



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 07:39 PM EDT

PDB ID : 6Z86 / pdb_00006z86
Title : human GTP cyclohydrolase I in complex with 7-deaza-GTP
Authors : Ebenhoch, R.; Nar, H.
Deposited on : 2020-06-02
Resolution : 2.21 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

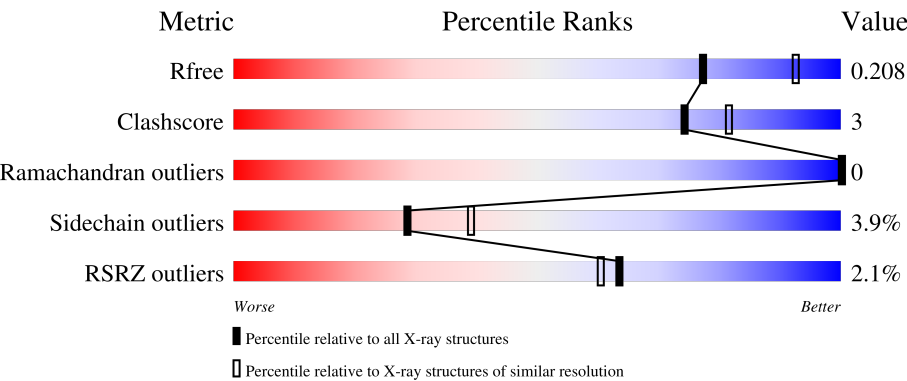
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal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

1 Overall quality at a glance i

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

The reported resolution of this entry is 2.21 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



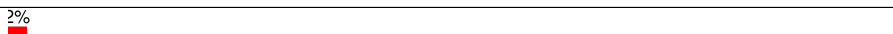
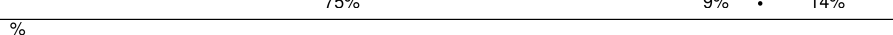
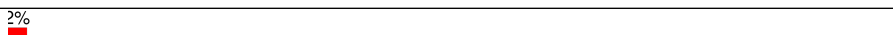
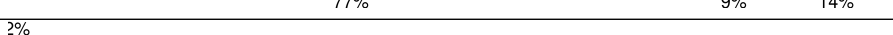
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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

Mol	Chain	Length	Quality of chain
1	A	224	3% 76% 8% 14%
1	B	224	76% 8% 15%
1	C	224	78% 7% 15%
1	D	224	2% 79% 6% 15%
1	E	224	74% 10% 15%

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Mol	Chain	Length	Quality of chain
1	F	224	
1	G	224	
1	H	224	
1	I	224	
1	J	224	
1	K	224	
1	L	224	
1	M	224	
1	N	224	
1	O	224	
1	P	224	
1	Q	224	
1	R	224	
1	S	224	
1	T	224	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 31346 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 GTP cyclohydrolase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	192	Total	C	N	O	S	0	0	0
			1515	956	265	283	11			
1	B	190	Total	C	N	O	S	0	0	0
			1497	945	260	281	11			
1	C	191	Total	C	N	O	S	0	0	0
			1504	950	261	282	11			
1	D	191	Total	C	N	O	S	0	0	0
			1508	951	264	282	11			
1	E	190	Total	C	N	O	S	0	0	0
			1497	945	260	281	11			
1	F	190	Total	C	N	O	S	0	0	0
			1497	945	260	281	11			
1	G	193	Total	C	N	O	S	0	0	0
			1521	959	266	285	11			
1	H	192	Total	C	N	O	S	0	1	0
			1525	960	269	285	11			
1	I	192	Total	C	N	O	S	0	0	0
			1515	956	265	283	11			
1	J	190	Total	C	N	O	S	0	0	0
			1497	945	260	281	11			
1	K	192	Total	C	N	O	S	0	0	0
			1514	954	265	284	11			
1	L	191	Total	C	N	O	S	0	0	0
			1504	950	261	282	11			
1	M	191	Total	C	N	O	S	0	0	0
			1504	950	261	282	11			
1	N	193	Total	C	N	O	S	0	0	0
			1521	959	266	285	11			
1	O	191	Total	C	N	O	S	0	0	0
			1503	948	261	283	11			
1	P	189	Total	C	N	O	S	0	0	0
			1486	939	256	280	11			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	190	Total	C	N	O	S	0	0	0
			1497	945	260	281	11			
1	R	191	Total	C	N	O	S	0	0	0
			1503	948	261	283	11			
1	S	188	Total	C	N	O	S	0	0	0
			1477	933	255	278	11			
1	T	186	Total	C	N	O	S	0	0	0
			1465	928	256	270	11			

There are 280 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	27	MET	-	initiating methionine	UNP P30793
A	28	HIS	-	expression tag	UNP P30793
A	29	HIS	-	expression tag	UNP P30793
A	30	HIS	-	expression tag	UNP P30793
A	31	HIS	-	expression tag	UNP P30793
A	32	HIS	-	expression tag	UNP P30793
A	33	HIS	-	expression tag	UNP P30793
A	34	GLY	-	expression tag	UNP P30793
A	35	SER	-	expression tag	UNP P30793
A	36	ASP	-	expression tag	UNP P30793
A	37	ASP	-	expression tag	UNP P30793
A	38	ASP	-	expression tag	UNP P30793
A	39	ASP	-	expression tag	UNP P30793
A	40	LYS	-	expression tag	UNP P30793
B	27	MET	-	initiating methionine	UNP P30793
B	28	HIS	-	expression tag	UNP P30793
B	29	HIS	-	expression tag	UNP P30793
B	30	HIS	-	expression tag	UNP P30793
B	31	HIS	-	expression tag	UNP P30793
B	32	HIS	-	expression tag	UNP P30793
B	33	HIS	-	expression tag	UNP P30793
B	34	GLY	-	expression tag	UNP P30793
B	35	SER	-	expression tag	UNP P30793
B	36	ASP	-	expression tag	UNP P30793
B	37	ASP	-	expression tag	UNP P30793
B	38	ASP	-	expression tag	UNP P30793
B	39	ASP	-	expression tag	UNP P30793
B	40	LYS	-	expression tag	UNP P30793
C	27	MET	-	initiating methionine	UNP P30793
C	28	HIS	-	expression tag	UNP P30793
C	29	HIS	-	expression tag	UNP P30793

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Chain	Residue	Modelled	Actual	Comment	Reference
C	30	HIS	-	expression tag	UNP P30793
C	31	HIS	-	expression tag	UNP P30793
C	32	HIS	-	expression tag	UNP P30793
C	33	HIS	-	expression tag	UNP P30793
C	34	GLY	-	expression tag	UNP P30793
C	35	SER	-	expression tag	UNP P30793
C	36	ASP	-	expression tag	UNP P30793
C	37	ASP	-	expression tag	UNP P30793
C	38	ASP	-	expression tag	UNP P30793
C	39	ASP	-	expression tag	UNP P30793
C	40	LYS	-	expression tag	UNP P30793
D	27	MET	-	initiating methionine	UNP P30793
D	28	HIS	-	expression tag	UNP P30793
D	29	HIS	-	expression tag	UNP P30793
D	30	HIS	-	expression tag	UNP P30793
D	31	HIS	-	expression tag	UNP P30793
D	32	HIS	-	expression tag	UNP P30793
D	33	HIS	-	expression tag	UNP P30793
D	34	GLY	-	expression tag	UNP P30793
D	35	SER	-	expression tag	UNP P30793
D	36	ASP	-	expression tag	UNP P30793
D	37	ASP	-	expression tag	UNP P30793
D	38	ASP	-	expression tag	UNP P30793
D	39	ASP	-	expression tag	UNP P30793
D	40	LYS	-	expression tag	UNP P30793
E	27	MET	-	initiating methionine	UNP P30793
E	28	HIS	-	expression tag	UNP P30793
E	29	HIS	-	expression tag	UNP P30793
E	30	HIS	-	expression tag	UNP P30793
E	31	HIS	-	expression tag	UNP P30793
E	32	HIS	-	expression tag	UNP P30793
E	33	HIS	-	expression tag	UNP P30793
E	34	GLY	-	expression tag	UNP P30793
E	35	SER	-	expression tag	UNP P30793
E	36	ASP	-	expression tag	UNP P30793
E	37	ASP	-	expression tag	UNP P30793
E	38	ASP	-	expression tag	UNP P30793
E	39	ASP	-	expression tag	UNP P30793
E	40	LYS	-	expression tag	UNP P30793
F	27	MET	-	initiating methionine	UNP P30793
F	28	HIS	-	expression tag	UNP P30793
F	29	HIS	-	expression tag	UNP P30793

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Chain	Residue	Modelled	Actual	Comment	Reference
F	30	HIS	-	expression tag	UNP P30793
F	31	HIS	-	expression tag	UNP P30793
F	32	HIS	-	expression tag	UNP P30793
F	33	HIS	-	expression tag	UNP P30793
F	34	GLY	-	expression tag	UNP P30793
F	35	SER	-	expression tag	UNP P30793
F	36	ASP	-	expression tag	UNP P30793
F	37	ASP	-	expression tag	UNP P30793
F	38	ASP	-	expression tag	UNP P30793
F	39	ASP	-	expression tag	UNP P30793
F	40	LYS	-	expression tag	UNP P30793
G	27	MET	-	initiating methionine	UNP P30793
G	28	HIS	-	expression tag	UNP P30793
G	29	HIS	-	expression tag	UNP P30793
G	30	HIS	-	expression tag	UNP P30793
G	31	HIS	-	expression tag	UNP P30793
G	32	HIS	-	expression tag	UNP P30793
G	33	HIS	-	expression tag	UNP P30793
G	34	GLY	-	expression tag	UNP P30793
G	35	SER	-	expression tag	UNP P30793
G	36	ASP	-	expression tag	UNP P30793
G	37	ASP	-	expression tag	UNP P30793
G	38	ASP	-	expression tag	UNP P30793
G	39	ASP	-	expression tag	UNP P30793
G	40	LYS	-	expression tag	UNP P30793
H	27	MET	-	initiating methionine	UNP P30793
H	28	HIS	-	expression tag	UNP P30793
H	29	HIS	-	expression tag	UNP P30793
H	30	HIS	-	expression tag	UNP P30793
H	31	HIS	-	expression tag	UNP P30793
H	32	HIS	-	expression tag	UNP P30793
H	33	HIS	-	expression tag	UNP P30793
H	34	GLY	-	expression tag	UNP P30793
H	35	SER	-	expression tag	UNP P30793
H	36	ASP	-	expression tag	UNP P30793
H	37	ASP	-	expression tag	UNP P30793
H	38	ASP	-	expression tag	UNP P30793
H	39	ASP	-	expression tag	UNP P30793
H	40	LYS	-	expression tag	UNP P30793
I	27	MET	-	initiating methionine	UNP P30793
I	28	HIS	-	expression tag	UNP P30793
I	29	HIS	-	expression tag	UNP P30793

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Chain	Residue	Modelled	Actual	Comment	Reference
I	30	HIS	-	expression tag	UNP P30793
I	31	HIS	-	expression tag	UNP P30793
I	32	HIS	-	expression tag	UNP P30793
I	33	HIS	-	expression tag	UNP P30793
I	34	GLY	-	expression tag	UNP P30793
I	35	SER	-	expression tag	UNP P30793
I	36	ASP	-	expression tag	UNP P30793
I	37	ASP	-	expression tag	UNP P30793
I	38	ASP	-	expression tag	UNP P30793
I	39	ASP	-	expression tag	UNP P30793
I	40	LYS	-	expression tag	UNP P30793
J	27	MET	-	initiating methionine	UNP P30793
J	28	HIS	-	expression tag	UNP P30793
J	29	HIS	-	expression tag	UNP P30793
J	30	HIS	-	expression tag	UNP P30793
J	31	HIS	-	expression tag	UNP P30793
J	32	HIS	-	expression tag	UNP P30793
J	33	HIS	-	expression tag	UNP P30793
J	34	GLY	-	expression tag	UNP P30793
J	35	SER	-	expression tag	UNP P30793
J	36	ASP	-	expression tag	UNP P30793
J	37	ASP	-	expression tag	UNP P30793
J	38	ASP	-	expression tag	UNP P30793
J	39	ASP	-	expression tag	UNP P30793
J	40	LYS	-	expression tag	UNP P30793
K	27	MET	-	initiating methionine	UNP P30793
K	28	HIS	-	expression tag	UNP P30793
K	29	HIS	-	expression tag	UNP P30793
K	30	HIS	-	expression tag	UNP P30793
K	31	HIS	-	expression tag	UNP P30793
K	32	HIS	-	expression tag	UNP P30793
K	33	HIS	-	expression tag	UNP P30793
K	34	GLY	-	expression tag	UNP P30793
K	35	SER	-	expression tag	UNP P30793
K	36	ASP	-	expression tag	UNP P30793
K	37	ASP	-	expression tag	UNP P30793
K	38	ASP	-	expression tag	UNP P30793
K	39	ASP	-	expression tag	UNP P30793
K	40	LYS	-	expression tag	UNP P30793
L	27	MET	-	initiating methionine	UNP P30793
L	28	HIS	-	expression tag	UNP P30793
L	29	HIS	-	expression tag	UNP P30793

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Chain	Residue	Modelled	Actual	Comment	Reference
L	30	HIS	-	expression tag	UNP P30793
L	31	HIS	-	expression tag	UNP P30793
L	32	HIS	-	expression tag	UNP P30793
L	33	HIS	-	expression tag	UNP P30793
L	34	GLY	-	expression tag	UNP P30793
L	35	SER	-	expression tag	UNP P30793
L	36	ASP	-	expression tag	UNP P30793
L	37	ASP	-	expression tag	UNP P30793
L	38	ASP	-	expression tag	UNP P30793
L	39	ASP	-	expression tag	UNP P30793
L	40	LYS	-	expression tag	UNP P30793
M	27	MET	-	initiating methionine	UNP P30793
M	28	HIS	-	expression tag	UNP P30793
M	29	HIS	-	expression tag	UNP P30793
M	30	HIS	-	expression tag	UNP P30793
M	31	HIS	-	expression tag	UNP P30793
M	32	HIS	-	expression tag	UNP P30793
M	33	HIS	-	expression tag	UNP P30793
M	34	GLY	-	expression tag	UNP P30793
M	35	SER	-	expression tag	UNP P30793
M	36	ASP	-	expression tag	UNP P30793
M	37	ASP	-	expression tag	UNP P30793
M	38	ASP	-	expression tag	UNP P30793
M	39	ASP	-	expression tag	UNP P30793
M	40	LYS	-	expression tag	UNP P30793
N	27	MET	-	initiating methionine	UNP P30793
N	28	HIS	-	expression tag	UNP P30793
N	29	HIS	-	expression tag	UNP P30793
N	30	HIS	-	expression tag	UNP P30793
N	31	HIS	-	expression tag	UNP P30793
N	32	HIS	-	expression tag	UNP P30793
N	33	HIS	-	expression tag	UNP P30793
N	34	GLY	-	expression tag	UNP P30793
N	35	SER	-	expression tag	UNP P30793
N	36	ASP	-	expression tag	UNP P30793
N	37	ASP	-	expression tag	UNP P30793
N	38	ASP	-	expression tag	UNP P30793
N	39	ASP	-	expression tag	UNP P30793
N	40	LYS	-	expression tag	UNP P30793
O	27	MET	-	initiating methionine	UNP P30793
O	28	HIS	-	expression tag	UNP P30793
O	29	HIS	-	expression tag	UNP P30793

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Chain	Residue	Modelled	Actual	Comment	Reference
O	30	HIS	-	expression tag	UNP P30793
O	31	HIS	-	expression tag	UNP P30793
O	32	HIS	-	expression tag	UNP P30793
O	33	HIS	-	expression tag	UNP P30793
O	34	GLY	-	expression tag	UNP P30793
O	35	SER	-	expression tag	UNP P30793
O	36	ASP	-	expression tag	UNP P30793
O	37	ASP	-	expression tag	UNP P30793
O	38	ASP	-	expression tag	UNP P30793
O	39	ASP	-	expression tag	UNP P30793
O	40	LYS	-	expression tag	UNP P30793
P	27	MET	-	initiating methionine	UNP P30793
P	28	HIS	-	expression tag	UNP P30793
P	29	HIS	-	expression tag	UNP P30793
P	30	HIS	-	expression tag	UNP P30793
P	31	HIS	-	expression tag	UNP P30793
P	32	HIS	-	expression tag	UNP P30793
P	33	HIS	-	expression tag	UNP P30793
P	34	GLY	-	expression tag	UNP P30793
P	35	SER	-	expression tag	UNP P30793
P	36	ASP	-	expression tag	UNP P30793
P	37	ASP	-	expression tag	UNP P30793
P	38	ASP	-	expression tag	UNP P30793
P	39	ASP	-	expression tag	UNP P30793
P	40	LYS	-	expression tag	UNP P30793
Q	27	MET	-	initiating methionine	UNP P30793
Q	28	HIS	-	expression tag	UNP P30793
Q	29	HIS	-	expression tag	UNP P30793
Q	30	HIS	-	expression tag	UNP P30793
Q	31	HIS	-	expression tag	UNP P30793
Q	32	HIS	-	expression tag	UNP P30793
Q	33	HIS	-	expression tag	UNP P30793
Q	34	GLY	-	expression tag	UNP P30793
Q	35	SER	-	expression tag	UNP P30793
Q	36	ASP	-	expression tag	UNP P30793
Q	37	ASP	-	expression tag	UNP P30793
Q	38	ASP	-	expression tag	UNP P30793
Q	39	ASP	-	expression tag	UNP P30793
Q	40	LYS	-	expression tag	UNP P30793
R	27	MET	-	initiating methionine	UNP P30793
R	28	HIS	-	expression tag	UNP P30793
R	29	HIS	-	expression tag	UNP P30793

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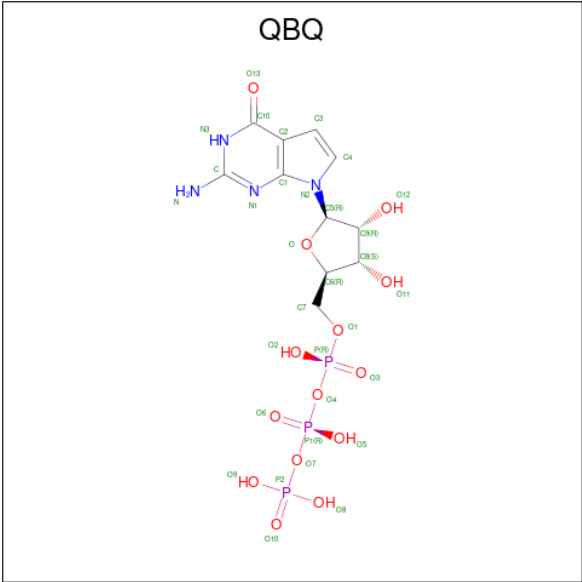
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Chain	Residue	Modelled	Actual	Comment	Reference
R	30	HIS	-	expression tag	UNP P30793
R	31	HIS	-	expression tag	UNP P30793
R	32	HIS	-	expression tag	UNP P30793
R	33	HIS	-	expression tag	UNP P30793
R	34	GLY	-	expression tag	UNP P30793
R	35	SER	-	expression tag	UNP P30793
R	36	ASP	-	expression tag	UNP P30793
R	37	ASP	-	expression tag	UNP P30793
R	38	ASP	-	expression tag	UNP P30793
R	39	ASP	-	expression tag	UNP P30793
R	40	LYS	-	expression tag	UNP P30793
S	27	MET	-	initiating methionine	UNP P30793
S	28	HIS	-	expression tag	UNP P30793
S	29	HIS	-	expression tag	UNP P30793
S	30	HIS	-	expression tag	UNP P30793
S	31	HIS	-	expression tag	UNP P30793
S	32	HIS	-	expression tag	UNP P30793
S	33	HIS	-	expression tag	UNP P30793
S	34	GLY	-	expression tag	UNP P30793
S	35	SER	-	expression tag	UNP P30793
S	36	ASP	-	expression tag	UNP P30793
S	37	ASP	-	expression tag	UNP P30793
S	38	ASP	-	expression tag	UNP P30793
S	39	ASP	-	expression tag	UNP P30793
S	40	LYS	-	expression tag	UNP P30793
T	27	MET	-	initiating methionine	UNP P30793
T	28	HIS	-	expression tag	UNP P30793
T	29	HIS	-	expression tag	UNP P30793
T	30	HIS	-	expression tag	UNP P30793
T	31	HIS	-	expression tag	UNP P30793
T	32	HIS	-	expression tag	UNP P30793
T	33	HIS	-	expression tag	UNP P30793
T	34	GLY	-	expression tag	UNP P30793
T	35	SER	-	expression tag	UNP P30793
T	36	ASP	-	expression tag	UNP P30793
T	37	ASP	-	expression tag	UNP P30793
T	38	ASP	-	expression tag	UNP P30793
T	39	ASP	-	expression tag	UNP P30793
T	40	LYS	-	expression tag	UNP P30793

