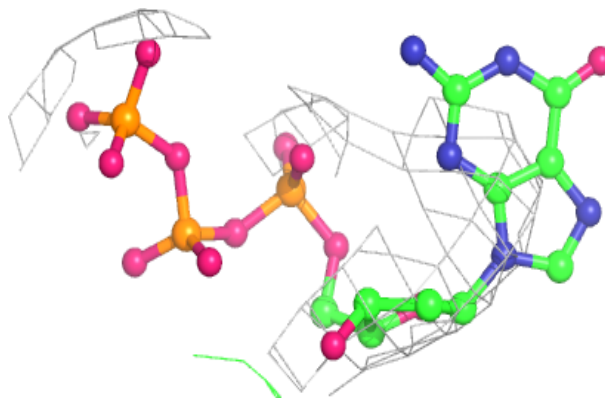


**Electron density around DGT D 701:**

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

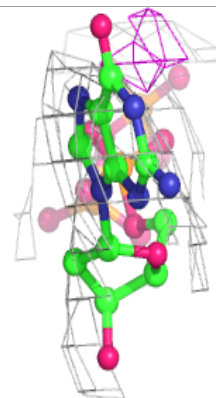
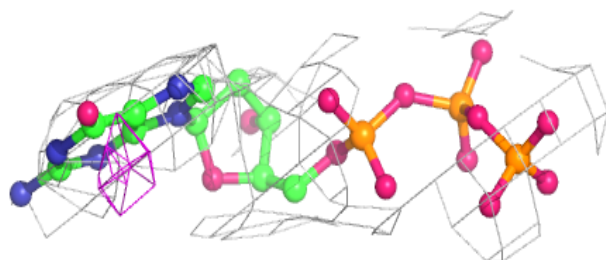
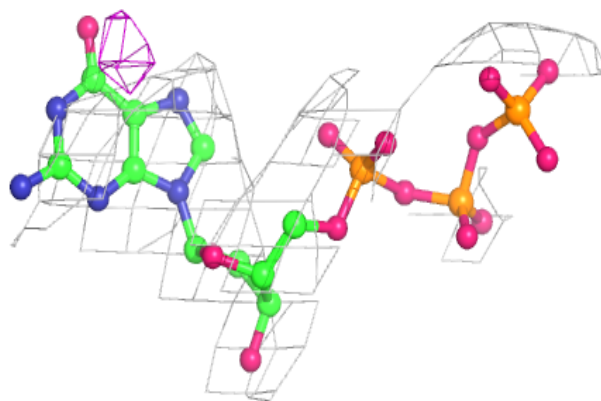
**Electron density around DGT E 701:**

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



**Electron density around DGT E 702:**

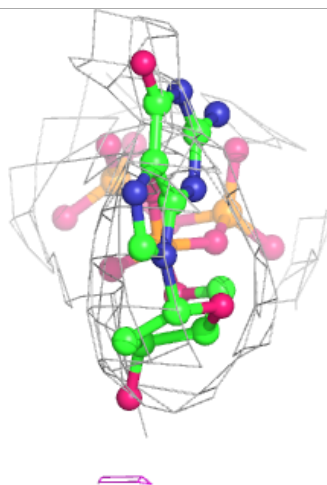
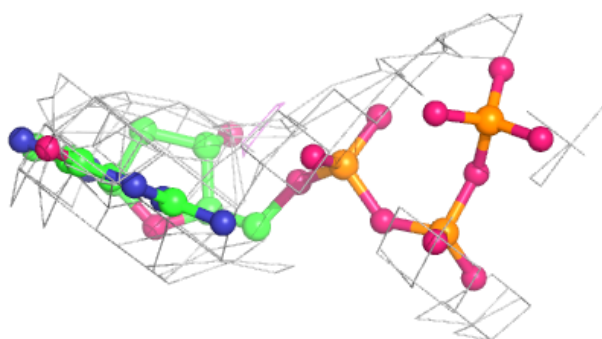
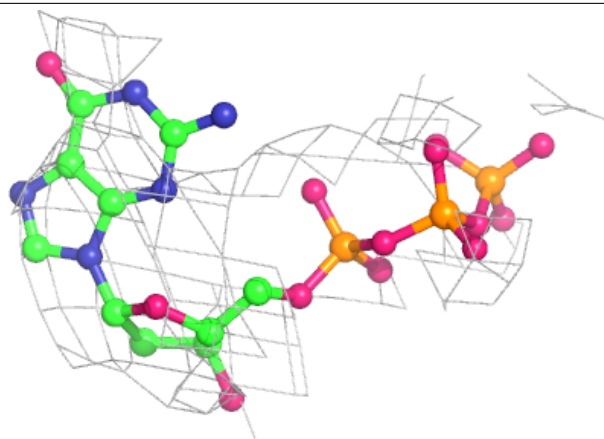
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





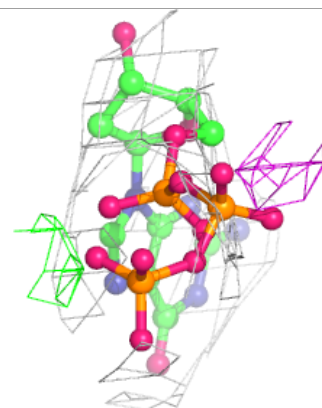
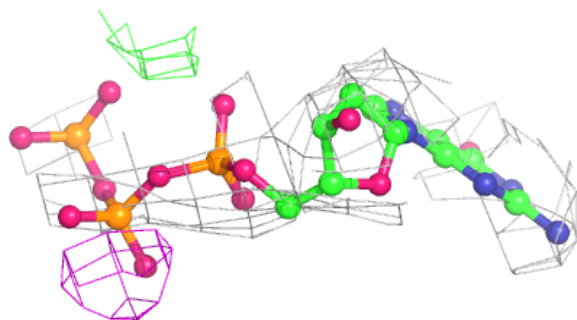
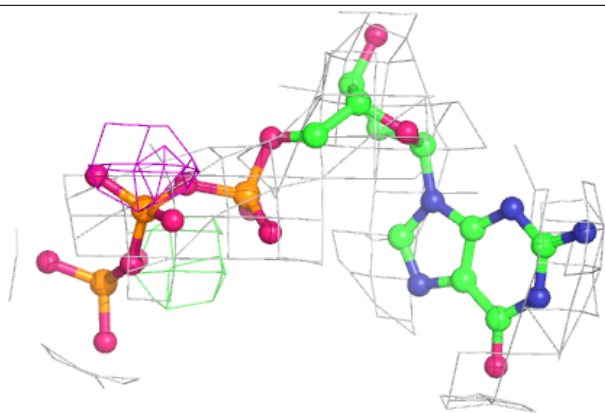
**Electron density around DGT E 703:**