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0
2	F	1	Total Zn 1 1	0	0
2	G	1	Total Zn 1 1	0	0
2	H	1	Total Zn 1 1	0	0
2	I	1	Total Zn 1 1	0	0
2	J	1	Total Zn 1 1	0	0
2	K	1	Total Zn 1 1	0	0
2	L	1	Total Zn 1 1	0	0
2	M	1	Total Zn 1 1	0	0
2	N	1	Total Zn 1 1	0	0
2	O	1	Total Zn 1 1	0	0
2	P	1	Total Zn 1 1	0	0
2	Q	1	Total Zn 1 1	0	0
2	R	1	Total Zn 1 1	0	0
2	S	1	Total Zn 1 1	0	0
2	T	1	Total Zn 1 1	0	0

- Molecule 3 is 7-deaza-GTP (CCD ID: QBQ) (formula: $C_{11}H_{17}N_4O_{14}P_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	B	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	C	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	D	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	E	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	F	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	G	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	H	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	I	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	J	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	K	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	L	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	M	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	N	1	Total	C	N	O	P	0	0
			32	11	4	14	3		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	O	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	P	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	Q	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	R	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	S	1	Total	C	N	O	P	0	0
			32	11	4	14	3		
3	T	1	Total	C	N	O	P	0	0
			32	11	4	14	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	34	Total	O	0	0
			34	34		
4	B	46	Total	O	0	0
			46	46		
4	C	30	Total	O	0	0
			30	30		
4	D	32	Total	O	0	0
			32	32		
4	E	35	Total	O	0	0
			35	35		
4	F	38	Total	O	0	0
			38	38		
4	G	33	Total	O	0	0
			33	33		
4	H	29	Total	O	0	0
			29	29		
4	I	34	Total	O	0	0
			34	34		
4	J	33	Total	O	0	0
			33	33		
4	K	34	Total	O	0	0
			34	34		
4	L	24	Total	O	0	0
			24	24		
4	M	31	Total	O	0	0
			31	31		

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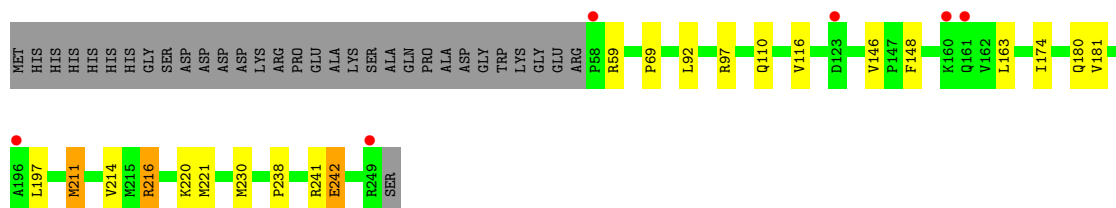
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	N	34	Total 34	O 34	0	0
4	O	30	Total 30	O 30	0	0
4	P	19	Total 19	O 19	0	0
4	Q	33	Total 33	O 33	0	0
4	R	32	Total 32	O 32	0	0
4	S	24	Total 24	O 24	0	0
4	T	31	Total 31	O 31	0	0

3 Residue-property plots [i](#)

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

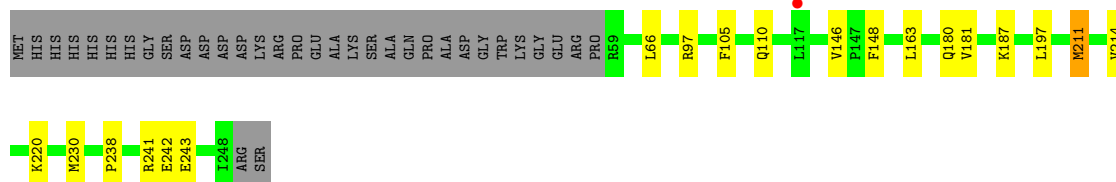
• Molecule 1: GTP cyclohydrolase 1

Chain A: 



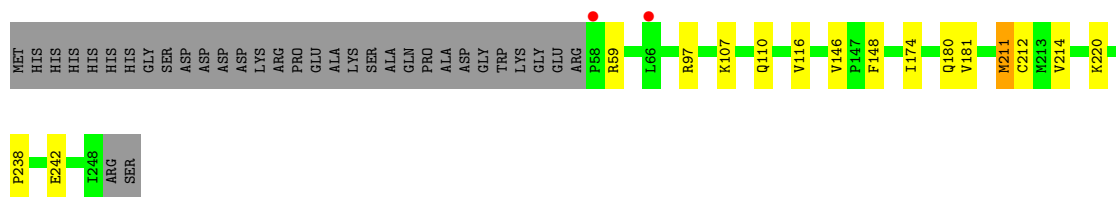
• Molecule 1: GTP cyclohydrolase 1

Chain B: 



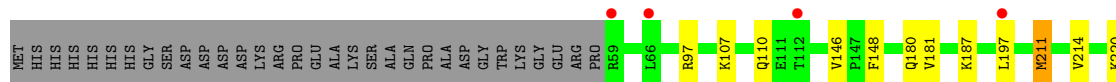
• Molecule 1: GTP cyclohydrolase 1

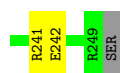
Chain C: 



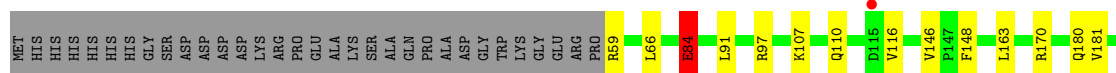
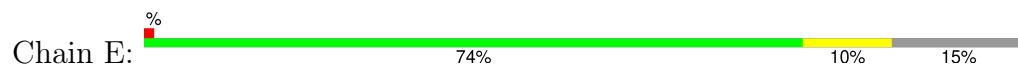
• Molecule 1: GTP cyclohydrolase 1

Chain D: 

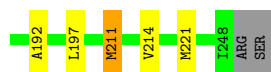
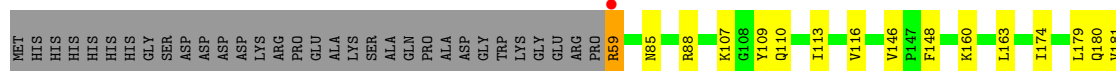




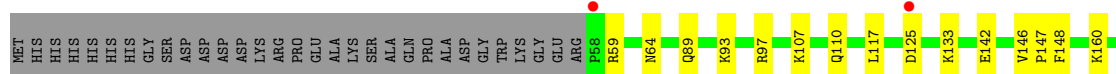
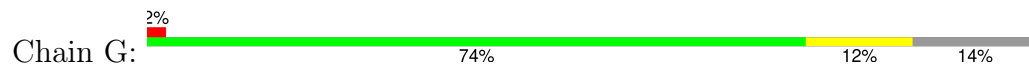
- Molecule 1: GTP cyclohydrolase 1



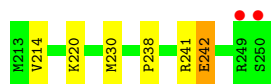
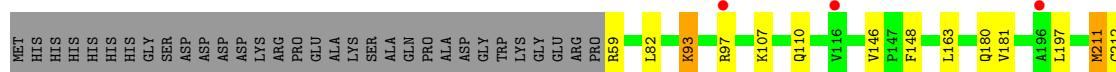
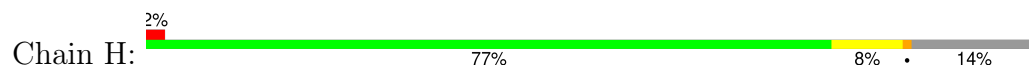
- Molecule 1: GTP cyclohydrolase 1



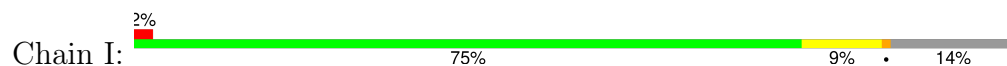
- Molecule 1: GTP cyclohydrolase 1

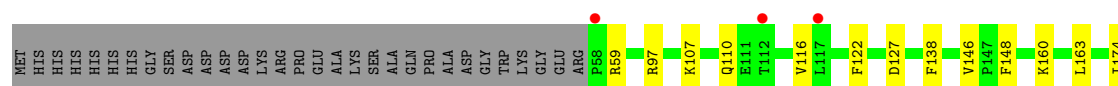


- Molecule 1: GTP cyclohydrolase 1

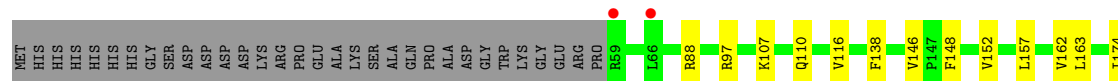
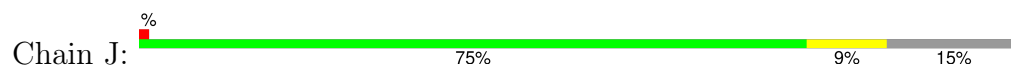


- Molecule 1: GTP cyclohydrolase 1

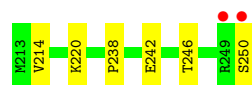
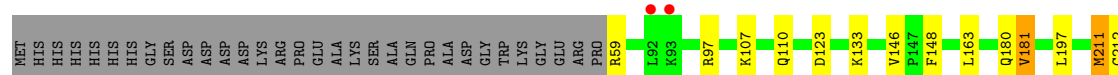
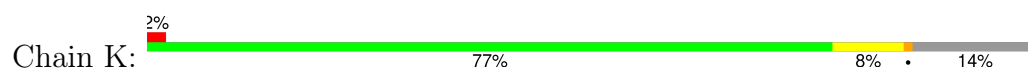




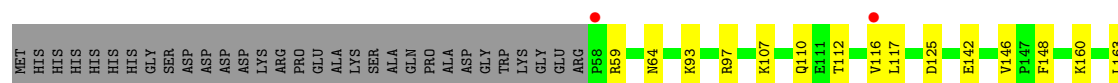
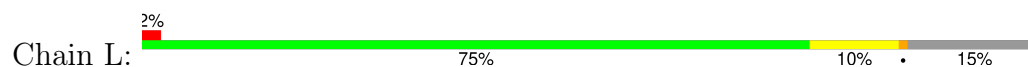
• Molecule 1: GTP cyclohydrolase 1



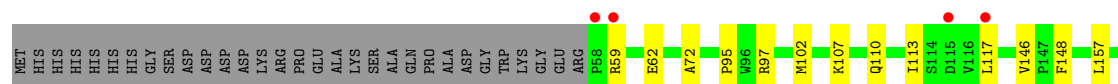
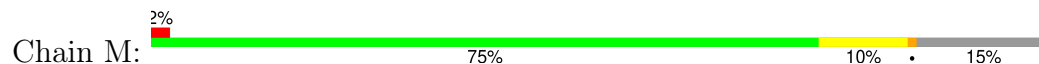
• Molecule 1: GTP cyclohydrolase 1



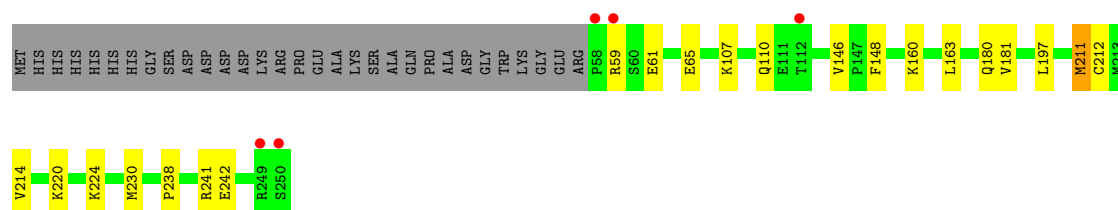
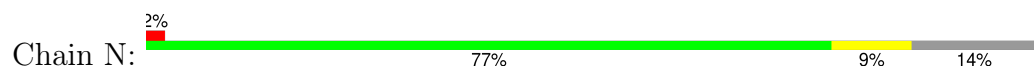
• Molecule 1: GTP cyclohydrolase 1



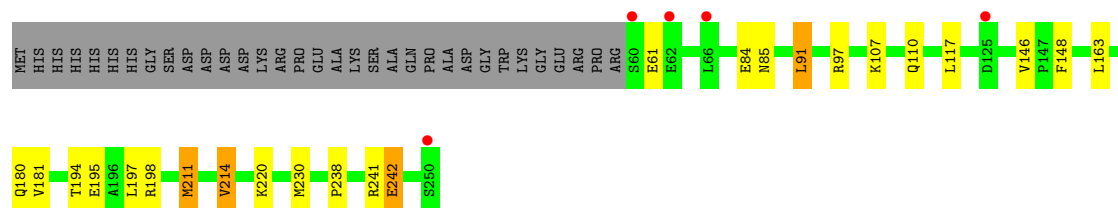
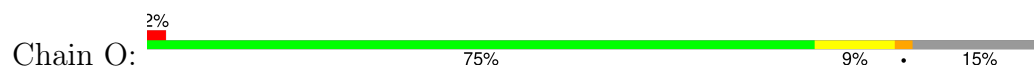
• Molecule 1: GTP cyclohydrolase 1



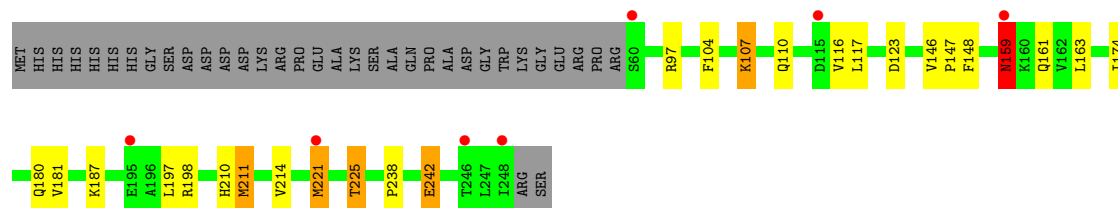
- Molecule 1: GTP cyclohydrolase 1



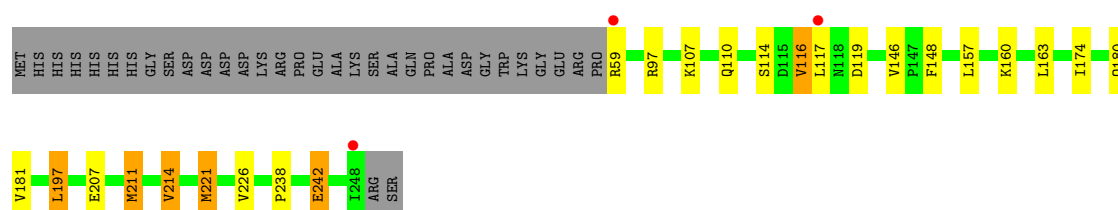
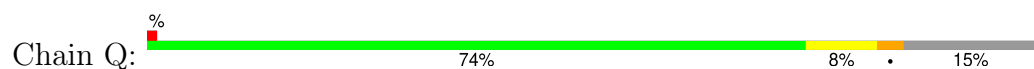
- Molecule 1: GTP cyclohydrolase 1



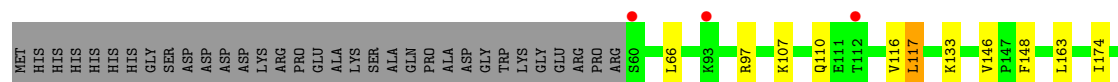
- Molecule 1: GTP cyclohydrolase 1

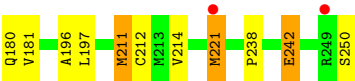


- Molecule 1: GTP cyclohydrolase 1

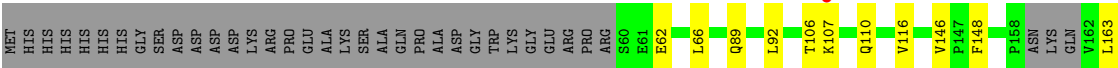


- Molecule 1: GTP cyclohydrolase 1

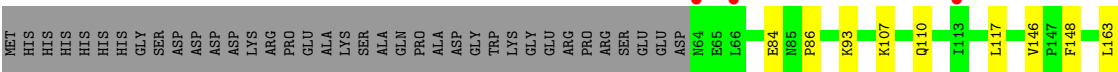




● Molecule 1: GTP cyclohydrolase 1



● Molecule 1: GTP cyclohydrolase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	87.86Å 88.59Å 163.57Å 85.22° 88.91° 83.55°	Depositor
Resolution (Å)	162.99 – 2.21 162.99 – 2.21	Depositor EDS
% Data completeness (in resolution range)	57.9 (162.99-2.21) 57.9 (162.99-2.21)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.29 (at 2.20Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.183 , 0.208 0.185 , 0.208	Depositor DCC
R_{free} test set	2913 reflections (2.04%)	wwPDB-VP
Wilson B-factor (Å ²)	42.4	Xtriage
Anisotropy	0.020	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.003 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	31346	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.76	1/1540 (0.1%)	1.09	4/2079 (0.2%)
1	B	0.75	0/1521	1.13	4/2054 (0.2%)
1	C	0.73	0/1529	1.10	4/2065 (0.2%)
1	D	0.72	0/1532	1.09	4/2068 (0.2%)
1	E	0.74	0/1521	1.11	5/2054 (0.2%)
1	F	0.79	0/1521	1.11	3/2054 (0.1%)
1	G	0.73	0/1546	1.10	6/2087 (0.3%)
1	H	0.73	0/1549	1.10	4/2090 (0.2%)
1	I	0.73	0/1540	1.09	3/2079 (0.1%)
1	J	0.74	0/1521	1.09	4/2054 (0.2%)
1	K	0.75	0/1538	1.10	5/2076 (0.2%)
1	L	0.70	0/1529	1.09	4/2065 (0.2%)
1	M	0.74	0/1529	1.12	4/2065 (0.2%)
1	N	0.74	0/1546	1.11	4/2087 (0.2%)
1	O	0.73	0/1527	1.11	5/2062 (0.2%)
1	P	0.69	0/1510	1.13	5/2040 (0.2%)
1	Q	0.69	0/1521	1.08	3/2054 (0.1%)
1	R	0.72	0/1527	1.10	3/2062 (0.1%)
1	S	0.75	0/1500	1.15	6/2025 (0.3%)
1	T	0.72	0/1489	1.09	5/2011 (0.2%)
All	All	0.73	1/30536 (0.0%)	1.11	85/41231 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	221	MET	SD-CE	5.59	1.93	1.79