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

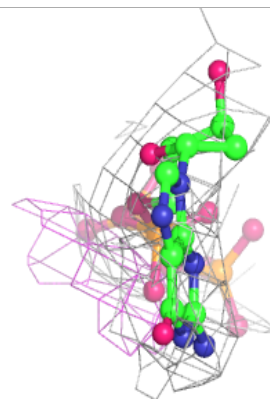
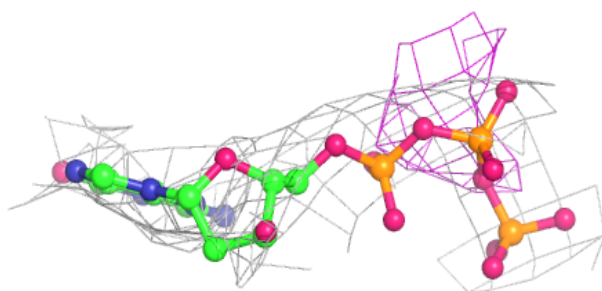
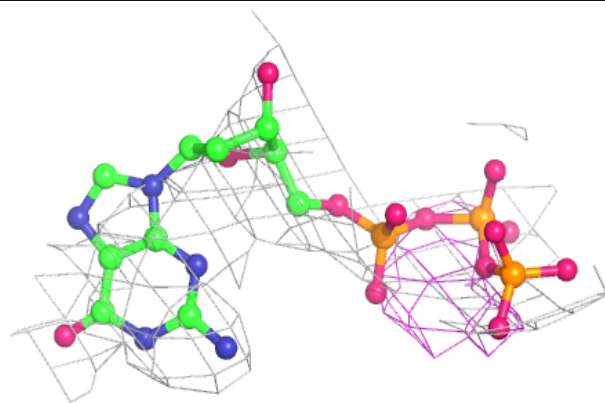


**Electron density around DGT F 701:**

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

**Electron density around DGT F 702:**

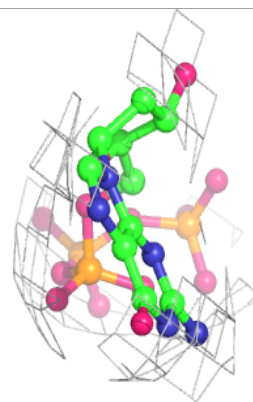
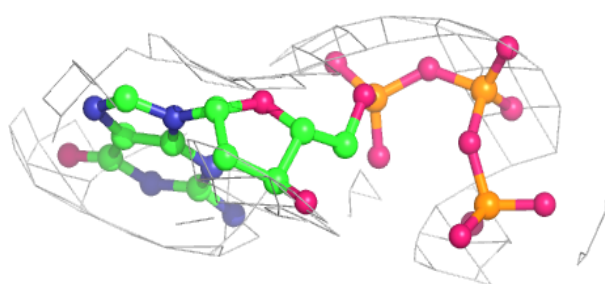
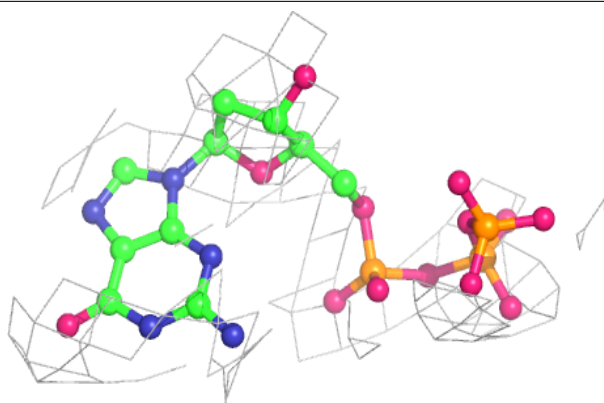
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



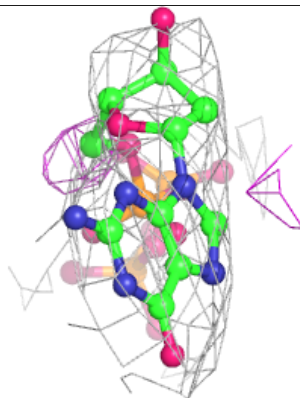
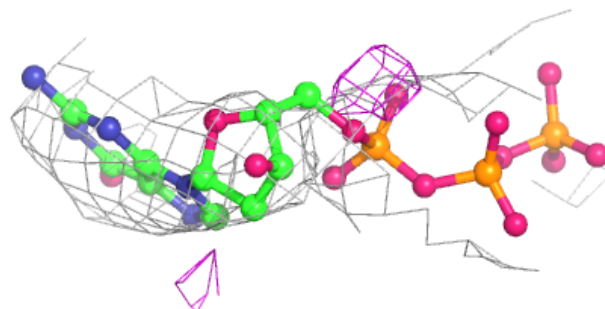
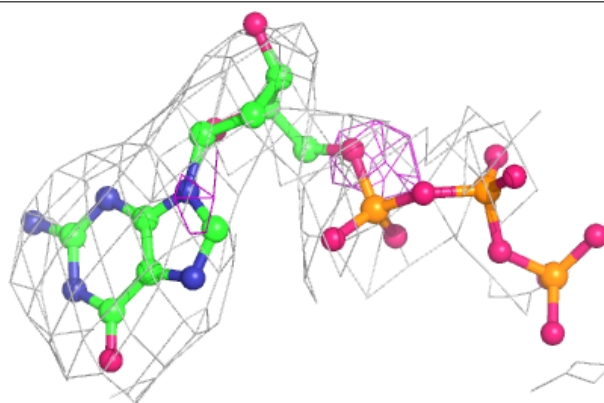