All (85) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	P	159	ASN	CA-CB-CG	8.42	121.02	112.60
1	C	220	LYS	CA-C-N	8.41	131.90	120.54
1	C	220	LYS	C-N-CA	8.41	131.90	120.54
1	O	220	LYS	CA-C-N	8.08	131.44	120.54
1	O	220	LYS	C-N-CA	8.08	131.44	120.54
1	B	220	LYS	CA-C-N	8.04	131.40	120.54
1	B	220	LYS	C-N-CA	8.04	131.40	120.54
1	M	220	LYS	CA-C-N	7.84	131.13	120.54
1	M	220	LYS	C-N-CA	7.84	131.13	120.54
1	G	220	LYS	CA-C-N	7.71	130.94	120.54
1	G	220	LYS	C-N-CA	7.71	130.94	120.54
1	E	220	LYS	CA-C-N	7.64	130.85	120.54
1	E	220	LYS	C-N-CA	7.64	130.85	120.54
1	N	220	LYS	CA-C-N	7.43	131.96	120.82
1	N	220	LYS	C-N-CA	7.43	131.96	120.82
1	S	220	LYS	CA-C-N	7.35	130.46	120.54
1	S	220	LYS	C-N-CA	7.35	130.46	120.54
1	H	220	LYS	CA-C-N	7.32	132.83	121.19
1	H	220	LYS	C-N-CA	7.32	132.83	121.19
1	A	220	LYS	CA-C-N	6.63	130.77	120.82
1	A	220	LYS	C-N-CA	6.63	130.77	120.82
1	L	220	LYS	CA-C-N	6.42	129.21	120.54
1	L	220	LYS	C-N-CA	6.42	129.21	120.54
1	E	148	PHE	CA-CB-CG	6.39	120.19	113.80
1	L	148	PHE	CA-CB-CG	6.35	120.15	113.80
1	F	148	PHE	CA-CB-CG	6.35	120.15	113.80
1	J	220	LYS	CA-C-N	6.34	129.11	120.54
1	J	220	LYS	C-N-CA	6.34	129.11	120.54
1	D	148	PHE	CA-CB-CG	6.21	120.01	113.80
1	E	84	GLU	CB-CG-CD	6.21	123.15	112.60
1	S	148	PHE	CA-CB-CG	6.16	119.96	113.80
1	C	148	PHE	CA-CB-CG	6.15	119.95	113.80
1	G	148	PHE	CA-CB-CG	6.12	119.92	113.80
1	I	148	PHE	CA-CB-CG	6.12	119.92	113.80
1	P	148	PHE	CA-CB-CG	6.10	119.90	113.80
1	H	148	PHE	CA-CB-CG	6.10	119.90	113.80
1	R	148	PHE	CA-CB-CG	6.08	119.88	113.80
1	K	148	PHE	CA-CB-CG	6.05	119.85	113.80
1	T	148	PHE	CA-CB-CG	6.02	119.82	113.80
1	G	180	GLN	N-CA-C	6.00	118.53	109.41
1	T	180	GLN	N-CA-C	5.99	118.51	109.41
1	H	180	GLN	N-CA-C	5.98	118.50	109.41
1	S	180	GLN	N-CA-C	5.96	118.47	109.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	180	GLN	N-CA-C	5.95	118.45	109.41
1	D	220	LYS	CA-C-N	5.92	129.69	120.82
1	D	220	LYS	C-N-CA	5.92	129.69	120.82
1	O	148	PHE	CA-CB-CG	5.85	119.65	113.80
1	I	180	GLN	N-CA-C	5.84	118.29	109.41
1	F	180	GLN	N-CA-C	5.81	118.23	109.41
1	P	180	GLN	N-CA-C	5.80	118.23	109.41
1	J	148	PHE	CA-CB-CG	5.79	119.59	113.80
1	E	180	GLN	N-CA-C	5.78	118.20	109.41
1	Q	148	PHE	CA-CB-CG	5.77	119.57	113.80
1	O	180	GLN	N-CA-C	5.77	118.18	109.41
1	C	180	GLN	N-CA-C	5.76	118.17	109.41
1	N	180	GLN	N-CA-C	5.71	118.09	109.41
1	D	180	GLN	N-CA-C	5.70	118.08	109.41
1	O	214	VAL	N-CA-CB	5.66	117.64	111.41
1	M	148	PHE	CA-CB-CG	5.64	119.44	113.80
1	R	180	GLN	N-CA-C	5.64	117.98	109.41
1	A	148	PHE	CA-CB-CG	5.64	119.44	113.80
1	T	216	ARG	N-CA-CB	-5.63	102.11	111.57
1	N	148	PHE	CA-CB-CG	5.62	119.42	113.80
1	Q	180	GLN	N-CA-C	5.62	117.95	109.41
1	B	148	PHE	CA-CB-CG	5.61	119.41	113.80
1	M	180	GLN	N-CA-C	5.59	117.90	109.41
1	A	180	GLN	N-CA-C	5.58	117.89	109.41
1	B	180	GLN	N-CA-C	5.55	117.84	109.41
1	R	221	MET	N-CA-C	5.53	117.31	111.28
1	T	220	LYS	CA-C-N	5.53	128.00	120.54
1	T	220	LYS	C-N-CA	5.53	128.00	120.54
1	K	180	GLN	N-CA-C	5.49	117.75	109.41
1	S	106	THR	CA-C-N	5.43	128.31	120.38
1	S	106	THR	C-N-CA	5.43	128.31	120.38
1	Q	221	MET	CB-CG-SD	-5.38	96.57	112.70
1	K	220	LYS	CA-C-N	5.37	128.88	120.82
1	K	220	LYS	C-N-CA	5.37	128.88	120.82
1	J	180	GLN	N-CA-C	5.35	117.54	109.41
1	K	181	VAL	CB-CA-C	5.33	118.12	110.33
1	G	180	GLN	CA-C-N	5.31	130.37	122.99
1	G	180	GLN	C-N-CA	5.31	130.37	122.99
1	I	221	MET	CB-CG-SD	-5.08	97.47	112.70
1	P	221	MET	N-CA-C	5.06	116.80	111.28
1	P	159	ASN	CB-CA-C	5.05	119.56	111.73
1	F	179	LEU	N-CA-C	-5.05	101.52	109.25

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	1515	0	1539	9	0
1	B	1497	0	1518	6	0
1	C	1504	0	1526	7	0
1	D	1508	0	1531	3	0
1	E	1497	0	1518	14	0
1	F	1497	0	1518	12	0
1	G	1521	0	1544	15	0
1	H	1525	0	1548	11	0
1	I	1515	0	1539	12	0
1	J	1497	0	1518	9	0
1	K	1514	0	1536	12	0
1	L	1504	0	1526	12	0
1	M	1504	0	1526	11	0
1	N	1521	0	1544	8	0
1	O	1503	0	1523	11	0
1	P	1486	0	1505	13	0
1	Q	1497	0	1518	11	0
1	R	1503	0	1523	10	0
1	S	1477	0	1495	8	0
1	T	1465	0	1497	8	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0
2	I	1	0	0	0	0
2	J	1	0	0	0	0
2	K	1	0	0	0	0
2	L	1	0	0	0	0
2	M	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	N	1	0	0	0	0
2	O	1	0	0	0	0
2	P	1	0	0	0	0
2	Q	1	0	0	0	0
2	R	1	0	0	0	0
2	S	1	0	0	0	0
2	T	1	0	0	0	0
3	A	32	0	0	0	0
3	B	32	0	0	0	0
3	C	32	0	0	1	0
3	D	32	0	0	0	0
3	E	32	0	0	0	0
3	F	32	0	0	0	0
3	G	32	0	0	0	0
3	H	32	0	0	0	0
3	I	32	0	0	0	0
3	J	32	0	0	0	0
3	K	32	0	0	0	0
3	L	32	0	0	0	0
3	M	32	0	0	0	0
3	N	32	0	0	0	0
3	O	32	0	0	0	0
3	P	32	0	0	0	0
3	Q	32	0	0	1	0
3	R	32	0	0	0	0
3	S	32	0	0	0	0
3	T	32	0	0	0	0
4	A	34	0	0	1	0
4	B	46	0	0	0	0
4	C	30	0	0	0	0
4	D	32	0	0	0	0
4	E	35	0	0	1	0
4	F	38	0	0	0	0
4	G	33	0	0	1	0
4	H	29	0	0	0	0
4	I	34	0	0	2	0
4	J	33	0	0	1	0
4	K	34	0	0	0	0
4	L	24	0	0	1	0
4	M	31	0	0	1	0
4	N	34	0	0	2	0
4	O	30	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	P	19	0	0	1	0
4	Q	33	0	0	0	0
4	R	32	0	0	0	0
4	S	24	0	0	0	0
4	T	31	0	0	1	0
All	All	31346	0	30492	190	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (190) 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:R:133:LYS:NZ	1:R:250:SER:HB3	1.84	0.92
1:E:84:GLU:HG3	1:E:91:LEU:HD11	1.58	0.85
1:R:133:LYS:HZ2	1:R:250:SER:HB3	1.41	0.84
1:O:238:PRO:O	1:O:242:GLU:HG2	1.86	0.76
1:E:163:LEU:HB2	1:E:197:LEU:HD11	1.66	0.75
1:G:238:PRO:O	1:G:242:GLU:HG2	1.87	0.74
1:M:238:PRO:O	1:M:242:GLU:HG2	1.87	0.74
1:T:238:PRO:O	1:T:242:GLU:HG2	1.87	0.74
1:P:238:PRO:O	1:P:242:GLU:HG2	1.86	0.74
1:A:238:PRO:O	1:A:242:GLU:HG2	1.88	0.73
1:Q:238:PRO:O	1:Q:242:GLU:HG2	1.87	0.73
1:H:238:PRO:O	1:H:242:GLU:HG2	1.87	0.73
1:G:163:LEU:HB2	1:G:197:LEU:HD11	1.71	0.73
1:O:84:GLU:HG2	1:O:91:LEU:HD22	1.71	0.73
1:O:84:GLU:HG3	1:O:85:ASN:H	1.54	0.72
1:E:84:GLU:HG3	1:E:91:LEU:CD1	2.20	0.71
1:N:163:LEU:HB2	1:N:197:LEU:HD11	1.73	0.71
1:R:238:PRO:O	1:R:242:GLU:HG2	1.89	0.71
1:F:59:ARG:HD2	1:F:109:TYR:HB2	1.72	0.70
1:F:113:ILE:HD12	1:F:192:ALA:HB1	1.73	0.70
1:K:133:LYS:HZ2	1:K:250:SER:HB3	1.56	0.70
1:M:163:LEU:HB2	1:M:197:LEU:HD11	1.74	0.69
1:H:93:LYS:NZ	1:H:97[A]:ARG:HH22	1.90	0.69
1:M:113:ILE:HD12	1:M:192:ALA:HB1	1.75	0.69
1:H:93:LYS:HZ1	1:H:97[A]:ARG:HH22	1.40	0.68
1:A:163:LEU:HB2	1:A:197:LEU:HD11	1.75	0.67
1:L:163:LEU:HB2	1:L:197:LEU:HD11	1.76	0.67
1:F:85:ASN:O	1:F:88:ARG:HG2	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:107:LYS:O	1:D:110:GLN:HG2	1.94	0.67
1:O:84:GLU:HG2	1:O:91:LEU:CD2	2.25	0.67
1:O:163:LEU:HB2	1:O:197:LEU:HD11	1.77	0.66
1:P:104:PHE:O	1:P:107:LYS:HG3	1.97	0.65
1:P:163:LEU:HB2	1:P:197:LEU:HD11	1.79	0.65
1:E:238:PRO:O	1:E:242:GLU:HG3	1.97	0.64
1:R:133:LYS:HZ1	1:R:250:SER:HB3	1.62	0.64
1:B:238:PRO:O	1:B:242:GLU:HG3	1.98	0.64
1:O:84:GLU:HG3	1:O:85:ASN:N	2.12	0.64
1:F:88:ARG:NH2	1:G:142:GLU:OE2	2.28	0.64
1:R:163:LEU:HB2	1:R:197:LEU:HD11	1.81	0.63
1:F:163:LEU:HB2	1:F:197:LEU:HD11	1.80	0.63
1:M:107:LYS:O	1:M:110:GLN:HG2	1.98	0.62
4:I:420:HOH:O	1:J:221:MET:HE3	1.99	0.61
1:K:163:LEU:HB2	1:K:197:LEU:HD11	1.81	0.61
1:C:238:PRO:O	1:C:242:GLU:HG3	2.01	0.60
1:H:59:ARG:HB3	1:H:110:GLN:HB3	1.84	0.60
1:N:238:PRO:O	1:N:242:GLU:HG3	2.01	0.60
1:K:133:LYS:NZ	1:K:250:SER:HB3	2.17	0.60
1:O:195:GLU:HA	1:O:198:ARG:HE	1.66	0.60
1:J:163:LEU:HB2	1:J:197:LEU:HD11	1.84	0.59
1:E:84:GLU:CG	1:E:91:LEU:HD11	2.32	0.59
1:M:59:ARG:HB3	1:M:110:GLN:HB3	1.84	0.59
1:B:163:LEU:HB2	1:B:197:LEU:HD11	1.84	0.58
1:O:117:LEU:HD22	1:O:163:LEU:HD11	1.85	0.58
1:C:59:ARG:HB3	1:C:110:GLN:HB3	1.86	0.58
1:I:163:LEU:HB2	1:I:197:LEU:HD11	1.86	0.58
1:K:59:ARG:HB3	1:K:110:GLN:HB3	1.86	0.58
1:Q:59:ARG:HB3	1:Q:110:GLN:HB3	1.86	0.57
1:L:59:ARG:O	1:L:110:GLN:HB3	2.04	0.57
1:S:163:LEU:HB2	1:S:197:LEU:HD11	1.87	0.56
1:E:107:LYS:O	1:E:110:GLN:HG2	2.06	0.55
4:E:401:HOH:O	1:J:187:LYS:HE2	2.06	0.55
1:G:107:LYS:O	1:G:110:GLN:HG2	2.06	0.55
1:P:225:THR:HG21	1:S:165:LEU:HD13	1.88	0.55
1:L:107:LYS:O	1:L:110:GLN:HG2	2.07	0.55
1:C:107:LYS:O	1:C:110:GLN:HG2	2.07	0.55
1:T:163:LEU:HB2	1:T:197:LEU:HD11	1.89	0.55
1:L:59:ARG:HD3	1:L:64:ASN:OD1	2.06	0.54
1:N:107:LYS:O	1:N:110:GLN:HG2	2.08	0.54
1:H:163:LEU:HB2	1:H:197:LEU:HD11	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:107:LYS:O	1:I:110:GLN:HG2	2.08	0.54
1:A:59:ARG:HB3	1:A:110:GLN:HG2	1.89	0.54
1:E:59:ARG:HB3	1:E:110:GLN:HB3	1.88	0.54
1:L:59:ARG:O	1:L:110:GLN:CB	2.56	0.53
1:P:107:LYS:O	1:P:110:GLN:HG2	2.08	0.53
1:S:107:LYS:O	1:S:110:GLN:HG2	2.09	0.53
4:G:432:HOH:O	1:N:212:CYS:HB2	2.09	0.53
1:G:59:ARG:HD2	1:G:110:GLN:OE1	2.08	0.52
1:P:147:PRO:HG2	1:T:211:MET:HE3	1.92	0.52
1:H:107:LYS:O	1:H:110:GLN:HG2	2.08	0.52
1:G:146:VAL:HG12	1:G:211:MET:HB2	1.91	0.52
1:Q:116:VAL:HG21	1:Q:174:ILE:HD11	1.92	0.52
1:I:59:ARG:HB3	1:I:110:GLN:HB3	1.91	0.52
1:J:138:PHE:HZ	1:L:221:MET:SD	2.33	0.52
1:G:59:ARG:HB3	1:G:110:GLN:HB3	1.91	0.51
1:P:117:LEU:HD22	1:P:163:LEU:HD11	1.91	0.51
1:J:107:LYS:O	1:J:110:GLN:HG2	2.11	0.51
1:S:239:LYS:HA	1:S:242:GLU:HG2	1.92	0.51
1:H:146:VAL:HG12	1:H:211:MET:HB2	1.92	0.51
1:T:84:GLU:O	1:T:86:PRO:HD3	2.11	0.51
1:O:107:LYS:O	1:O:110:GLN:HG2	2.11	0.51
1:F:107:LYS:O	1:F:110:GLN:HG2	2.11	0.51
1:F:113:ILE:CD1	1:F:192:ALA:HB1	2.40	0.51
1:Q:107:LYS:O	1:Q:110:GLN:HG2	2.11	0.50
1:R:107:LYS:O	1:R:110:GLN:HG2	2.10	0.50
1:K:107:LYS:O	1:K:110:GLN:HG2	2.11	0.50
1:C:212:CYS:HB2	4:P:418:HOH:O	2.12	0.50
1:B:146:VAL:HG12	1:B:211:MET:HB2	1.94	0.50
1:R:146:VAL:HG12	1:R:211:MET:HB2	1.94	0.50
1:T:107:LYS:O	1:T:110:GLN:HG2	2.11	0.50
1:I:146:VAL:HG12	1:I:211:MET:HB2	1.93	0.49
1:I:214:VAL:HB	1:I:221:MET:HA	1.95	0.49
1:T:146:VAL:HG12	1:T:211:MET:HB2	1.93	0.49
1:E:187:LYS:HE2	4:N:412:HOH:O	2.12	0.49
1:K:238:PRO:O	1:K:242:GLU:HG2	2.12	0.49
1:L:146:VAL:HG12	1:L:211:MET:HB2	1.95	0.49
1:N:146:VAL:HG12	1:N:211:MET:HB2	1.93	0.49
1:J:88:ARG:NH2	1:L:142:GLU:OE2	2.44	0.49
1:F:146:VAL:HG12	1:F:211:MET:HB2	1.95	0.49
1:K:146:VAL:HB	1:K:212:CYS:SG	2.52	0.48
1:E:146:VAL:HG12	1:E:211:MET:HB2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:210:HIS:CD2	1:P:225:THR:CG2	2.96	0.48
1:G:133:LYS:NZ	1:G:250:SER:HB3	2.28	0.48
1:D:146:VAL:HG12	1:D:211:MET:HB2	1.95	0.48
1:G:89:GLN:NE2	1:K:123:ASP:OD2	2.38	0.48
1:M:146:VAL:HG12	1:M:211:MET:HB2	1.96	0.48
1:O:146:VAL:HG12	1:O:211:MET:HB2	1.96	0.48
3:Q:302:QBQ:O8	1:S:216:ARG:NH1	2.41	0.48
1:A:216:ARG:HG2	1:B:105:PHE:CE1	2.48	0.47
1:M:113:ILE:CD1	1:M:192:ALA:HB1	2.44	0.47
1:S:146:VAL:HG12	1:S:211:MET:HB2	1.96	0.47
1:A:146:VAL:HG12	1:A:211:MET:HB2	1.95	0.47
1:J:146:VAL:HG12	1:J:211:MET:HB2	1.96	0.47
1:E:212:CYS:HB2	4:N:430:HOH:O	2.15	0.47
1:P:146:VAL:HG12	1:P:211:MET:HB2	1.95	0.46
1:L:116:VAL:HG21	1:L:174:ILE:HD11	1.97	0.46
1:Q:146:VAL:HG12	1:Q:211:MET:HB2	1.96	0.46
1:P:159:ASN:O	1:P:159:ASN:OD1	2.33	0.46
1:S:89:GLN:HA	1:S:92:LEU:HD12	1.98	0.46
1:H:212:CYS:HB2	4:J:431:HOH:O	2.15	0.46
1:R:117:LEU:HD21	1:R:196:ALA:HB1	1.98	0.46
1:C:146:VAL:HG12	1:C:211:MET:HB2	1.96	0.46
1:N:61:GLU:O	1:N:65:GLU:HG2	2.16	0.45
1:Q:117:LEU:HD12	1:Q:119:ASP:H	1.81	0.45
1:I:216:ARG:HG2	1:I:217:GLY:H	1.82	0.45
1:K:146:VAL:HG12	1:K:211:MET:HB2	1.98	0.45
1:K:133:LYS:NZ	1:K:250:SER:CB	2.79	0.45
1:G:117:LEU:HD21	1:G:196:ALA:HB1	1.99	0.45
1:M:72:ALA:O	1:M:95:PRO:HB3	2.16	0.45
1:A:69:PRO:HG2	4:A:402:HOH:O	2.16	0.44
1:E:221:MET:SD	1:I:138:PHE:HZ	2.40	0.44
1:C:59:ARG:O	1:C:110:GLN:HB2	2.18	0.44
1:F:160:LYS:HA	1:F:160:LYS:HD3	1.90	0.43
1:I:59:ARG:O	1:I:110:GLN:HB2	2.19	0.43
1:L:211:MET:HA	1:L:214:VAL:HG13	2.00	0.43
1:G:211:MET:O	1:G:215:MET:HG2	2.19	0.43
1:L:125:ASP:HA	1:L:160:LYS:HD2	2.01	0.43
1:A:116:VAL:HG21	1:A:174:ILE:HD11	2.01	0.43
4:I:432:HOH:O	1:K:212:CYS:HB2	2.18	0.43
1:F:211:MET:HE3	1:G:147:PRO:HG2	2.00	0.43
1:F:211:MET:HE1	1:G:211:MET:HG3	2.01	0.42
1:G:59:ARG:HD3	1:G:64:ASN:OD1	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:243:GLU:OE2	1:D:241:ARG:HD2	2.19	0.42
1:L:212:CYS:HB2	4:M:425:HOH:O	2.19	0.42
1:J:116:VAL:HG21	1:J:174:ILE:HD11	2.01	0.42
1:K:242:GLU:O	1:K:246:THR:HG23	2.20	0.42
1:F:116:VAL:HG21	1:F:174:ILE:HD11	2.01	0.42
1:M:117:LEU:HD22	1:M:163:LEU:HD11	2.02	0.42
1:G:211:MET:HA	1:G:211:MET:HE2	2.00	0.42
1:N:160:LYS:HA	1:N:160:LYS:HD3	1.92	0.42
1:A:216:ARG:NH1	3:C:302:QBQ:O8	2.52	0.41
1:E:116:VAL:HG13	1:E:170:ARG:HB3	2.02	0.41
1:A:230:MET:HG3	1:A:241:ARG:CZ	2.51	0.41
1:I:127:ASP:HB3	1:I:160:LYS:NZ	2.35	0.41
1:R:212:CYS:HB2	4:T:429:HOH:O	2.19	0.41
1:I:116:VAL:HG21	1:I:174:ILE:HD11	2.02	0.41
1:Q:207:GLU:HG3	1:Q:226:VAL:HG22	2.02	0.41
1:H:230:MET:HG3	1:H:241:ARG:CZ	2.51	0.41
1:I:212:CYS:HB2	4:L:422:HOH:O	2.20	0.41
1:J:152:VAL:HG22	1:J:206:VAL:HG22	2.02	0.41
1:N:230:MET:HG3	1:N:241:ARG:CZ	2.50	0.41
1:O:230:MET:HG3	1:O:241:ARG:CZ	2.50	0.41
1:Q:214:VAL:HB	1:Q:221:MET:HA	2.02	0.41
1:C:116:VAL:HG21	1:C:174:ILE:HD11	2.03	0.41
1:I:221:MET:HE3	1:I:221:MET:HB2	1.85	0.41
1:B:230:MET:HG3	1:B:241:ARG:CZ	2.51	0.41
1:H:59:ARG:O	1:H:110:GLN:HB2	2.22	0.40
1:H:82:LEU:HD21	1:M:102:MET:HG3	2.03	0.40
1:M:160:LYS:HD3	1:M:160:LYS:HA	1.94	0.40
1:P:123:ASP:OD1	1:P:161:GLN:NE2	2.46	0.40
1:R:116:VAL:HG21	1:R:174:ILE:HD11	2.03	0.40
1:S:116:VAL:HG21	1:S:174:ILE:HD11	2.02	0.40
1:E:163:LEU:HB2	1:E:197:LEU:CD1	2.44	0.40
1:E:230:MET:HG3	1:E:241:ARG:CZ	2.51	0.40
1:P:116:VAL:HG21	1:P:174:ILE:HD11	2.03	0.40
1:P:210:HIS:CD2	1:P:225:THR:HG21	2.56	0.40
1:Q:163:LEU:HB2	1:Q:197:LEU:HD11	2.03	0.40
1:Q:221:MET:HE3	1:Q:221:MET:HB2	1.93	0.40
1:Q:160:LYS:HA	1:Q:160:LYS:HD3	1.92	0.40
1:T:117:LEU:HA	1:T:167:LYS:HG3	2.03	0.40
1:T:146:VAL:CG1	1:T:211:MET:HB2	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	190/224 (85%)	185 (97%)	5 (3%)	0	100	100
1	B	188/224 (84%)	186 (99%)	2 (1%)	0	100	100
1	C	189/224 (84%)	187 (99%)	2 (1%)	0	100	100
1	D	189/224 (84%)	187 (99%)	2 (1%)	0	100	100
1	E	188/224 (84%)	186 (99%)	2 (1%)	0	100	100
1	F	188/224 (84%)	186 (99%)	2 (1%)	0	100	100
1	G	191/224 (85%)	189 (99%)	2 (1%)	0	100	100
1	H	191/224 (85%)	189 (99%)	2 (1%)	0	100	100
1	I	190/224 (85%)	188 (99%)	2 (1%)	0	100	100
1	J	188/224 (84%)	186 (99%)	2 (1%)	0	100	100
1	K	190/224 (85%)	187 (98%)	3 (2%)	0	100	100
1	L	189/224 (84%)	187 (99%)	2 (1%)	0	100	100
1	M	189/224 (84%)	187 (99%)	2 (1%)	0	100	100
1	N	191/224 (85%)	189 (99%)	2 (1%)	0	100	100
1	O	189/224 (84%)	187 (99%)	2 (1%)	0	100	100
1	P	187/224 (84%)	185 (99%)	2 (1%)	0	100	100
1	Q	188/224 (84%)	186 (99%)	2 (1%)	0	100	100
1	R	189/224 (84%)	187 (99%)	2 (1%)	0	100	100
1	S	184/224 (82%)	184 (100%)	0	0	100	100
1	T	184/224 (82%)	181 (98%)	3 (2%)	0	100	100
All	All	3772/4480 (84%)	3729 (99%)	43 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	169/195 (87%)	162 (96%)	7 (4%)	27	37
1	B	167/195 (86%)	160 (96%)	7 (4%)	26	36
1	C	168/195 (86%)	164 (98%)	4 (2%)	43	58
1	D	168/195 (86%)	161 (96%)	7 (4%)	26	36
1	E	167/195 (86%)	161 (96%)	6 (4%)	31	42
1	F	167/195 (86%)	162 (97%)	5 (3%)	36	49
1	G	170/195 (87%)	163 (96%)	7 (4%)	27	37
1	H	170/195 (87%)	165 (97%)	5 (3%)	37	51
1	I	169/195 (87%)	163 (96%)	6 (4%)	31	42
1	J	167/195 (86%)	161 (96%)	6 (4%)	31	42
1	K	169/195 (87%)	165 (98%)	4 (2%)	43	58
1	L	168/195 (86%)	161 (96%)	7 (4%)	26	36
1	M	168/195 (86%)	161 (96%)	7 (4%)	26	36
1	N	170/195 (87%)	165 (97%)	5 (3%)	37	51
1	O	168/195 (86%)	160 (95%)	8 (5%)	23	30
1	P	166/195 (85%)	155 (93%)	11 (7%)	15	18
1	Q	167/195 (86%)	158 (95%)	9 (5%)	20	25
1	R	168/195 (86%)	160 (95%)	8 (5%)	23	30
1	S	165/195 (85%)	160 (97%)	5 (3%)	36	49
1	T	163/195 (84%)	157 (96%)	6 (4%)	30	41
All	All	3354/3900 (86%)	3224 (96%)	130 (4%)	28	39