**Electron density around DGT G 701:**

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

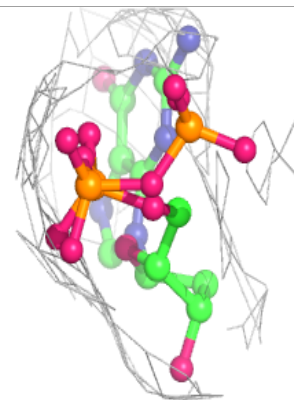
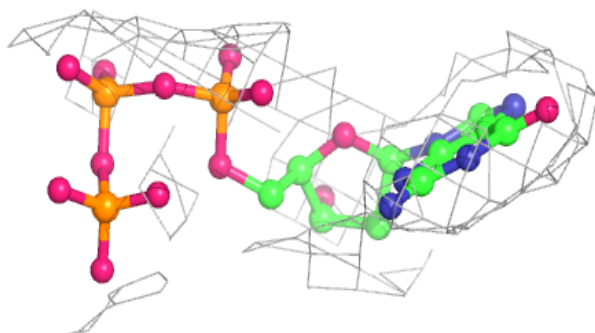
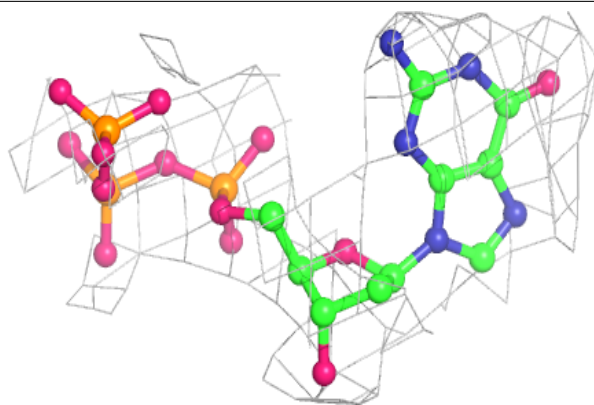
**Electron density around DGT G 702:**

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

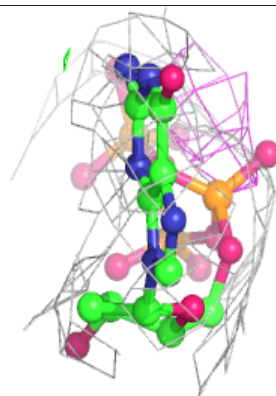
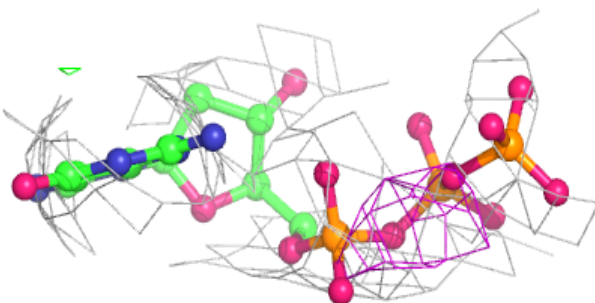
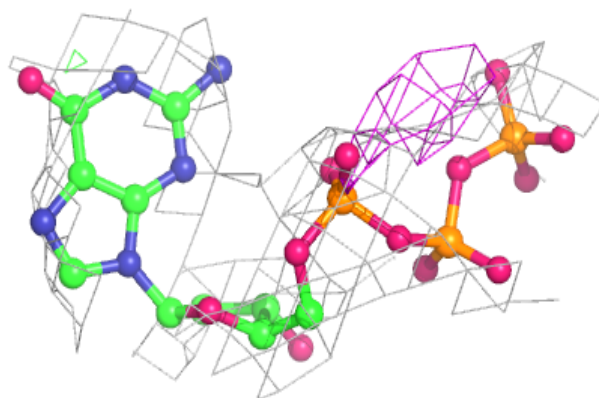


**Electron density around DGT H 701:**

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

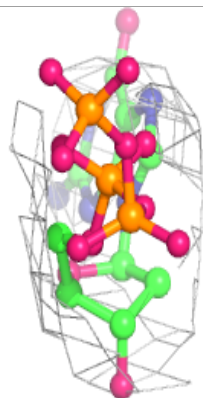
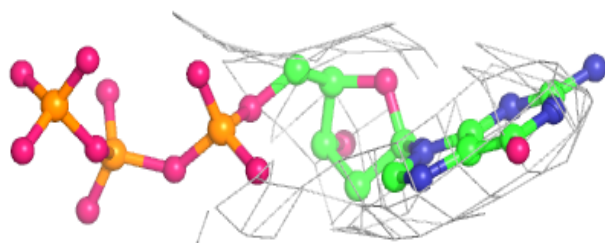
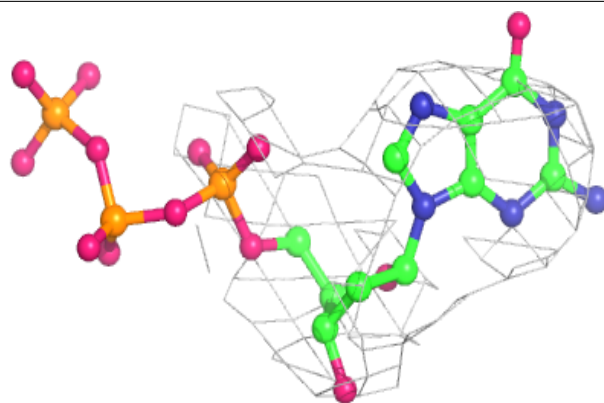
**Electron density around DGT I 701:**

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

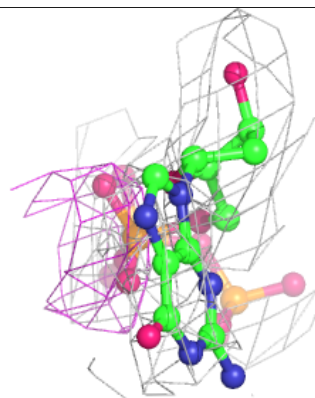
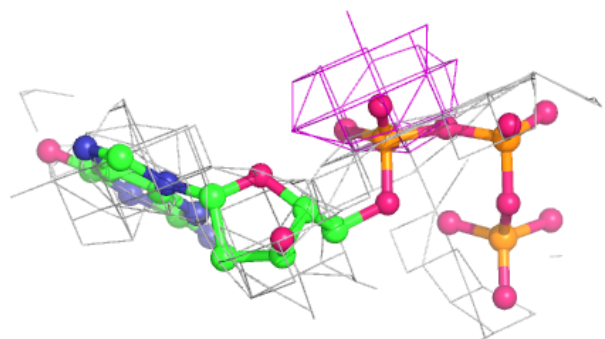
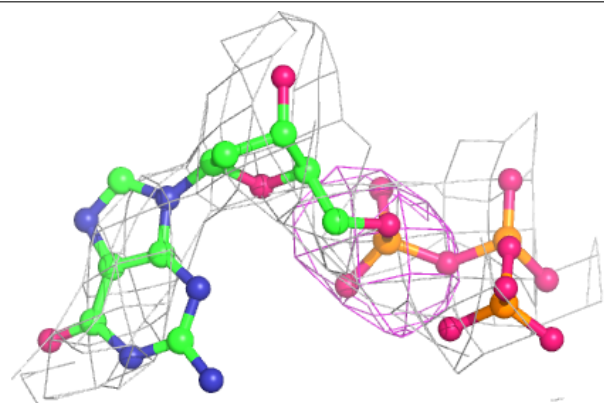


**Electron density around DGT I 702:**

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

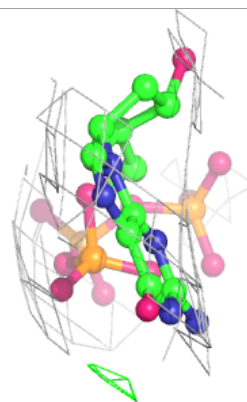
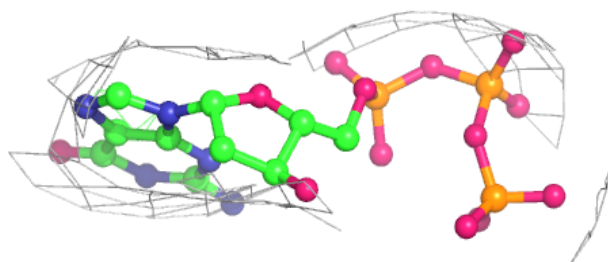
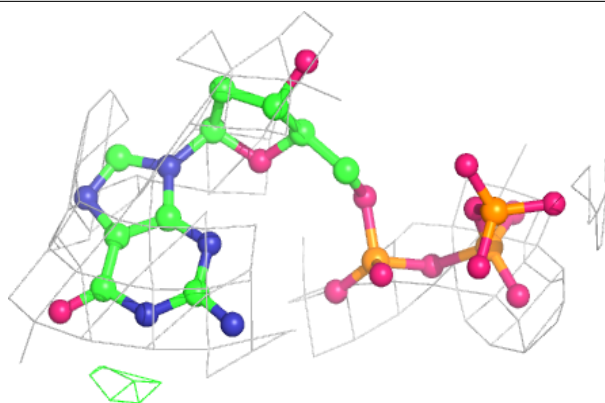
**Electron density around DGT J 701:**

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

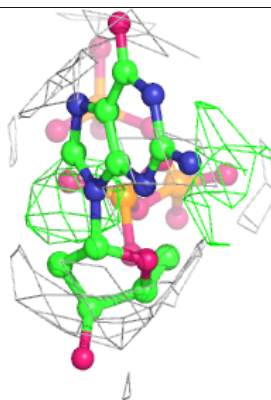
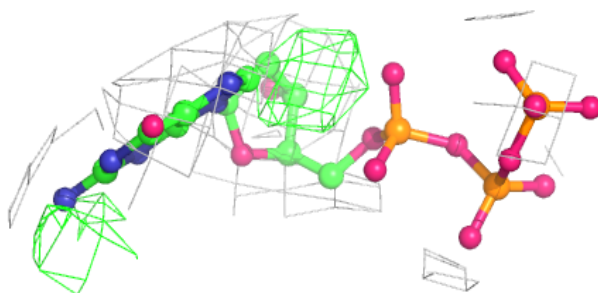
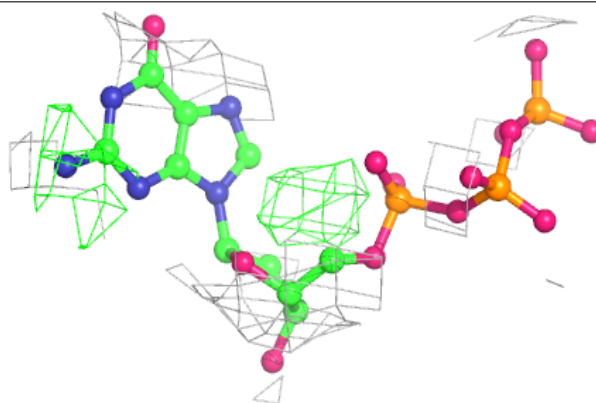