All (130) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	92	LEU
1	A	97	ARG
1	A	181	VAL

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Mol	Chain	Res	Type
1	A	211	MET
1	A	214	VAL
1	A	216	ARG
1	A	242	GLU
1	B	66	LEU
1	B	97	ARG
1	B	110	GLN
1	B	181	VAL
1	B	187	LYS
1	B	211	MET
1	B	214	VAL
1	C	97	ARG
1	C	181	VAL
1	C	211	MET
1	C	214	VAL
1	D	97	ARG
1	D	181	VAL
1	D	187	LYS
1	D	197	LEU
1	D	211	MET
1	D	214	VAL
1	D	242	GLU
1	E	66	LEU
1	E	84	GLU
1	E	97	ARG
1	E	181	VAL
1	E	211	MET
1	E	214	VAL
1	F	59	ARG
1	F	181	VAL
1	F	211	MET
1	F	214	VAL
1	F	221	MET
1	G	93	LYS
1	G	97	ARG
1	G	125	ASP
1	G	160	LYS
1	G	181	VAL
1	G	214	VAL
1	G	242	GLU
1	H	93	LYS
1	H	181	VAL

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Mol	Chain	Res	Type
1	H	211	MET
1	H	214	VAL
1	H	242	GLU
1	I	97	ARG
1	I	122	PHE
1	I	181	VAL
1	I	211	MET
1	I	214	VAL
1	I	239	LYS
1	J	97	ARG
1	J	157	LEU
1	J	162	VAL
1	J	181	VAL
1	J	211	MET
1	J	214	VAL
1	K	97	ARG
1	K	181	VAL
1	K	211	MET
1	K	214	VAL
1	L	93	LYS
1	L	97	ARG
1	L	112	THR
1	L	117	LEU
1	L	181	VAL
1	L	211	MET
1	L	214	VAL
1	M	62	GLU
1	M	97	ARG
1	M	157	LEU
1	M	181	VAL
1	M	211	MET
1	M	214	VAL
1	M	242	GLU
1	N	59	ARG
1	N	181	VAL
1	N	211	MET
1	N	214	VAL
1	N	224	LYS
1	O	61	GLU
1	O	91	LEU
1	O	97	ARG
1	O	181	VAL

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Mol	Chain	Res	Type
1	O	194	THR
1	O	211	MET
1	O	214	VAL
1	O	242	GLU
1	P	97	ARG
1	P	107	LYS
1	P	159	ASN
1	P	181	VAL
1	P	187	LYS
1	P	198	ARG
1	P	211	MET
1	P	214	VAL
1	P	221	MET
1	P	225	THR
1	P	242	GLU
1	Q	97	ARG
1	Q	114	SER
1	Q	116	VAL
1	Q	157	LEU
1	Q	181	VAL
1	Q	197	LEU
1	Q	211	MET
1	Q	214	VAL
1	Q	242	GLU
1	R	66	LEU
1	R	97	ARG
1	R	117	LEU
1	R	181	VAL
1	R	211	MET
1	R	214	VAL
1	R	221	MET
1	R	242	GLU
1	S	62	GLU
1	S	66	LEU
1	S	181	VAL
1	S	211	MET
1	S	214	VAL
1	T	93	LYS
1	T	181	VAL
1	T	211	MET
1	T	214	VAL
1	T	221	MET

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Mol	Chain	Res	Type
1	T	242	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	126	HIS
1	A	222	ASN
1	C	126	HIS
1	C	161	GLN
1	G	222	ASN
1	H	222	ASN
1	J	222	ASN
1	K	222	ASN
1	L	126	HIS
1	M	70	ASN
1	M	222	ASN
1	N	87	GLN
1	N	222	ASN
1	T	85	ASN
1	T	222	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 40 ligands modelled in this entry, 20 are monoatomic - leaving 20 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	QBQ	O	302	-	33,34,34	0.58	0	47,54,54	0.89	1 (2%)
3	QBQ	B	302	-	33,34,34	0.65	1 (3%)	47,54,54	0.94	1 (2%)
3	QBQ	I	302	-	33,34,34	0.57	0	47,54,54	0.94	2 (4%)
3	QBQ	D	302	-	33,34,34	0.47	0	47,54,54	0.83	0
3	QBQ	R	302	-	33,34,34	0.60	0	47,54,54	0.89	1 (2%)
3	QBQ	M	302	-	33,34,34	0.50	0	47,54,54	0.90	1 (2%)
3	QBQ	Q	302	-	33,34,34	0.52	0	47,54,54	0.81	0
3	QBQ	K	302	-	33,34,34	0.49	0	47,54,54	0.93	1 (2%)
3	QBQ	E	302	-	33,34,34	0.63	1 (3%)	47,54,54	0.88	1 (2%)
3	QBQ	N	302	-	33,34,34	0.58	1 (3%)	47,54,54	0.92	1 (2%)
3	QBQ	F	302	-	33,34,34	0.51	0	47,54,54	0.84	1 (2%)
3	QBQ	J	302	-	33,34,34	0.54	0	47,54,54	0.85	1 (2%)
3	QBQ	P	302	-	33,34,34	0.60	0	47,54,54	0.87	0
3	QBQ	H	302	-	33,34,34	0.57	0	47,54,54	0.84	1 (2%)
3	QBQ	S	302	-	33,34,34	0.60	0	47,54,54	0.84	0
3	QBQ	T	302	-	33,34,34	0.56	0	47,54,54	0.87	1 (2%)
3	QBQ	C	302	-	33,34,34	0.61	0	47,54,54	0.83	0
3	QBQ	G	302	-	33,34,34	0.57	0	47,54,54	0.91	1 (2%)
3	QBQ	A	302	-	33,34,34	0.49	0	47,54,54	0.90	1 (2%)
3	QBQ	L	302	-	33,34,34	0.51	0	47,54,54	0.87	1 (2%)

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
3	QBQ	O	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	B	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	I	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	D	302	-	-	7/22/38/38	0/3/3/3
3	QBQ	R	302	-	-	4/22/38/38	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	QBQ	M	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	Q	302	-	-	3/22/38/38	0/3/3/3
3	QBQ	K	302	-	-	2/22/38/38	0/3/3/3
3	QBQ	E	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	N	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	F	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	J	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	P	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	H	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	S	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	T	302	-	-	3/22/38/38	0/3/3/3
3	QBQ	C	302	-	-	4/22/38/38	0/3/3/3
3	QBQ	G	302	-	-	3/22/38/38	0/3/3/3
3	QBQ	A	302	-	-	7/22/38/38	0/3/3/3
3	QBQ	L	302	-	-	3/22/38/38	0/3/3/3

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	302	QBQ	P2-O8	-2.13	1.46	1.54
3	N	302	QBQ	C2-C1	-2.09	1.39	1.45
3	E	302	QBQ	P2-O8	-2.00	1.47	1.54

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	R	302	QBQ	O1-C7-C6	-2.52	100.42	108.99
3	M	302	QBQ	C10-C2-C1	-2.36	116.89	118.35
3	O	302	QBQ	C10-C2-C1	-2.31	116.92	118.35
3	L	302	QBQ	C10-C2-C1	-2.29	116.93	118.35
3	N	302	QBQ	O1-C7-C6	-2.28	101.25	108.99
3	A	302	QBQ	C10-C2-C1	-2.23	116.97	118.35
3	F	302	QBQ	C10-C2-C1	-2.22	116.97	118.35
3	I	302	QBQ	C10-C2-C1	-2.17	117.01	118.35
3	I	302	QBQ	O12-C9-C8	2.16	118.74	111.82
3	J	302	QBQ	C6-O-C5	-2.15	104.71	109.47
3	B	302	QBQ	C10-C2-C1	-2.10	117.05	118.35
3	G	302	QBQ	O11-C8-C6	-2.08	105.10	111.08
3	E	302	QBQ	O11-C8-C6	-2.05	105.20	111.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	K	302	QBQ	C10-C2-C1	-2.02	117.10	118.35
3	T	302	QBQ	C6-O-C5	-2.01	105.03	109.47
3	H	302	QBQ	O1-C7-C6	-2.00	102.18	108.99

There are no chirality outliers.

All (80) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	302	QBQ	P1-O7-P2-O9
3	B	302	QBQ	P1-O7-P2-O9
3	C	302	QBQ	P1-O7-P2-O9
3	D	302	QBQ	P1-O7-P2-O8
3	D	302	QBQ	P1-O7-P2-O9
3	E	302	QBQ	P1-O7-P2-O9
3	H	302	QBQ	P1-O7-P2-O9
3	J	302	QBQ	P1-O7-P2-O9
3	L	302	QBQ	P1-O7-P2-O9
3	O	302	QBQ	P1-O7-P2-O9
3	P	302	QBQ	P1-O7-P2-O9
3	R	302	QBQ	P1-O7-P2-O9
3	S	302	QBQ	P1-O7-P2-O9
3	Q	302	QBQ	P1-O7-P2-O10
3	A	302	QBQ	P1-O4-P-O3
3	C	302	QBQ	P1-O4-P-O3
3	E	302	QBQ	P1-O4-P-O3
3	G	302	QBQ	P1-O4-P-O3
3	H	302	QBQ	P1-O4-P-O3
3	N	302	QBQ	P1-O4-P-O3
3	O	302	QBQ	P1-O4-P-O3
3	P	302	QBQ	P1-O4-P-O3
3	S	302	QBQ	P1-O4-P-O3
3	T	302	QBQ	P1-O4-P-O3
3	G	302	QBQ	P1-O7-P2-O10
3	I	302	QBQ	P1-O7-P2-O10
3	F	302	QBQ	P1-O7-P2-O9
3	I	302	QBQ	P1-O7-P2-O9
3	F	302	QBQ	P1-O4-P-O3
3	K	302	QBQ	P1-O4-P-O2
3	M	302	QBQ	P1-O4-P-O2
3	D	302	QBQ	O-C6-C7-O1
3	M	302	QBQ	O-C6-C7-O1
3	D	302	QBQ	C7-O1-P-O3

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Mol	Chain	Res	Type	Atoms
3	M	302	QBQ	C7-O1-P-O3
3	F	302	QBQ	P1-O7-P2-O10
3	H	302	QBQ	P1-O7-P2-O10
3	J	302	QBQ	P1-O7-P2-O10
3	T	302	QBQ	P1-O7-P2-O10
3	B	302	QBQ	P1-O4-P-O3
3	C	302	QBQ	P1-O4-P-O2
3	D	302	QBQ	P1-O4-P-O3
3	D	302	QBQ	P1-O4-P-O2
3	E	302	QBQ	P1-O4-P-O2
3	G	302	QBQ	P1-O4-P-O2
3	H	302	QBQ	P1-O4-P-O2
3	J	302	QBQ	P1-O4-P-O3
3	N	302	QBQ	P1-O4-P-O2
3	O	302	QBQ	P1-O4-P-O2
3	Q	302	QBQ	P1-O4-P-O2
3	R	302	QBQ	P1-O4-P-O2
3	T	302	QBQ	P1-O4-P-O2
3	O	302	QBQ	P1-O7-P2-O10
3	A	302	QBQ	P1-O4-P-O2
3	F	302	QBQ	P1-O4-P-O2
3	I	302	QBQ	P1-O4-P-O3
3	Q	302	QBQ	P1-O4-P-O3
3	R	302	QBQ	P1-O4-P-O3
3	L	302	QBQ	C9-C5-N2-C4
3	A	302	QBQ	O-C6-C7-O1
3	A	302	QBQ	P1-O7-P2-O10
3	B	302	QBQ	P1-O7-P2-O10
3	C	302	QBQ	P1-O7-P2-O10
3	D	302	QBQ	P1-O7-P2-O10
3	E	302	QBQ	P1-O7-P2-O10
3	L	302	QBQ	P1-O7-P2-O10
3	P	302	QBQ	P1-O7-P2-O10
3	A	302	QBQ	P1-O7-P2-O8
3	B	302	QBQ	P1-O4-P-O2
3	I	302	QBQ	P1-O4-P-O2
3	J	302	QBQ	P1-O4-P-O2
3	K	302	QBQ	P1-O4-P-O3
3	M	302	QBQ	P1-O4-P-O3
3	P	302	QBQ	P1-O4-P-O2
3	S	302	QBQ	P1-O4-P-O2
3	A	302	QBQ	C9-C5-N2-C4

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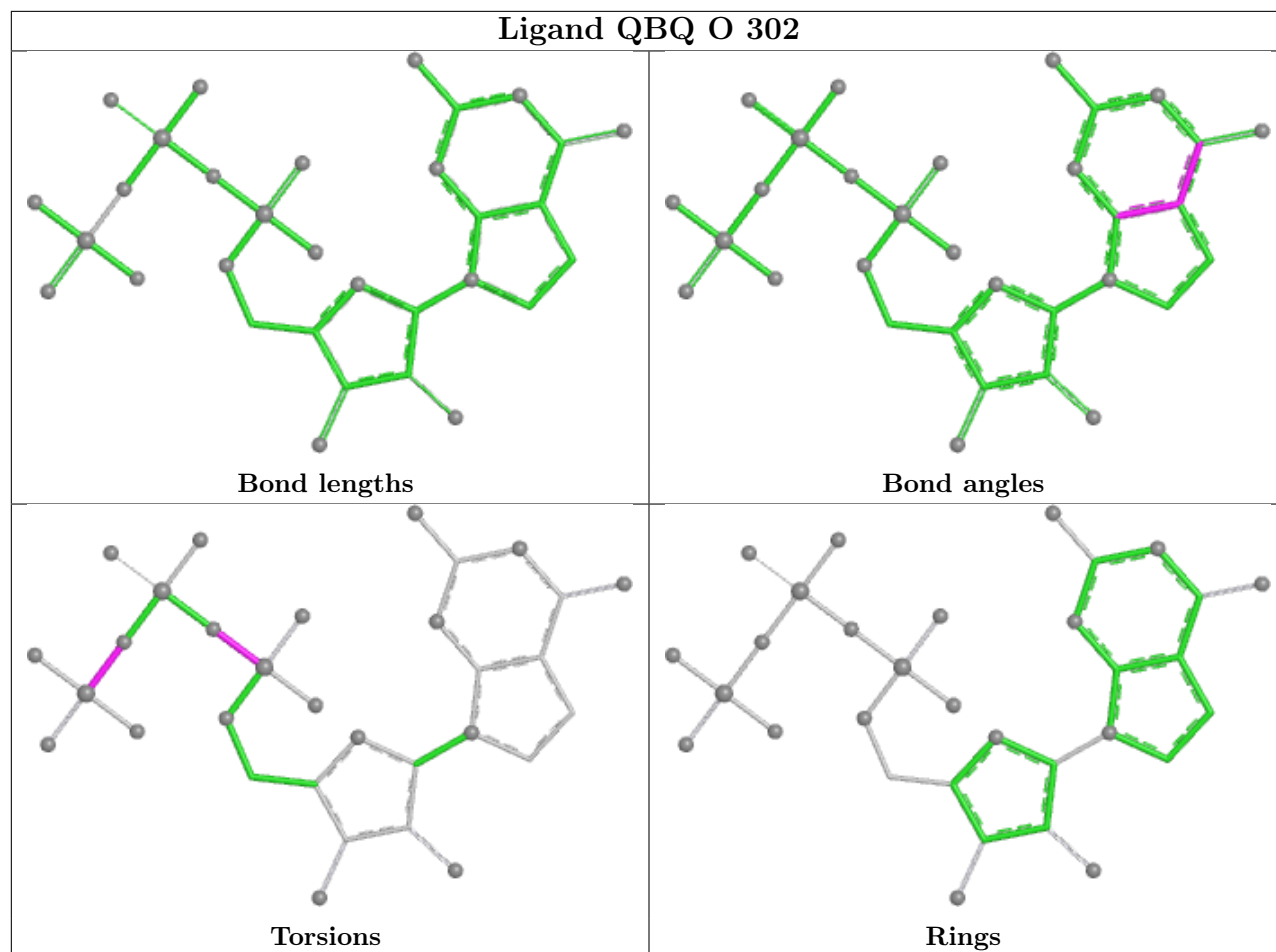
Mol	Chain	Res	Type	Atoms
3	N	302	QBQ	O-C6-C7-O1
3	N	302	QBQ	P1-O7-P2-O10
3	R	302	QBQ	P1-O7-P2-O10
3	S	302	QBQ	P1-O7-P2-O10