**Electron density around DGT J 702:**

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

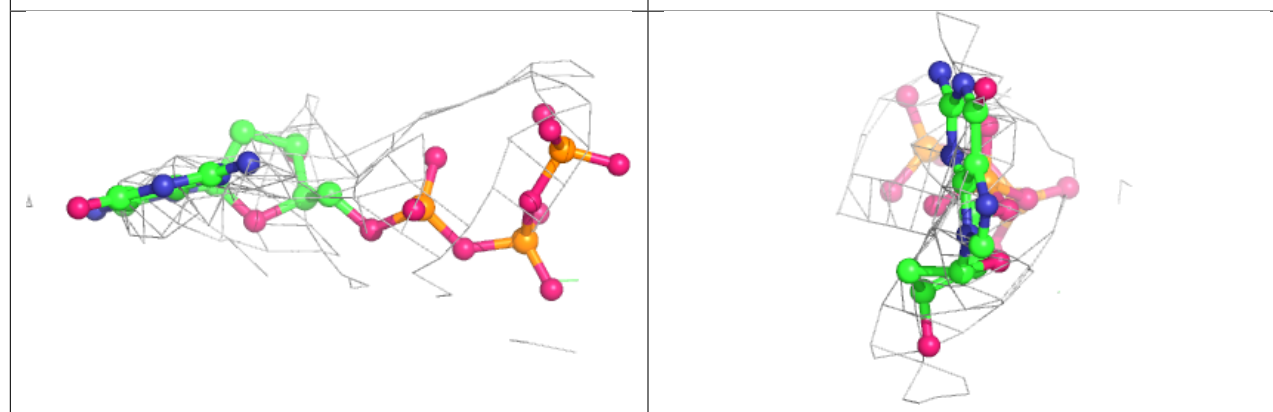
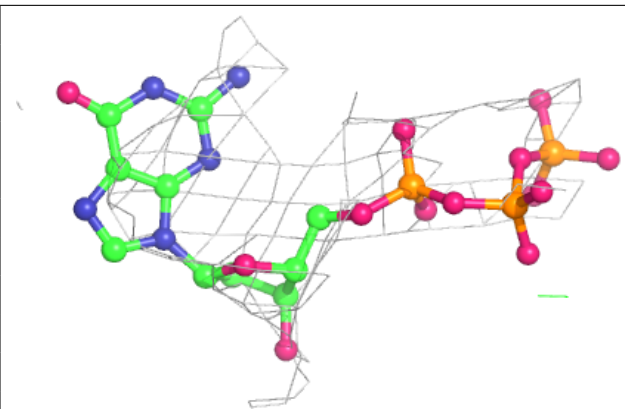
**Electron density around DGT J 703:**

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

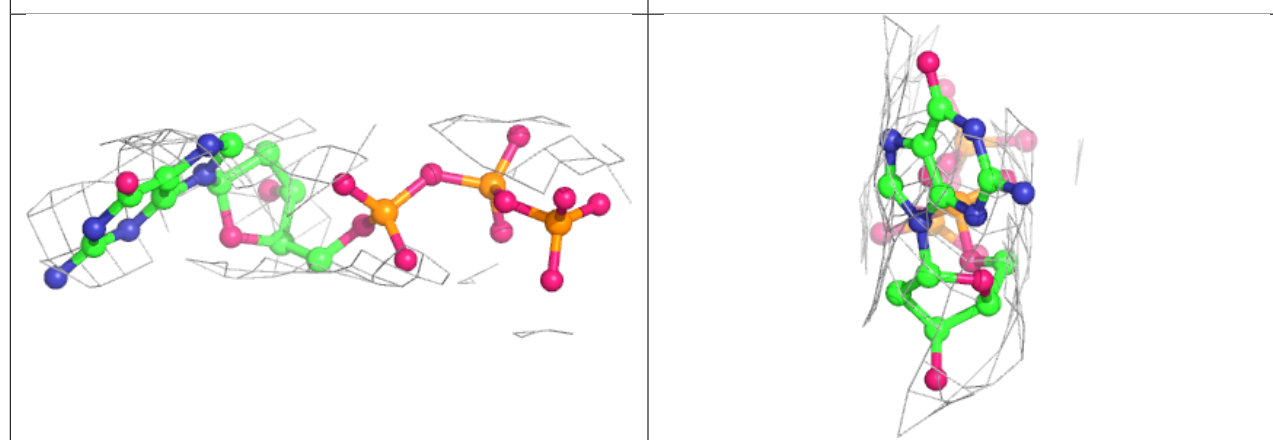
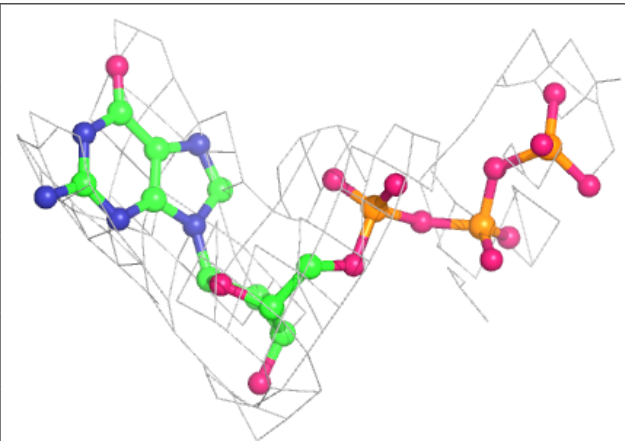


**Electron density around DGT K 701:**

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

**Electron density around DGT K 702:**

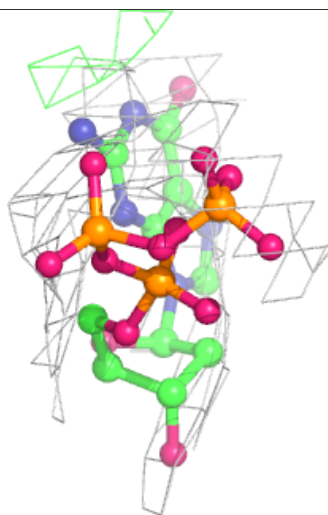
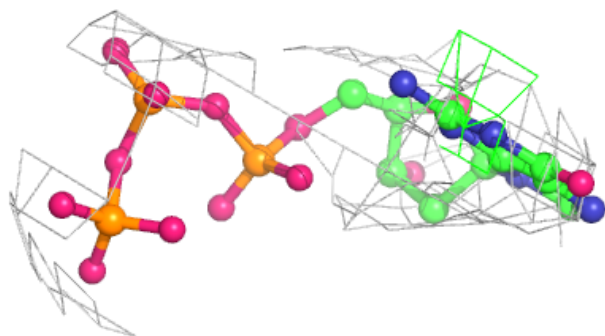
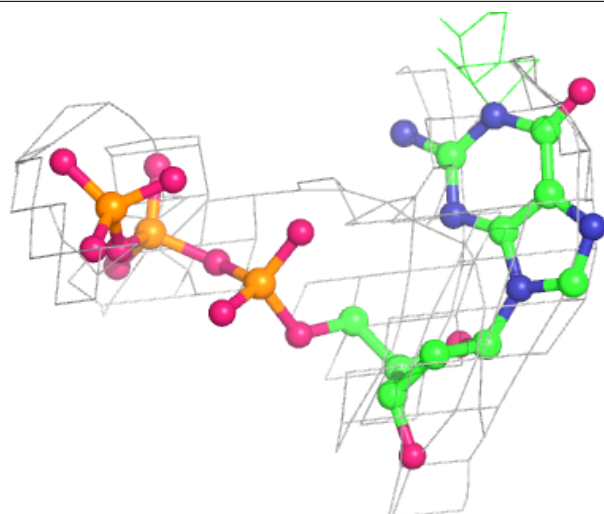
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





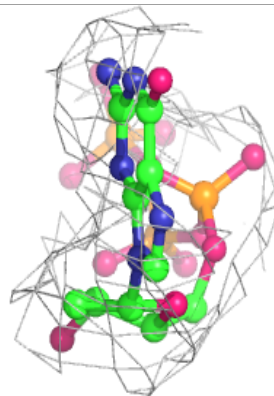
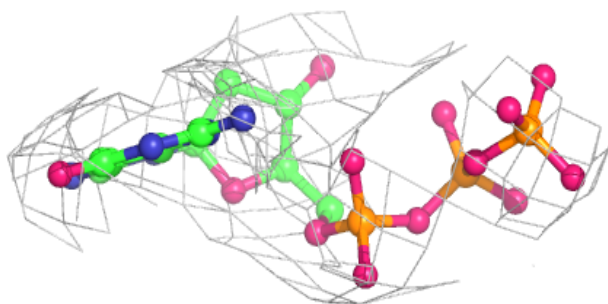
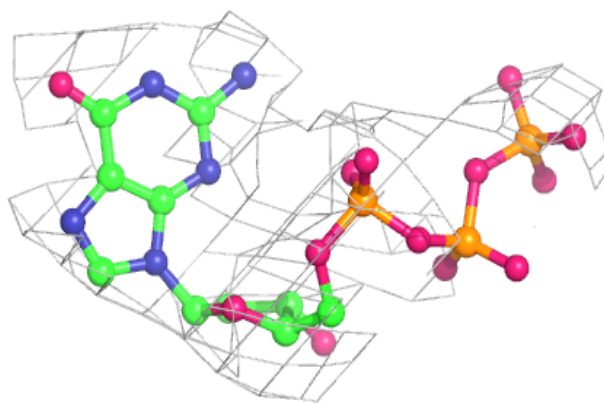
**Electron density around DGT L 701:**

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

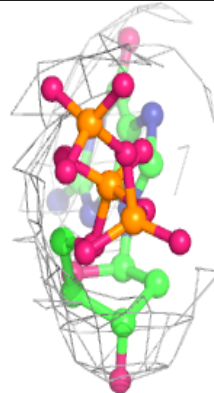
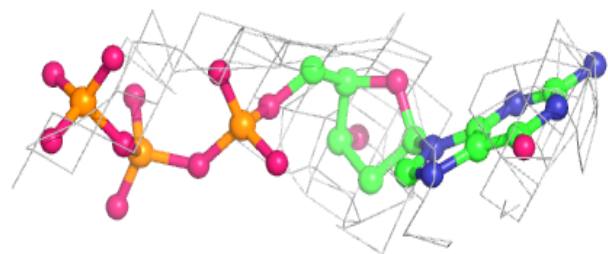
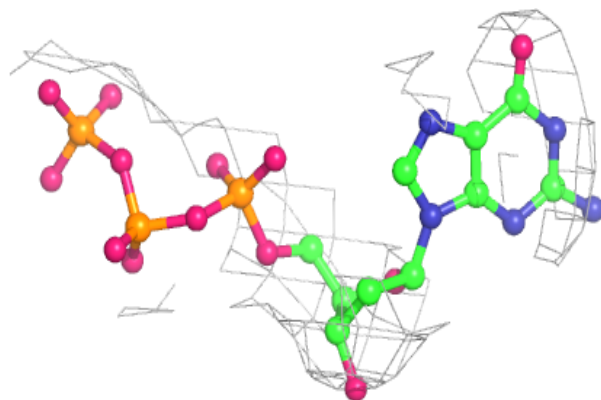