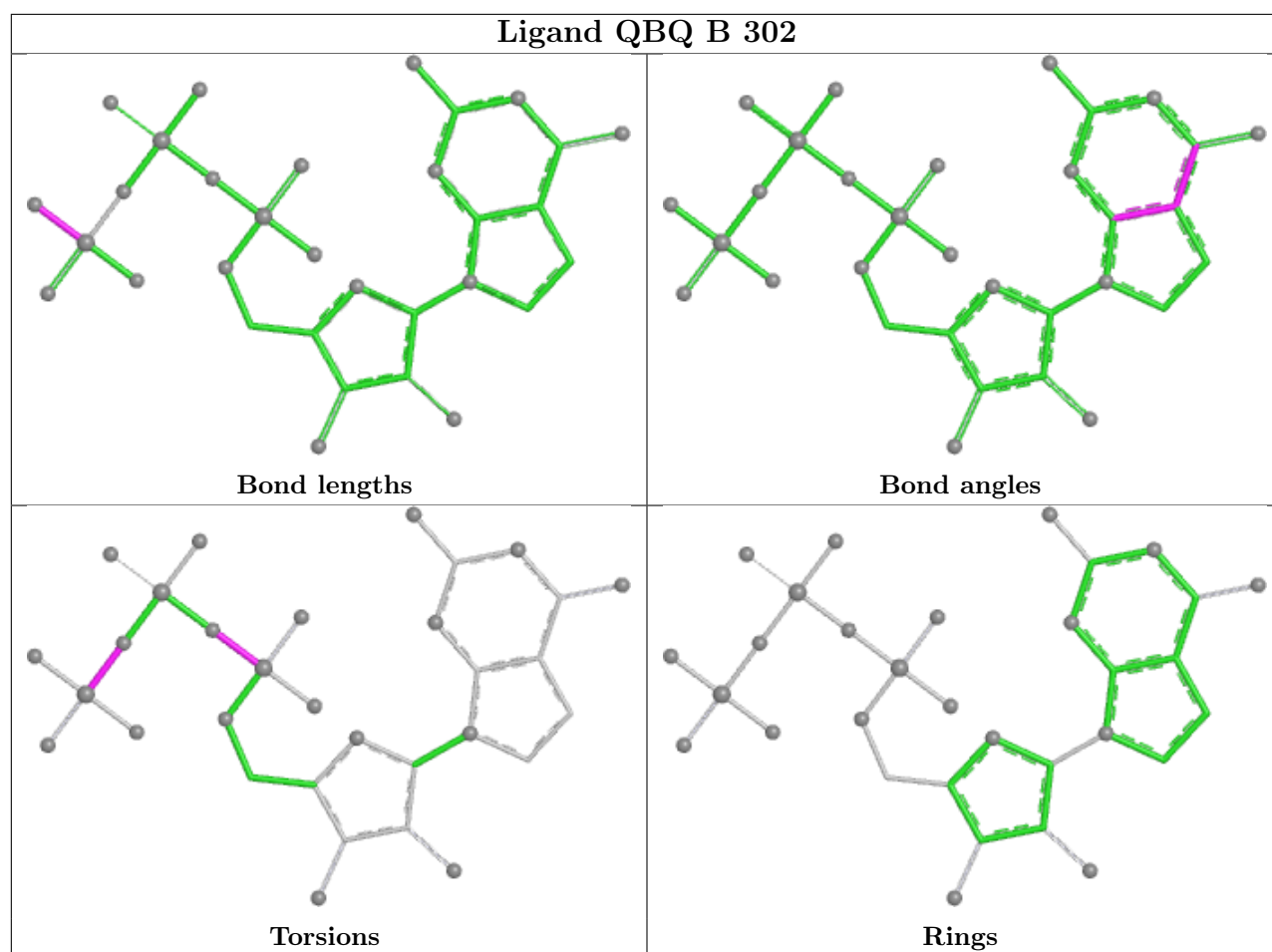
There are no ring outliers.

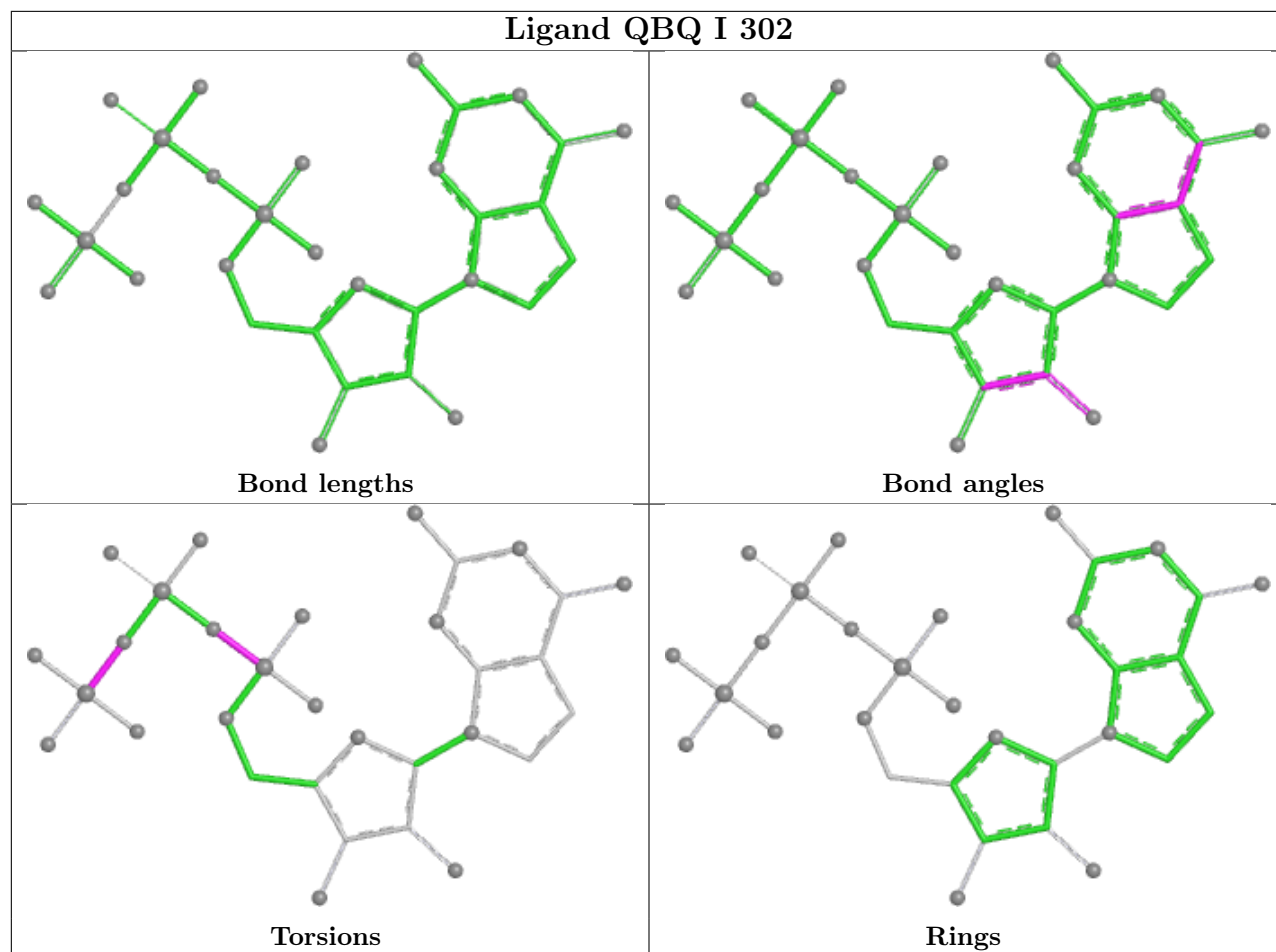
2 monomers are involved in 2 short contacts:

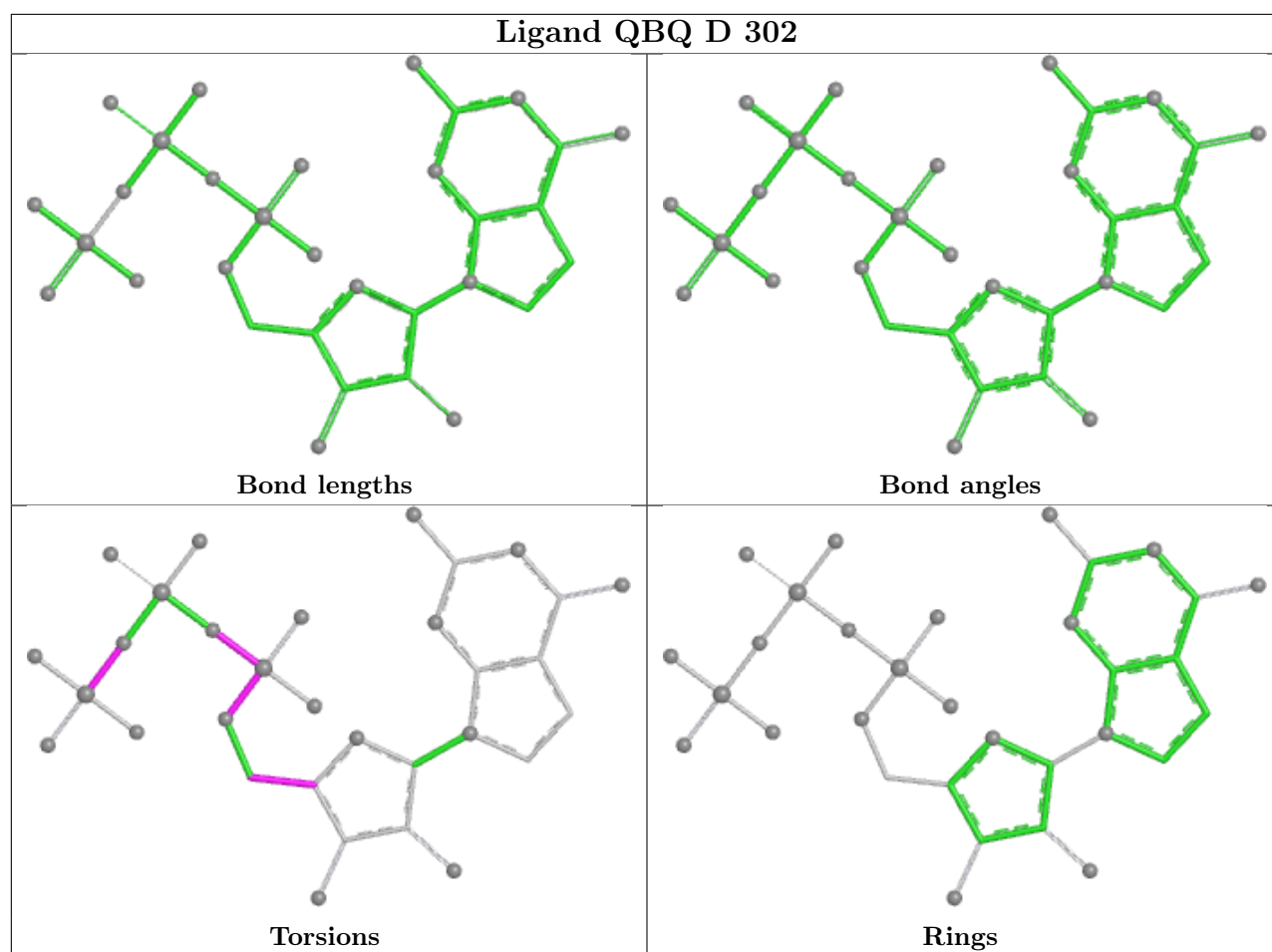
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	Q	302	QBQ	1	0
3	C	302	QBQ	1	0

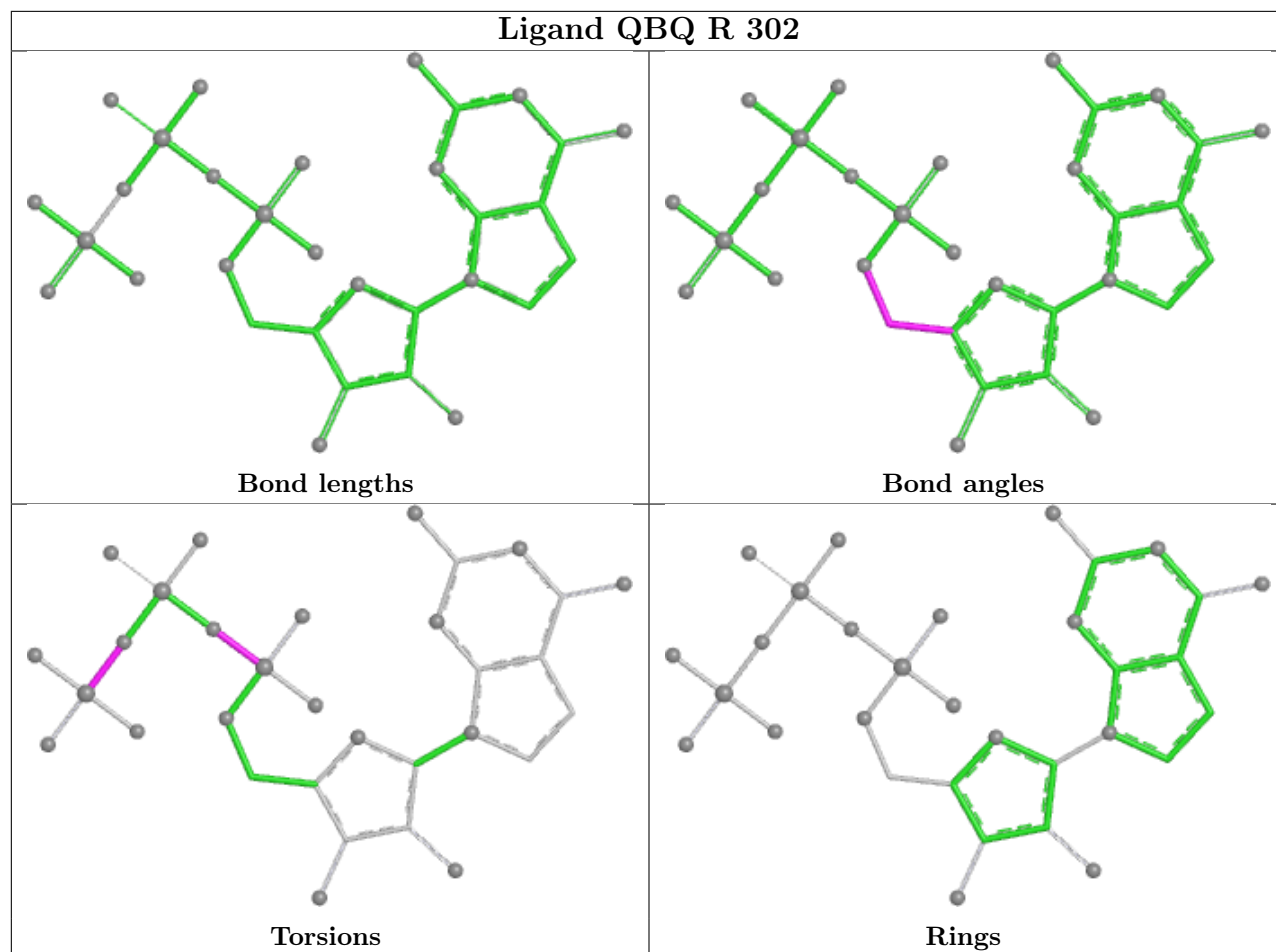
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