**Electron density around DGT M 701:**

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

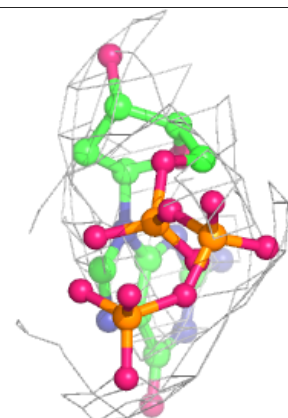
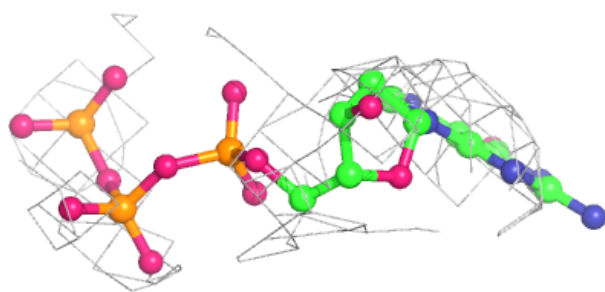
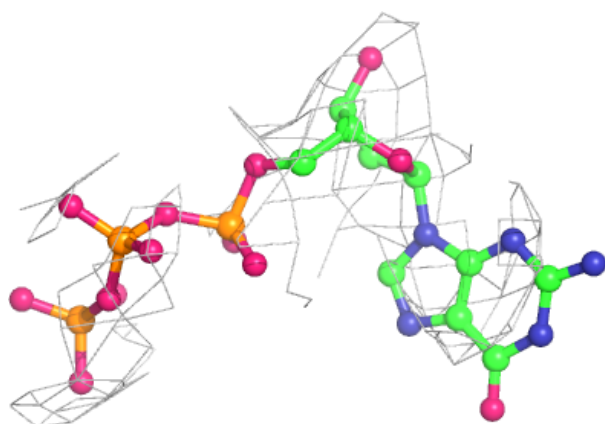
**Electron density around DGT M 702:**

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

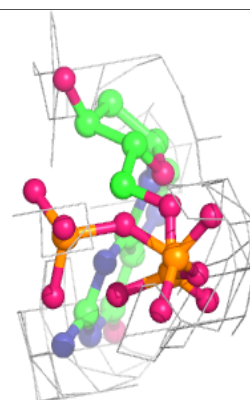
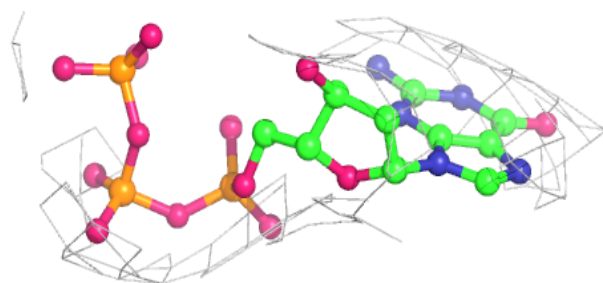
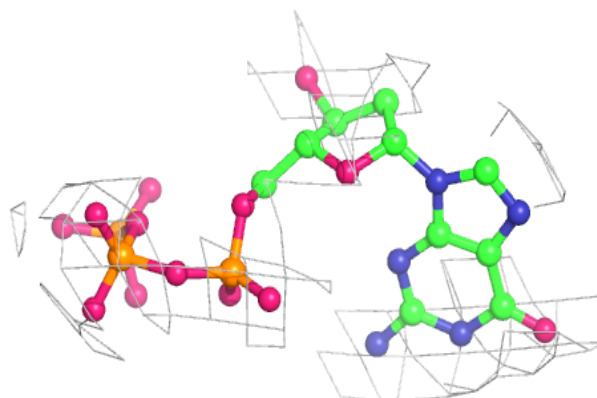


**Electron density around DGT N 701:**

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

**Electron density around DGT O 701:**

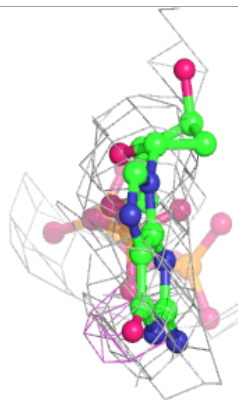
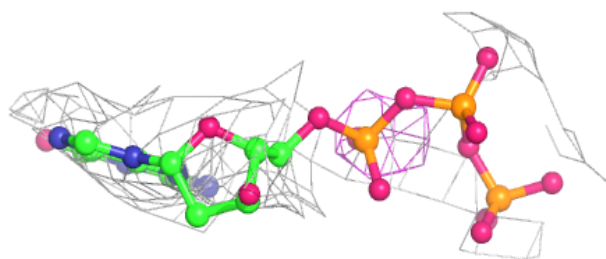
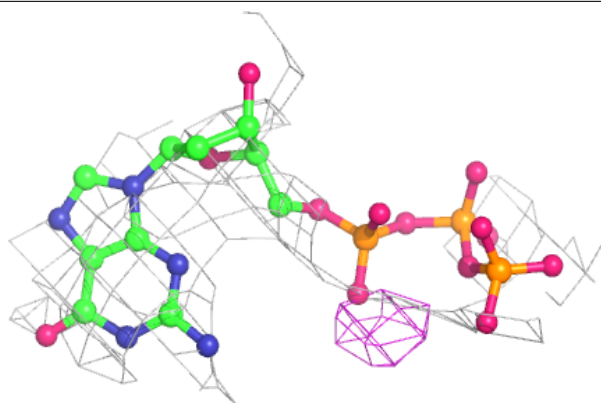
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