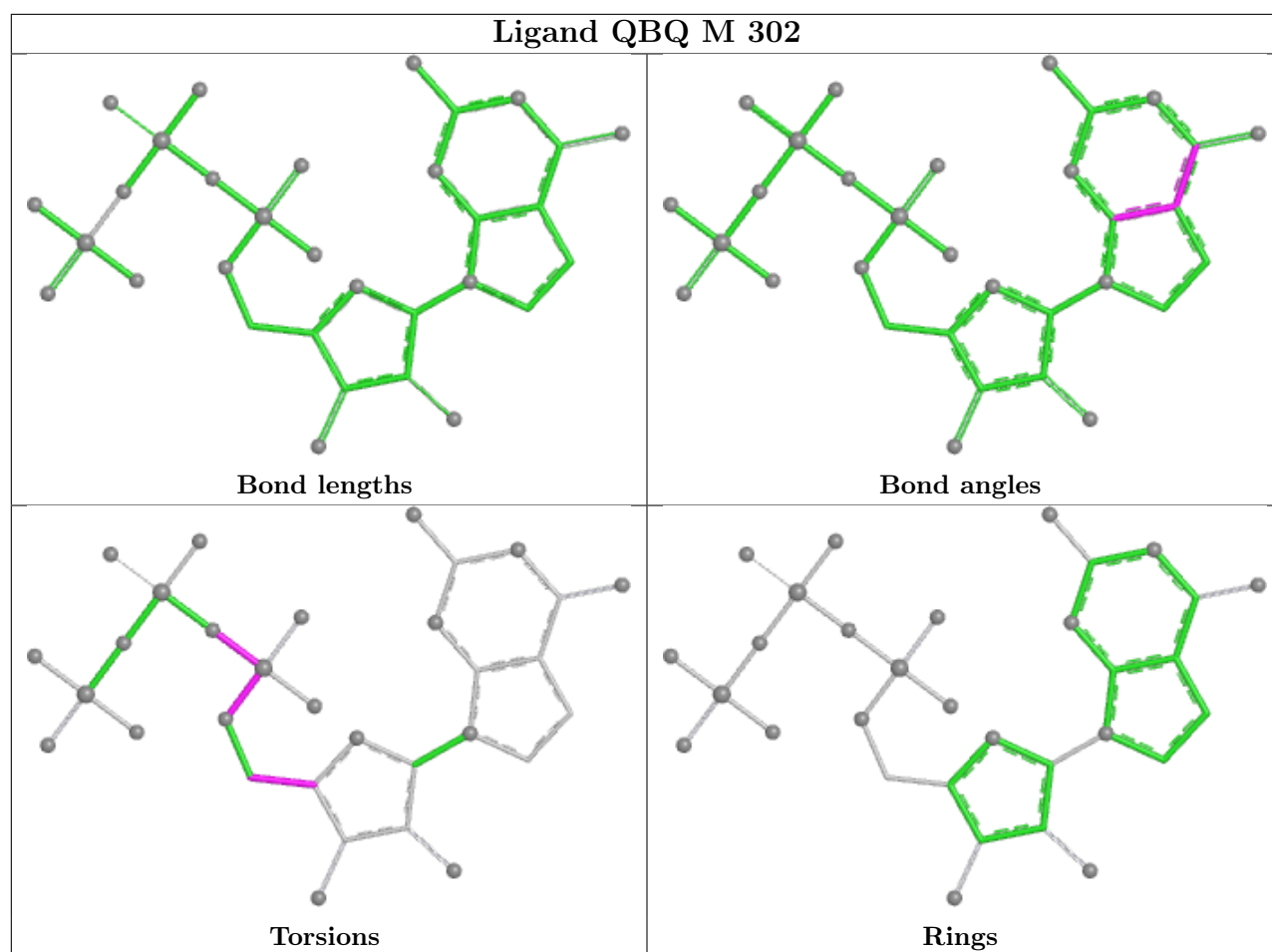


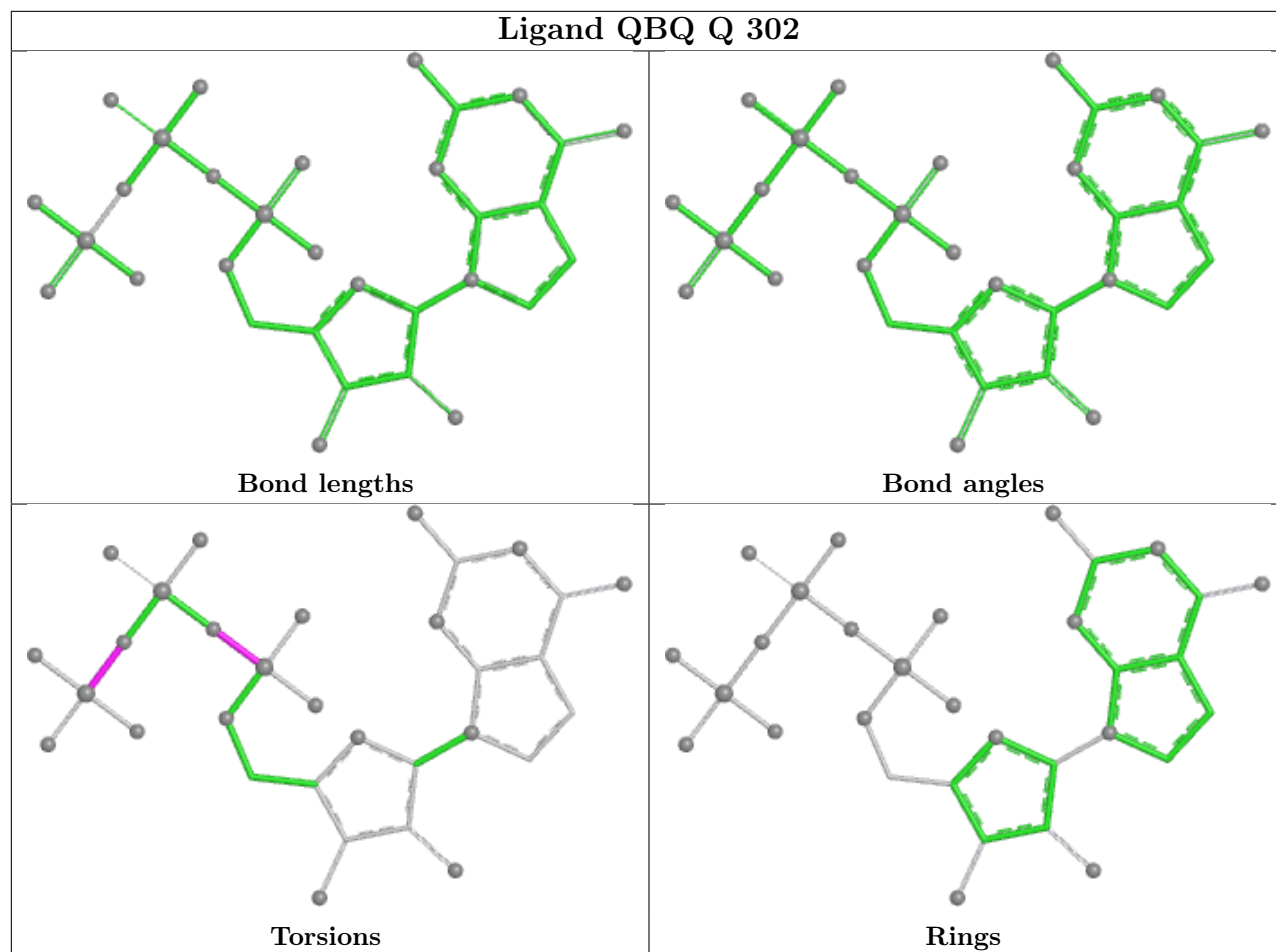


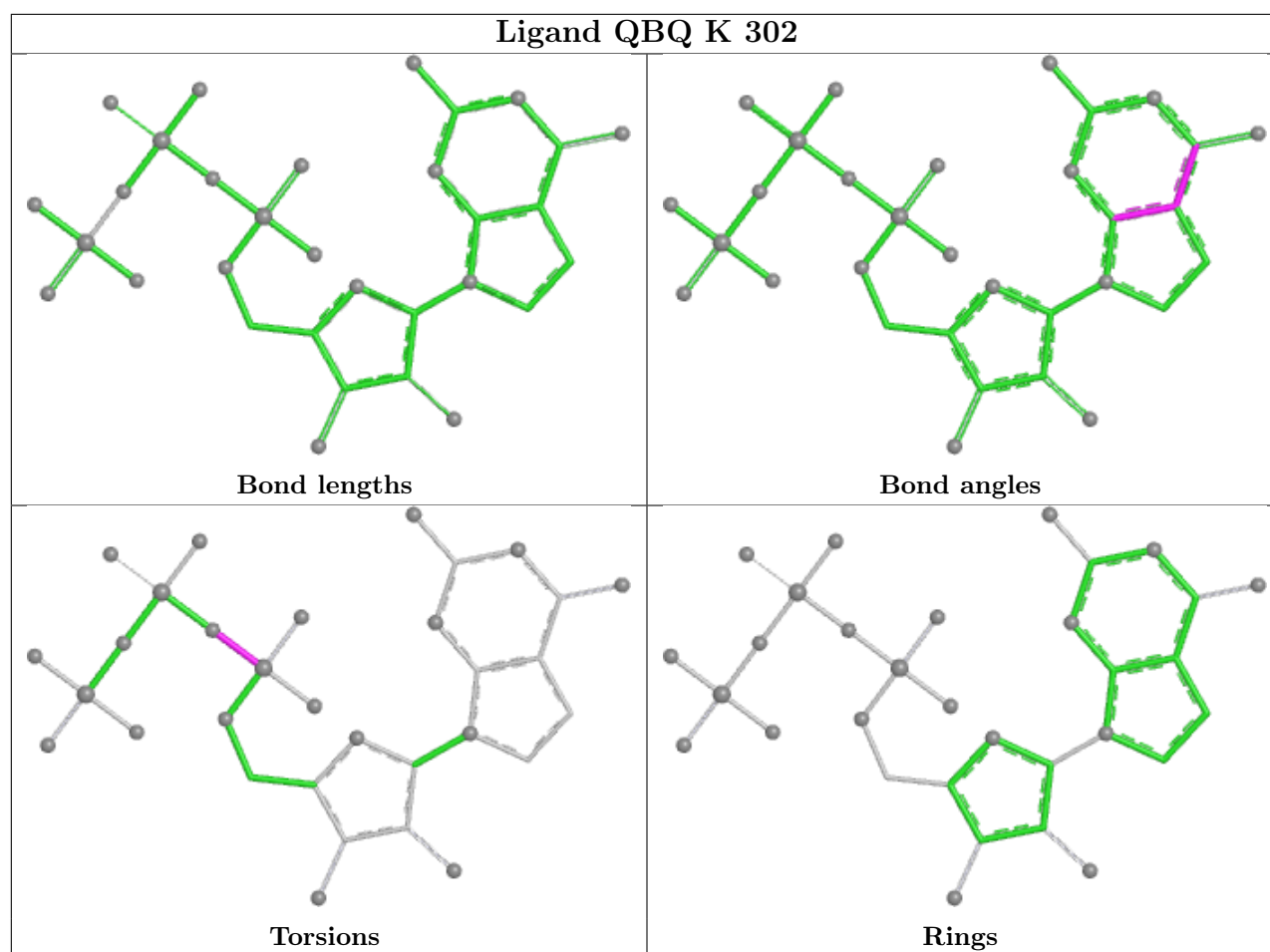


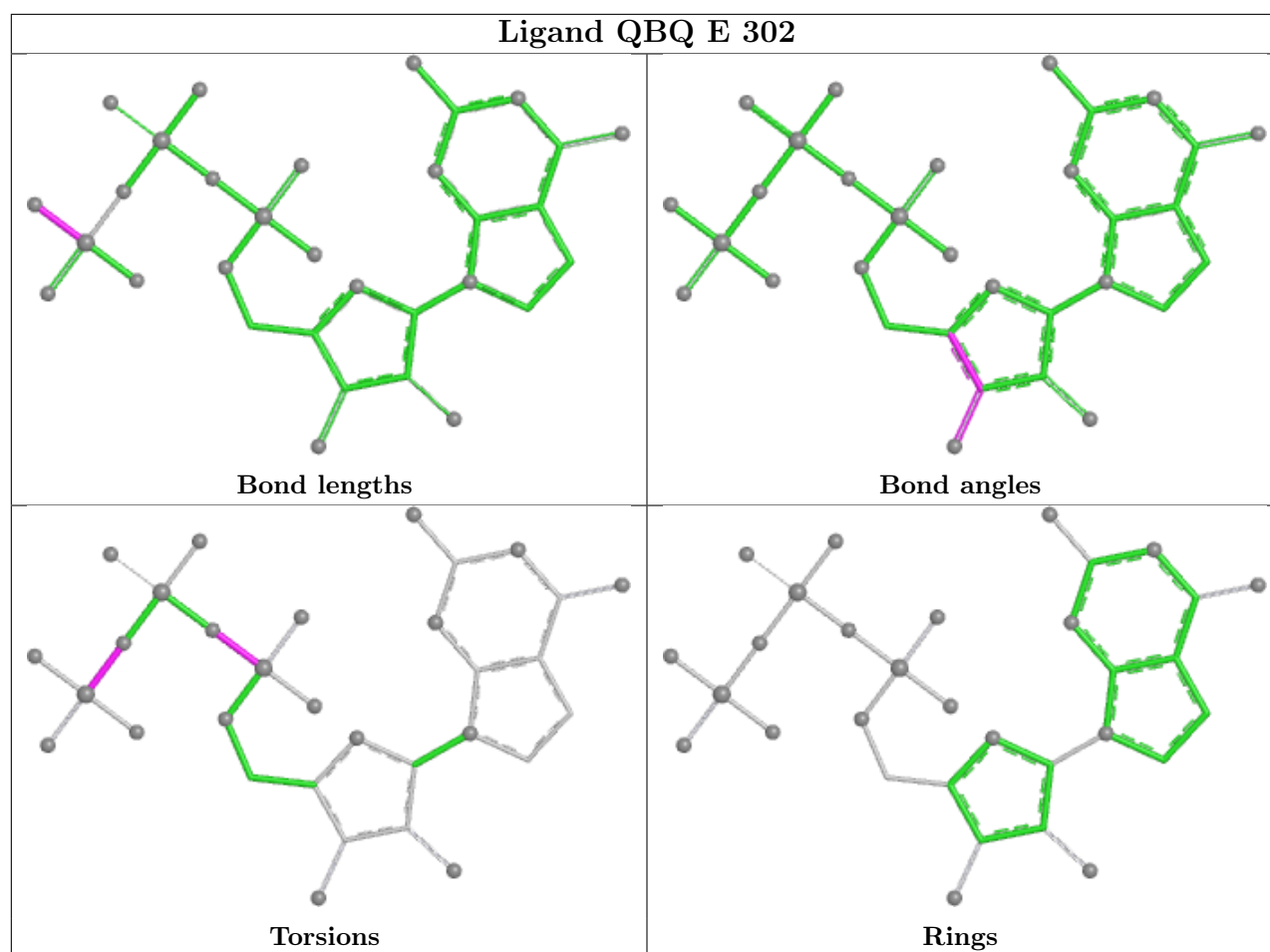


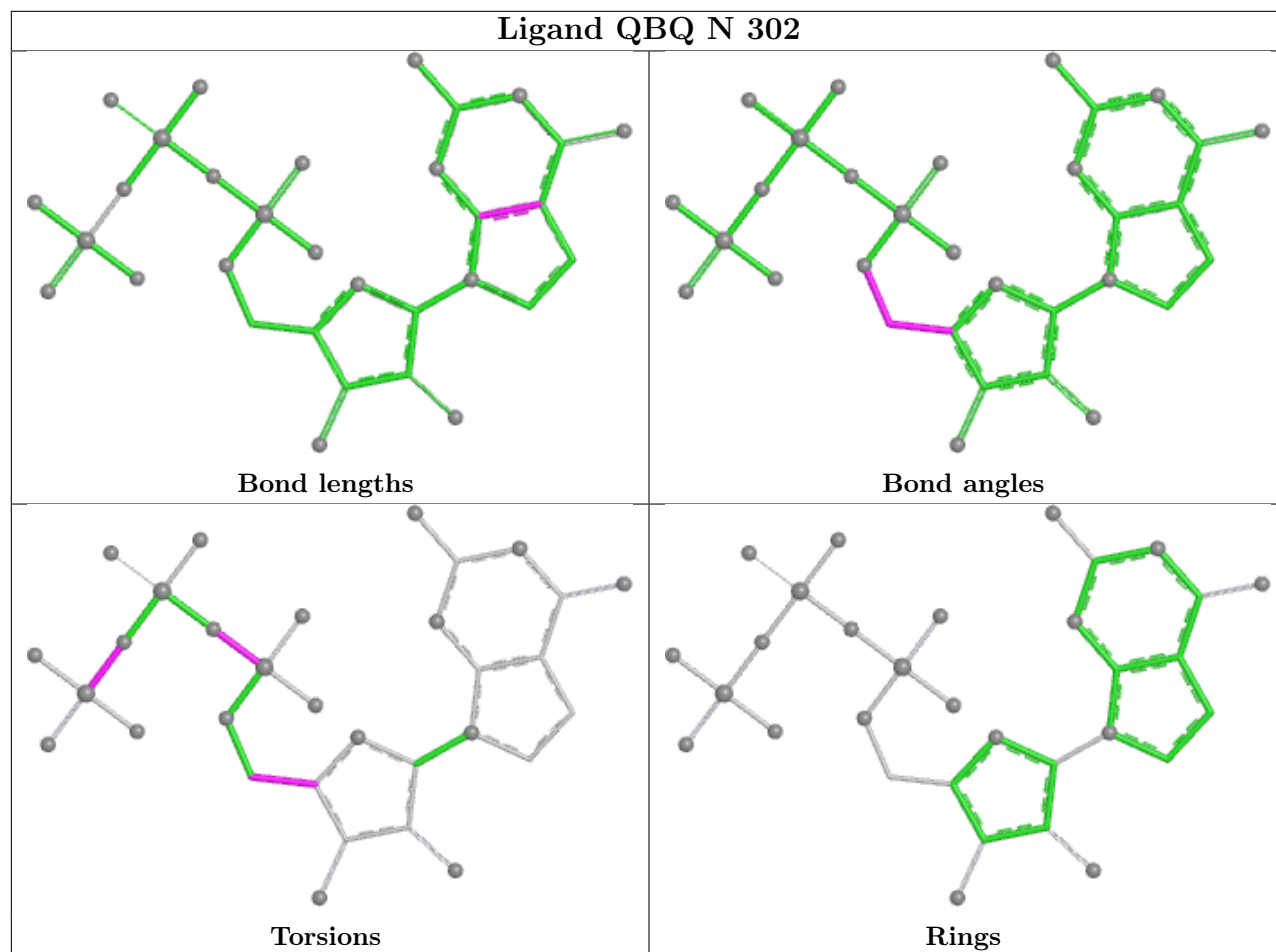


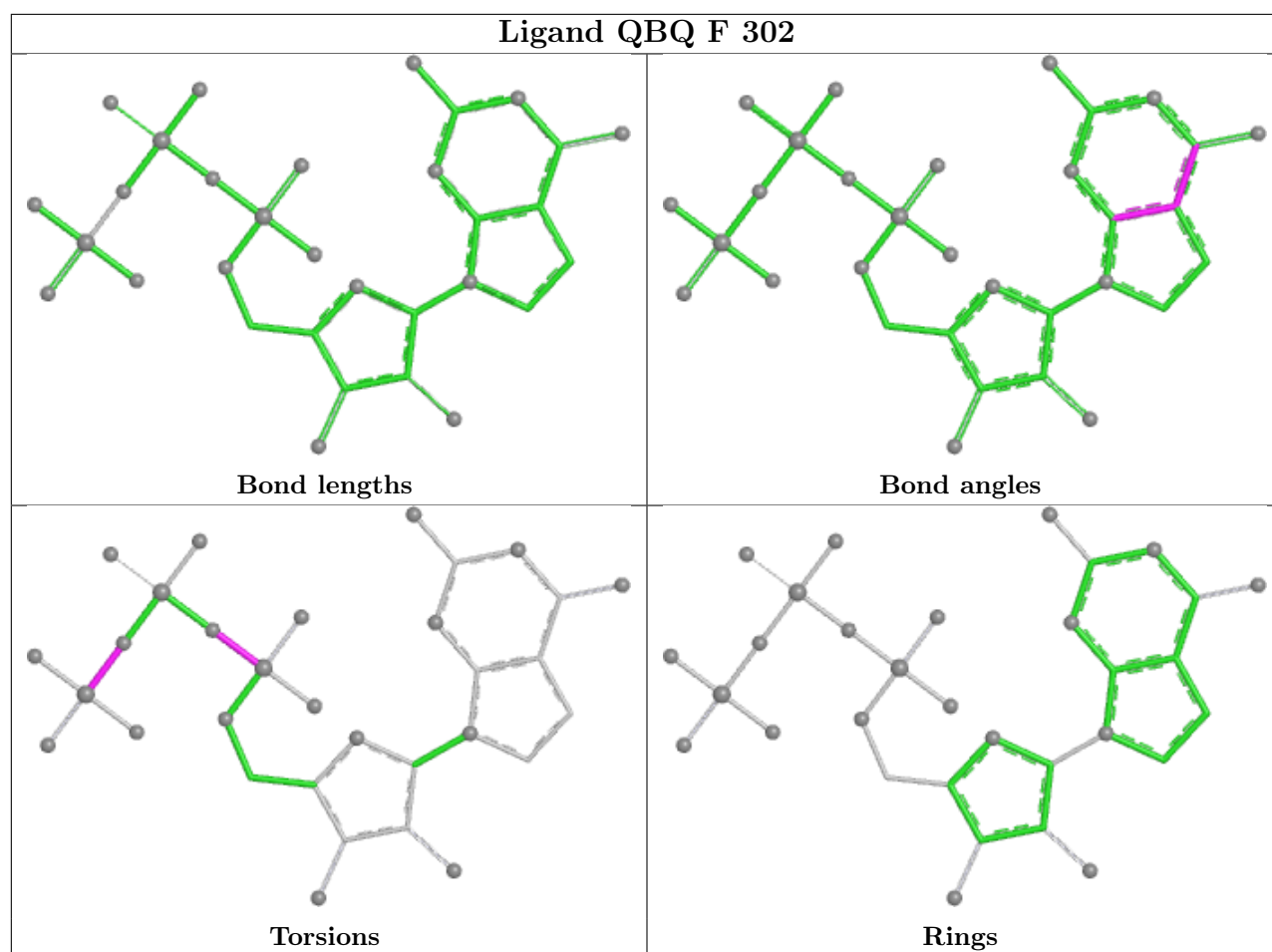


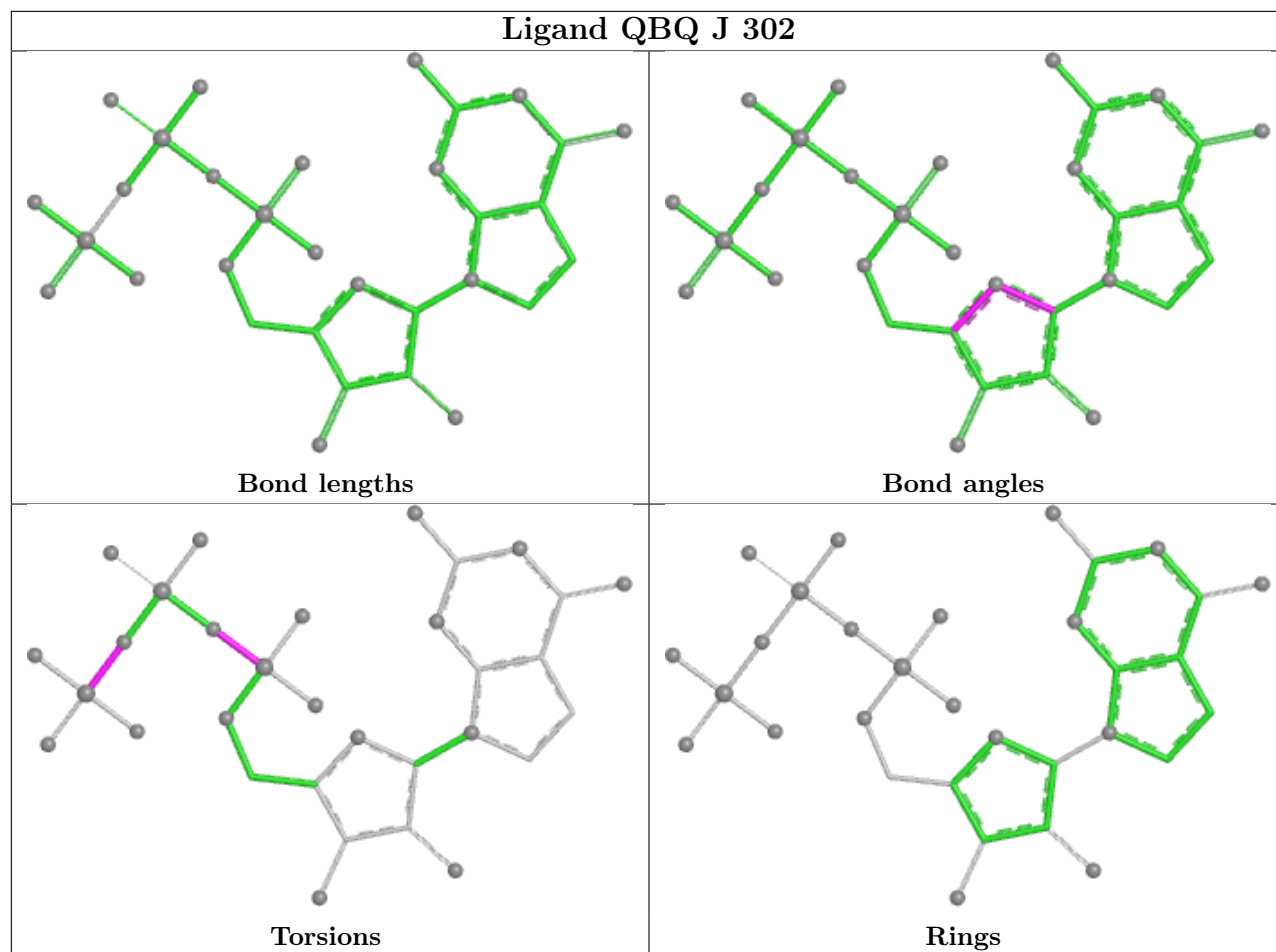


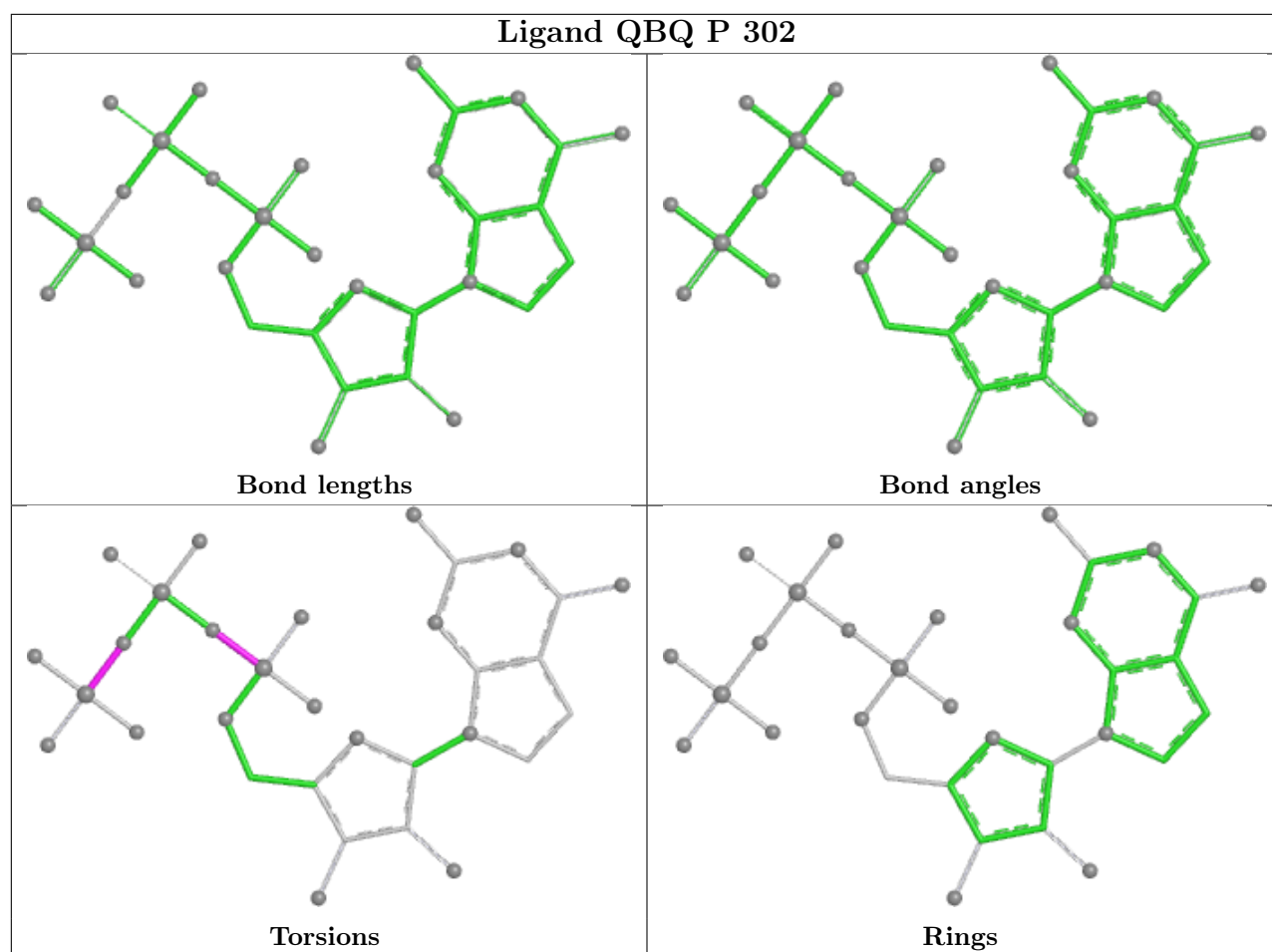


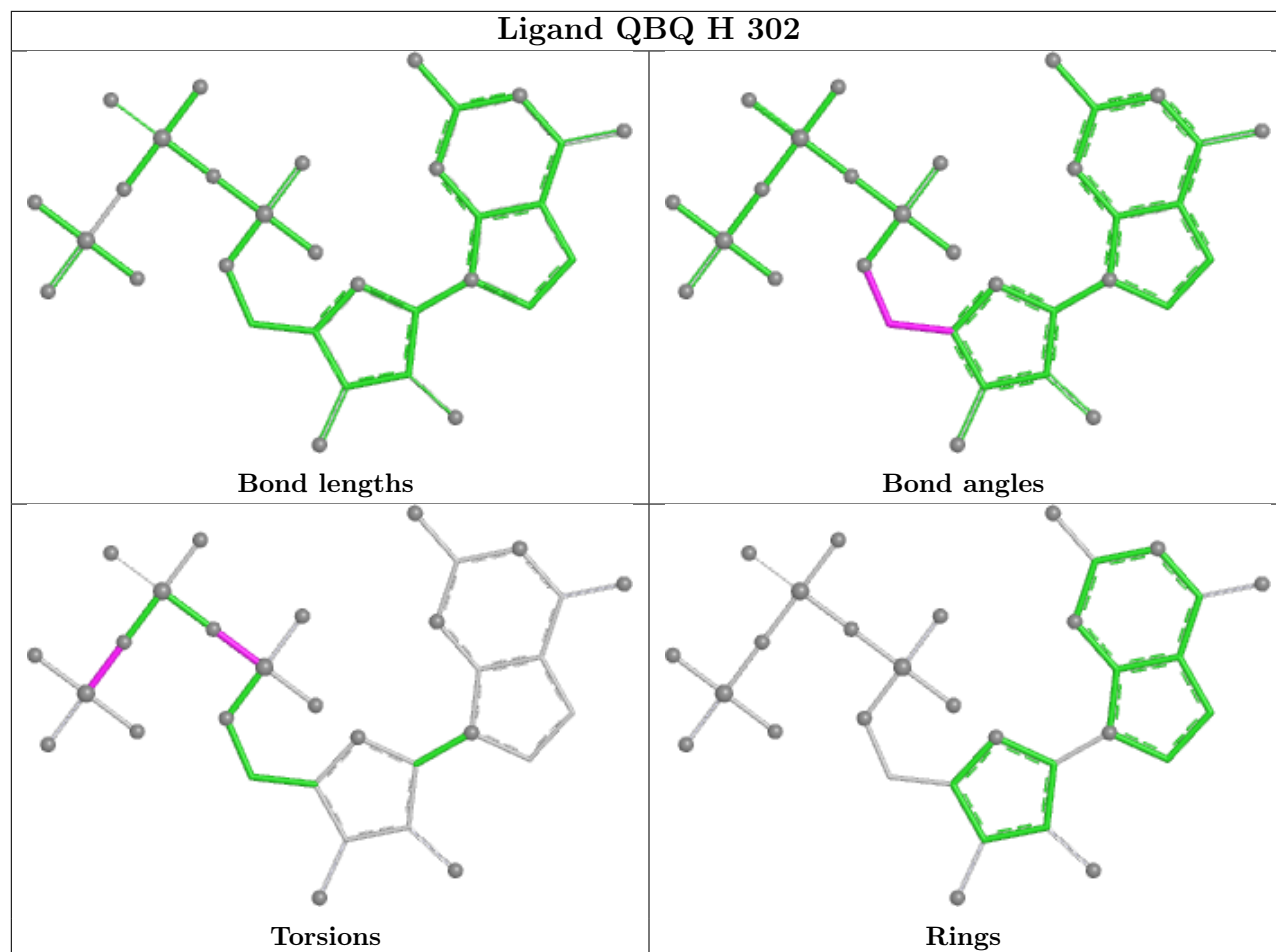


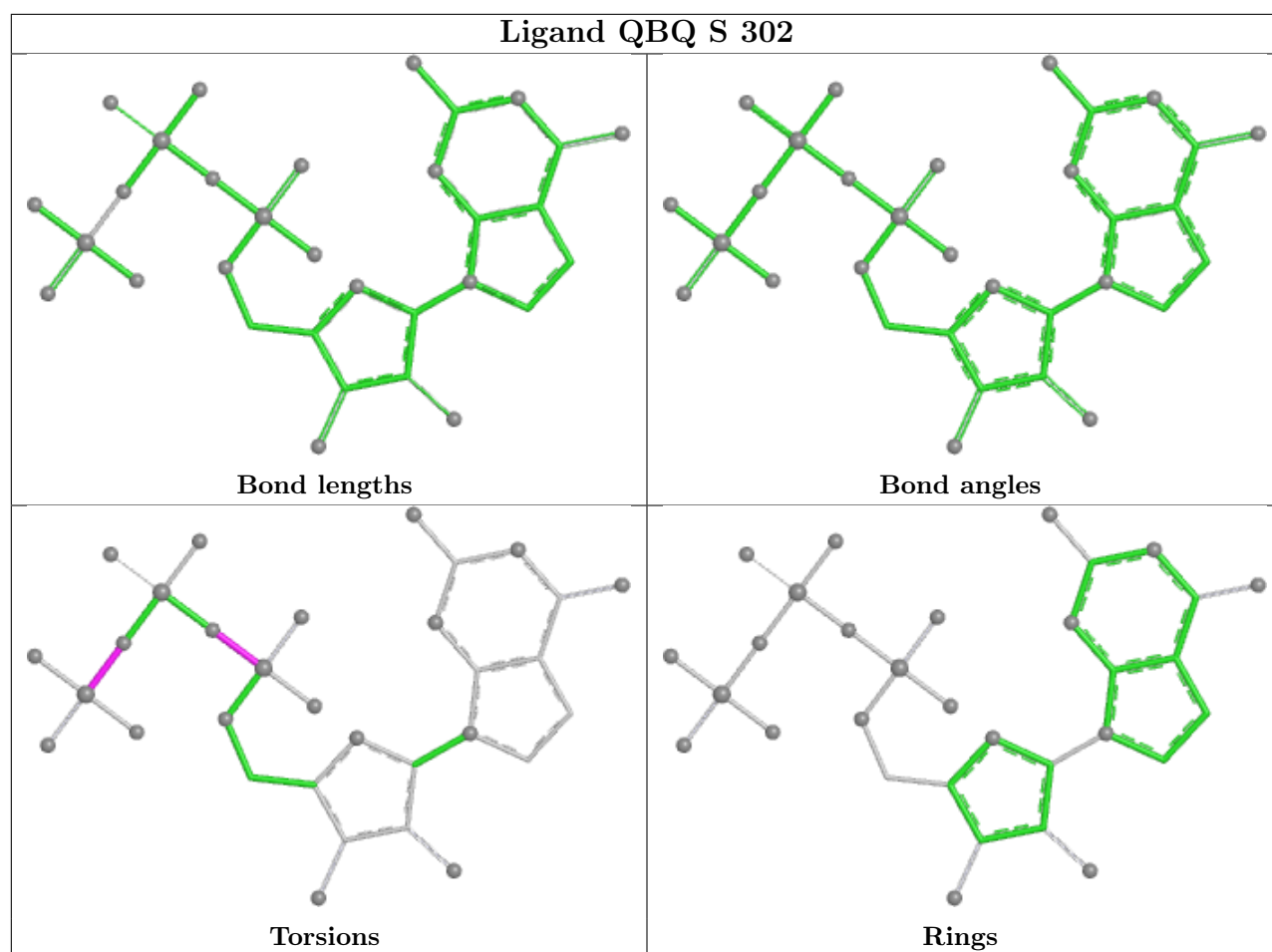


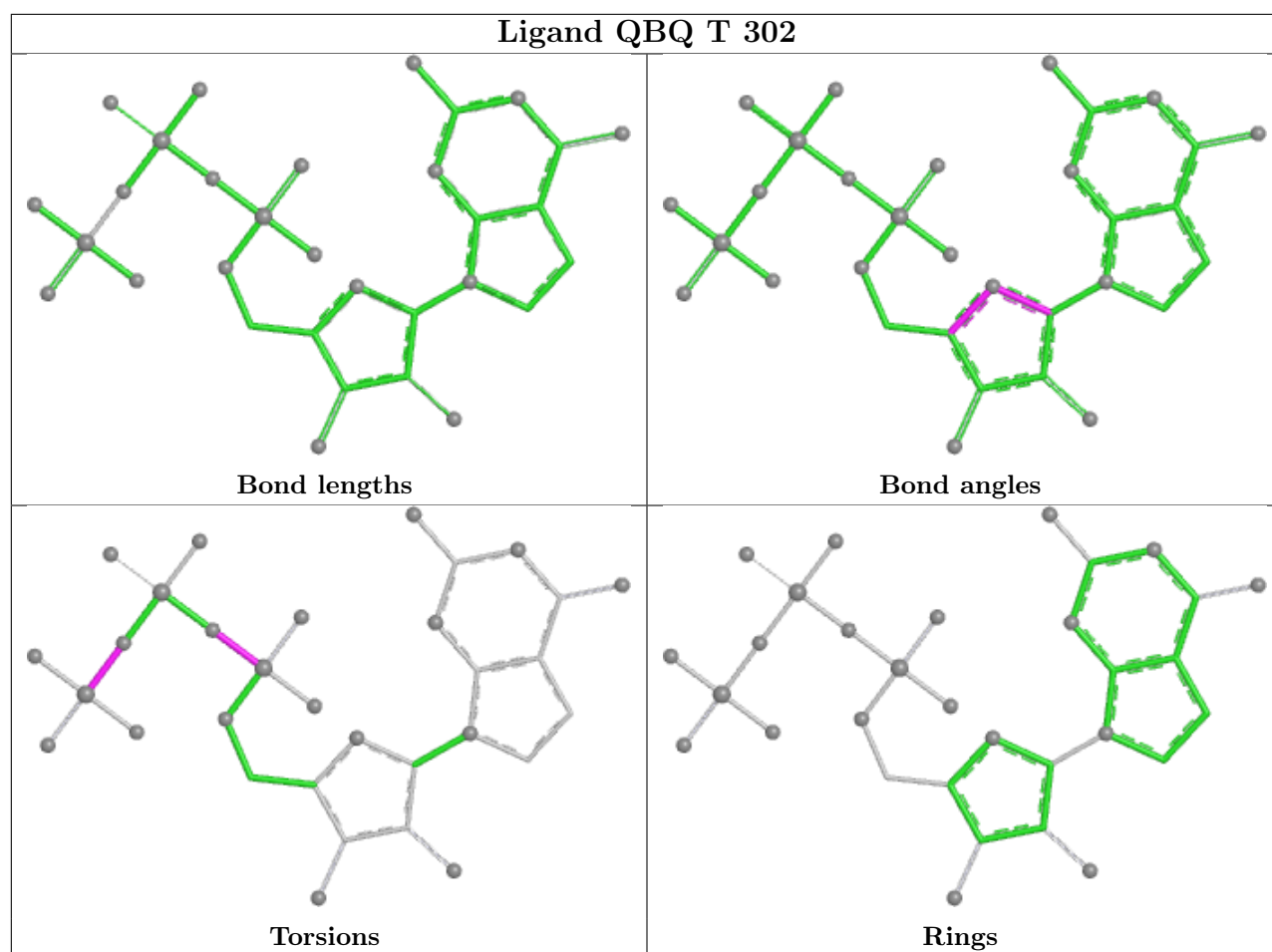


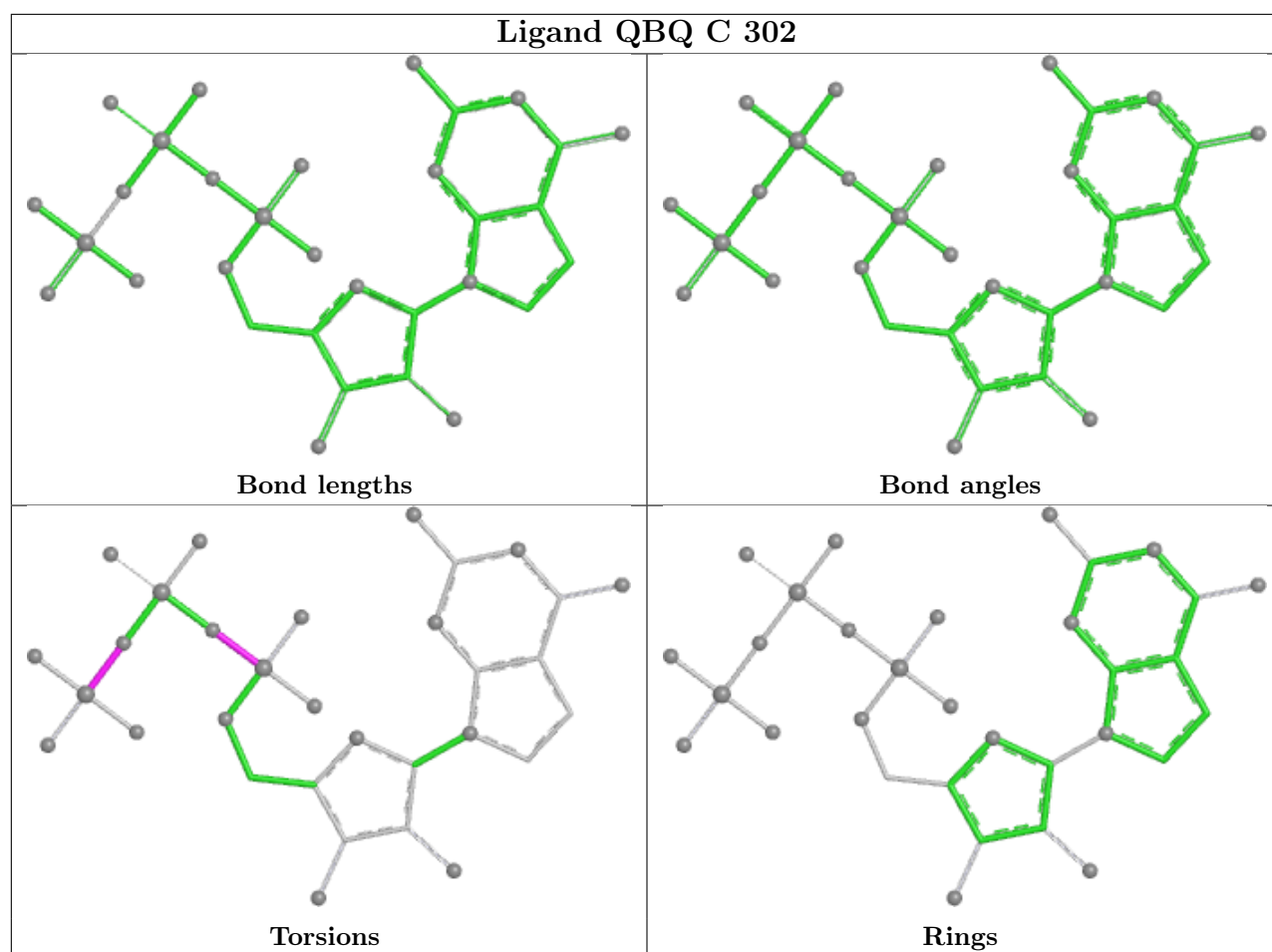


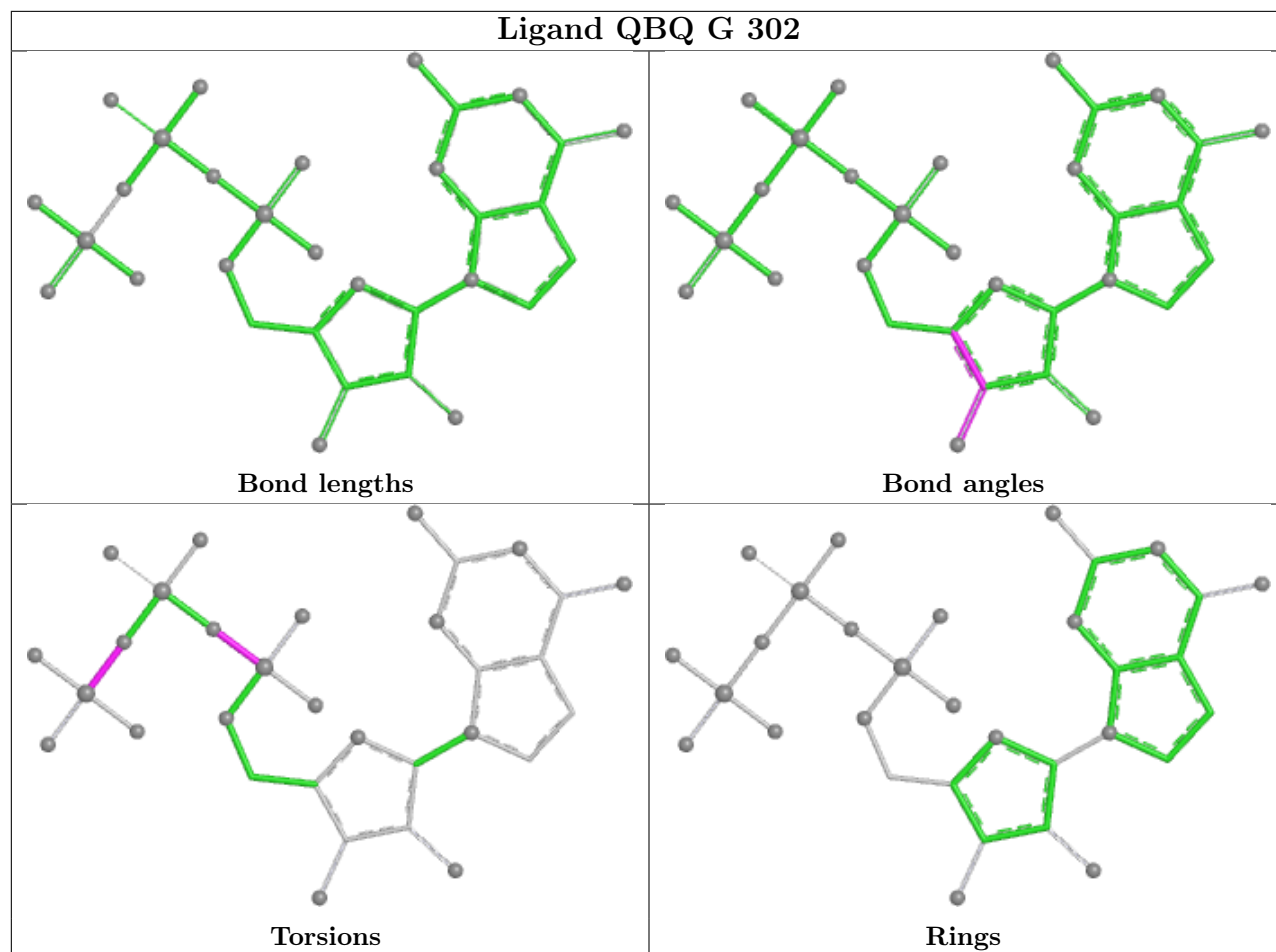


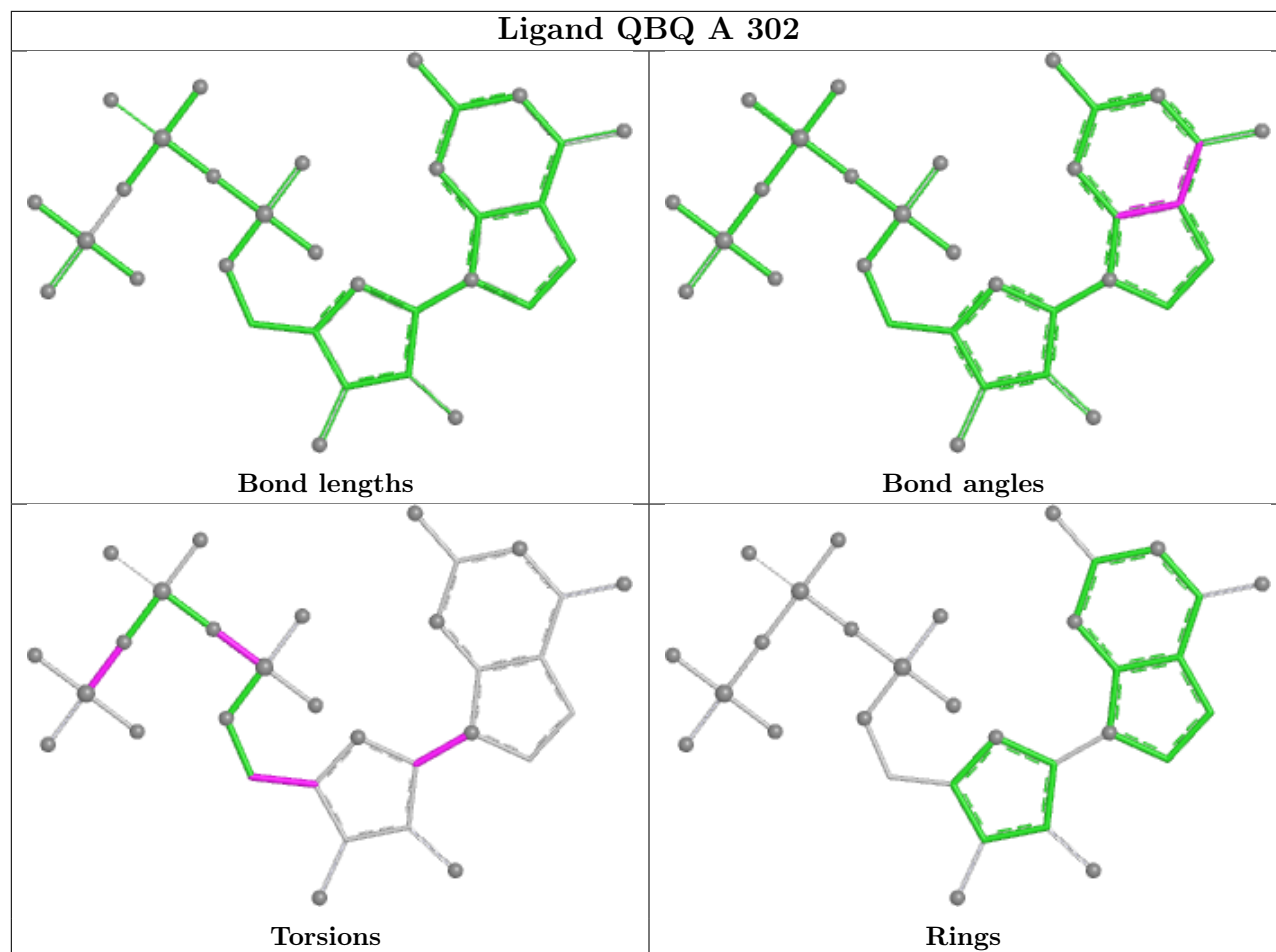


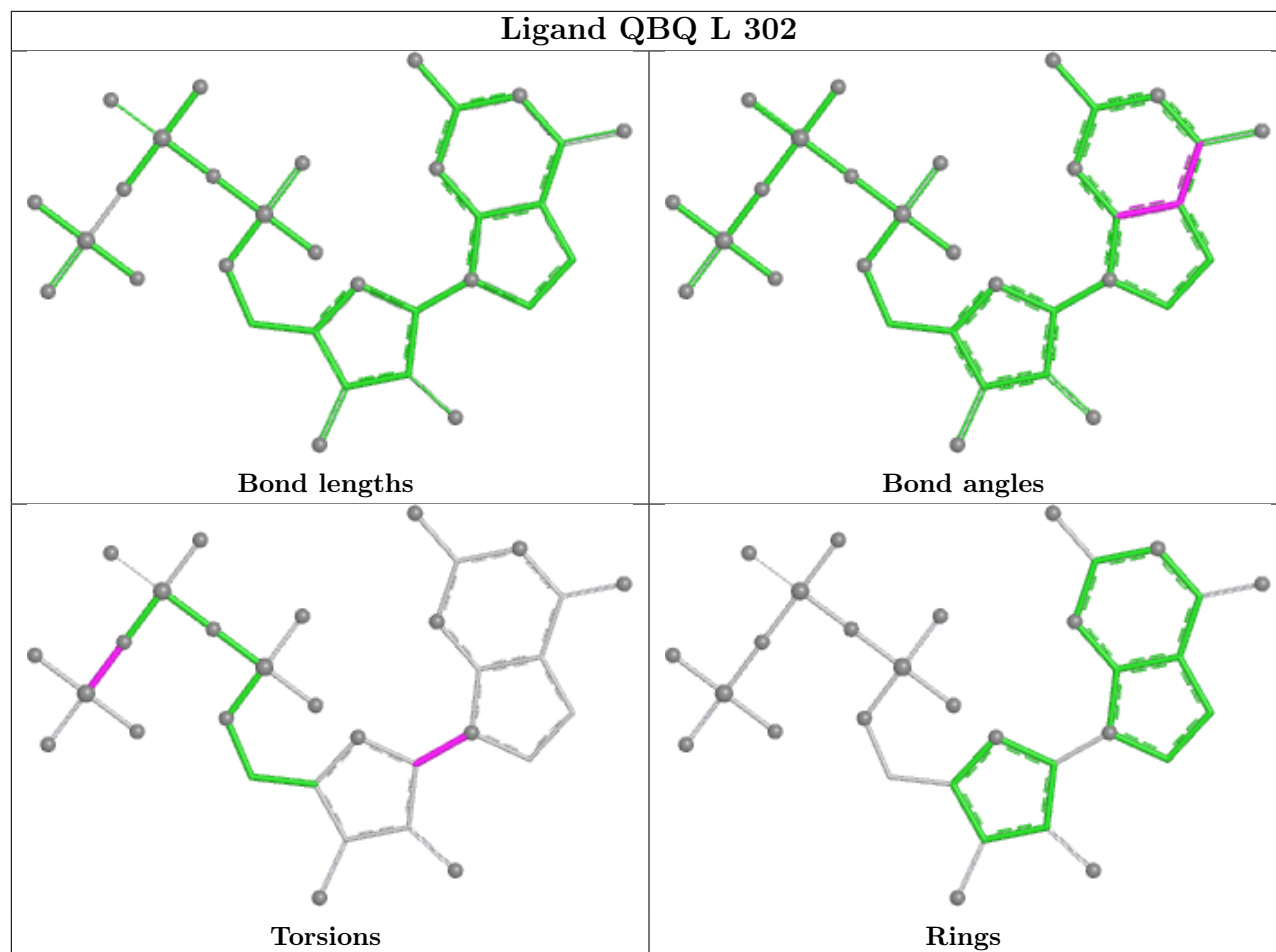












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2			OWAB(Å ²)	Q < 0.9
1	A	192/224 (85%)	0.11	6 (3%)	51	48	28, 40, 72, 90	0
1	B	190/224 (84%)	-0.09	1 (0%)	87	85	24, 39, 64, 82	0
1	C	191/224 (85%)	0.08	2 (1%)	79	77	28, 43, 70, 89	0
1	D	191/224 (85%)	0.01	4 (2%)	63	60	23, 43, 69, 98	0
1	E	190/224 (84%)	-0.07	3 (1%)	70	67	26, 40, 62, 86	0
1	F	190/224 (84%)	-0.10	1 (0%)	87	85	26, 38, 66, 86	0
1	G	193/224 (86%)	0.01	4 (2%)	63	60	27, 40, 74, 97	0
1	H	192/224 (85%)	0.03	5 (2%)	57	54	17, 41, 71, 88	1 (0%)
1	I	192/224 (85%)	0.09	5 (2%)	57	54	33, 45, 72, 85	0
1	J	190/224 (84%)	0.14	2 (1%)	78	76	27, 42, 75, 102	0
1	K	192/224 (85%)	0.09	4 (2%)	63	60	27, 45, 78, 108	0
1	L	191/224 (85%)	0.20	4 (2%)	63	60	33, 49, 78, 99	0
1	M	191/224 (85%)	0.10	5 (2%)	57	54	27, 42, 74, 98	0
1	N	193/224 (86%)	0.28	5 (2%)	57	54	29, 49, 79, 109	0
1	O	191/224 (85%)	0.24	5 (2%)	57	54	28, 48, 80, 110	0
1	P	189/224 (84%)	0.33	7 (3%)	45	42	29, 53, 79, 110	0
1	Q	190/224 (84%)	0.24	3 (1%)	70	67	33, 52, 78, 103	0
1	R	191/224 (85%)	0.23	5 (2%)	57	54	28, 44, 72, 98	0
1	S	188/224 (83%)	0.20	3 (1%)	70	67	32, 46, 78, 93	0
1	T	186/224 (83%)	0.30	6 (3%)	50	47	28, 47, 81, 93	0
All	All	3813/4480 (85%)	0.12	80 (2%)	63	60	17, 45, 76, 110	1 (0%)

All (80) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	N	58	PRO	4.6
1	H	250	SER	4.0
1	T	249	ARG	3.8
1	M	58	PRO	3.7
1	T	66	LEU	3.5
1	N	250	SER	3.5
1	R	60	SER	3.4
1	Q	117	LEU	3.4
1	G	249	ARG	3.3
1	N	249	ARG	3.3
1	O	250	SER	3.3
1	A	123	ASP	3.2
1	A	161	GLN	3.2
1	D	59	ARG	3.1
1	G	125	ASP	3.0
1	G	58	PRO	3.0
1	K	249	ARG	3.0
1	T	64	ASN	3.0
1	L	58	PRO	3.0
1	A	58	PRO	2.9
1	H	249	ARG	2.8
1	D	66	LEU	2.8
1	R	249	ARG	2.8
1	M	248	ILE	2.8
1	A	160	LYS	2.8
1	R	93	LYS	2.6
1	P	159	ASN	2.6
1	M	59	ARG	2.6
1	K	93	LYS	2.6
1	I	197	LEU	2.6
1	J	66	LEU	2.6
1	S	107	LYS	2.5
1	B	117	LEU	2.5
1	I	239	LYS	2.5
1	O	125	ASP	2.5
1	N	59	ARG	2.4
1	O	62	GLU	2.4
1	J	59	ARG	2.4
1	Q	59	ARG	2.4
1	K	250	SER	2.4
1	P	248	ILE	2.4
1	O	66	LEU	2.4
1	H	196	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	T	221	MET	2.4
1	I	117	LEU	2.3
1	A	196	ALA	2.3
1	I	112	THR	2.3
1	A	249	ARG	2.3