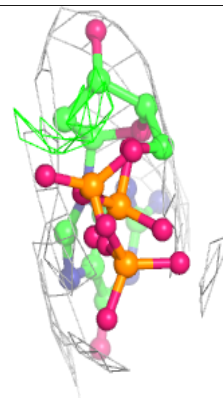
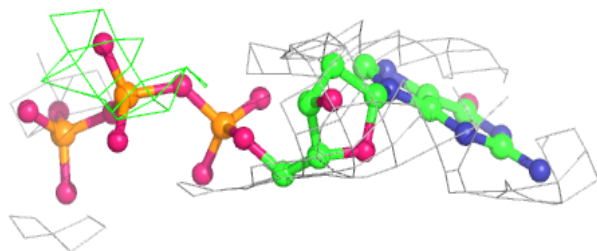
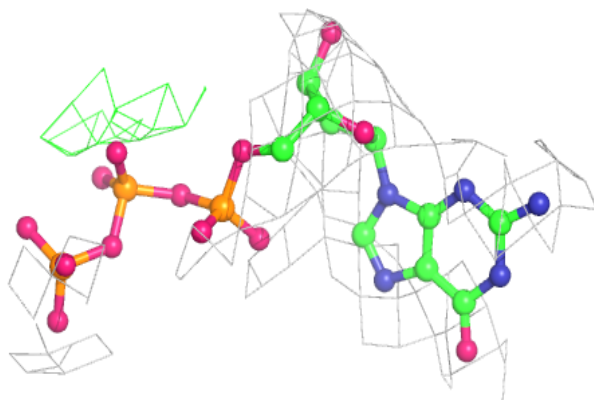


**Electron density around DGT O 702:**

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

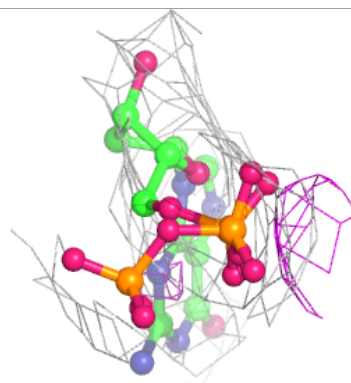
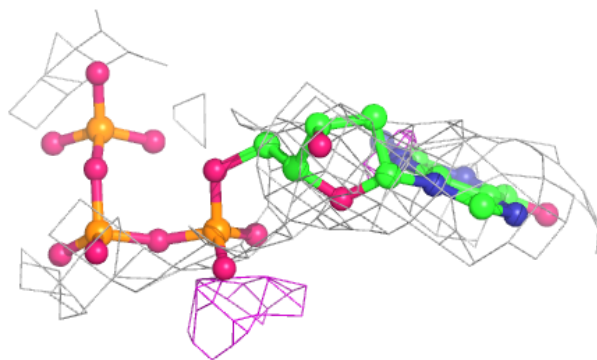
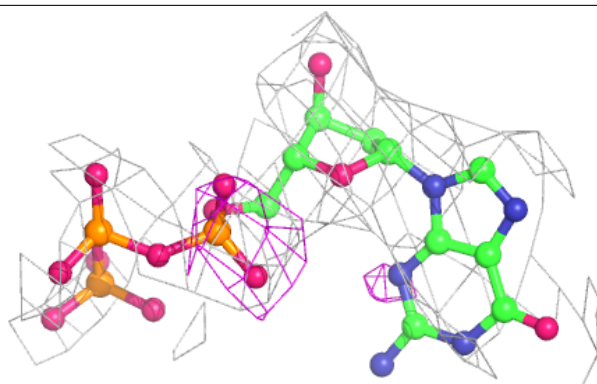
**Electron density around DGT O 703:**

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



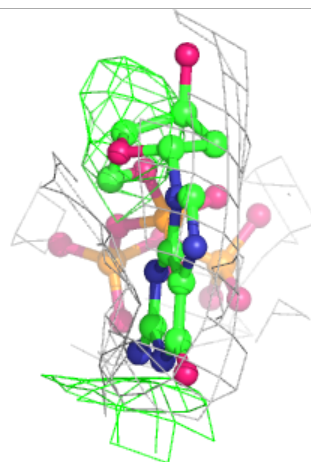
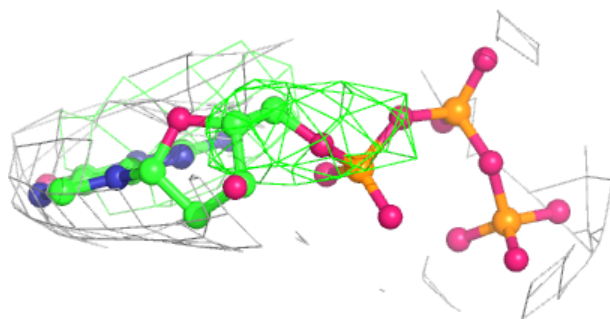
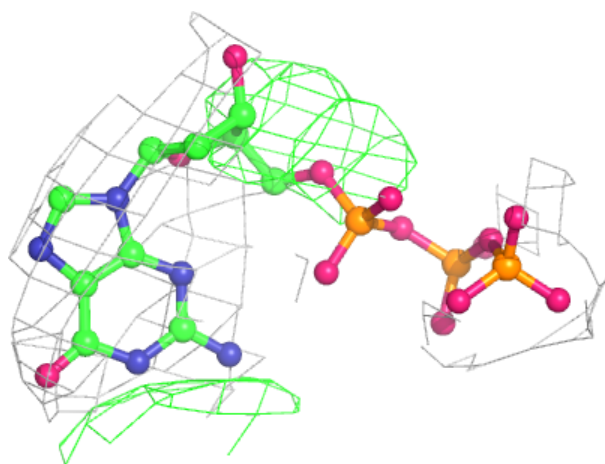
**Electron density around DGT P 701:**

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



**Electron density around DGT P 702:**

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



## 6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.