1	P	60	SER	2.3
1	M	115	ASP	2.3
1	O	60	SER	2.2
1	K	92	LEU	2.2
1	P	115	ASP	2.2
1	I	58	PRO	2.2
1	T	113	ILE	2.2
1	M	117	LEU	2.2
1	E	198	ARG	2.2
1	E	115	ASP	2.2
1	C	66	LEU	2.1
1	P	221	MET	2.1
1	G	250	SER	2.1
1	N	112	THR	2.1
1	P	195	GLU	2.1
1	H	116	VAL	2.1
1	S	250	SER	2.1
1	L	224	LYS	2.1
1	L	239	LYS	2.1
1	L	116	VAL	2.1
1	D	112	THR	2.1
1	R	112	THR	2.1
1	R	221	MET	2.1
1	D	197	LEU	2.1
1	Q	248	ILE	2.1
1	T	197	LEU	2.0
1	P	246	THR	2.0
1	E	197	LEU	2.0
1	C	58	PRO	2.0
1	F	59	ARG	2.0
1	H	97[A]	ARG	2.0
1	S	249	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

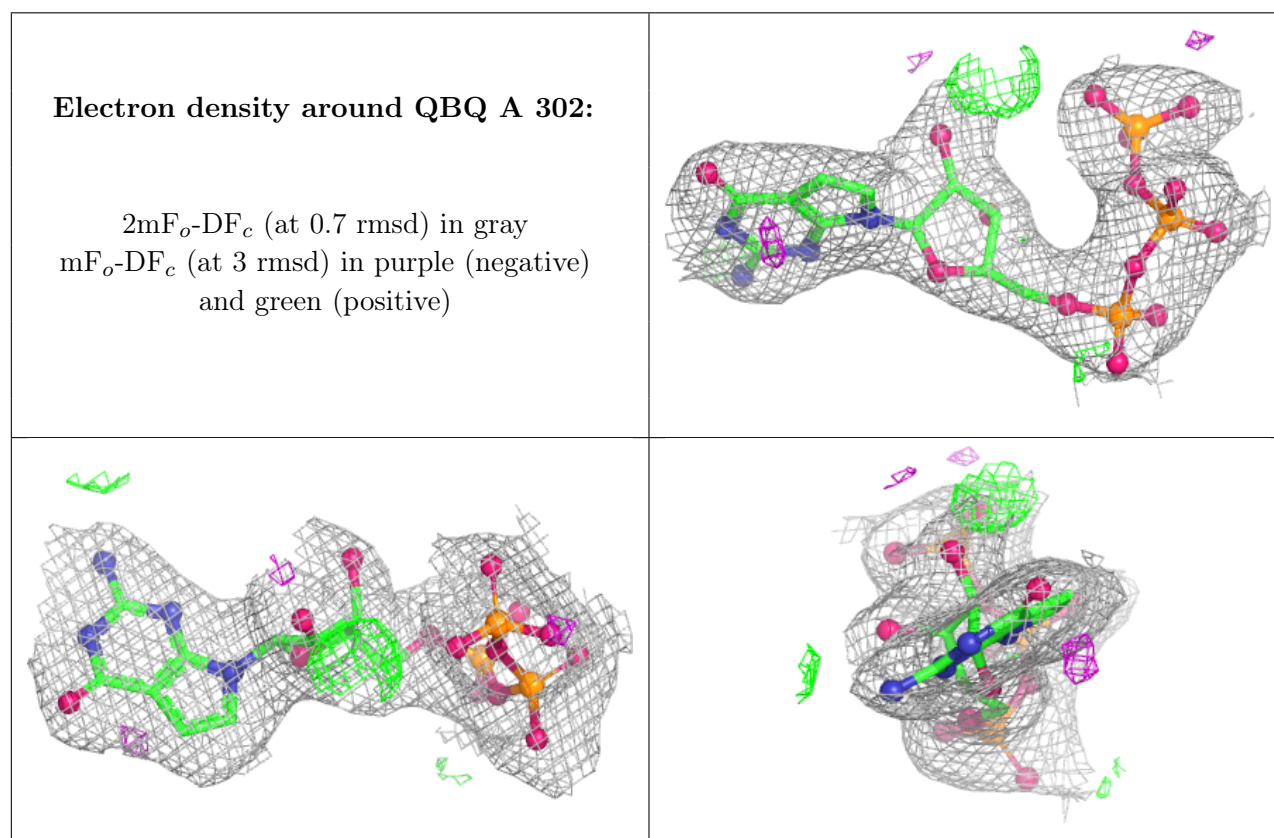
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	QBQ	A	302	32/32	0.96	0.07	44,48,52,54	0
3	QBQ	G	302	32/32	0.96	0.07	43,46,49,49	0
3	QBQ	I	302	32/32	0.96	0.06	44,47,50,52	0
3	QBQ	P	302	32/32	0.96	0.07	40,44,48,49	0
3	QBQ	S	302	32/32	0.96	0.07	51,53,56,56	0
3	QBQ	F	302	32/32	0.97	0.05	30,36,41,41	0
2	ZN	P	301	1/1	0.97	0.06	69,69,69,69	0
3	QBQ	H	302	32/32	0.97	0.06	33,36,38,39	0
2	ZN	L	301	1/1	0.97	0.05	67,67,67,67	0
3	QBQ	L	302	32/32	0.97	0.07	43,47,51,51	0
3	QBQ	M	302	32/32	0.97	0.06	47,49,49,49	0
3	QBQ	N	302	32/32	0.97	0.06	40,42,44,44	0
3	QBQ	D	302	32/32	0.97	0.06	37,43,50,51	0
3	QBQ	R	302	32/32	0.97	0.07	43,45,49,50	0
3	QBQ	E	302	32/32	0.97	0.06	32,38,41,42	0
3	QBQ	T	302	32/32	0.97	0.06	40,42,44,46	0
3	QBQ	J	302	32/32	0.98	0.04	32,37,40,41	0
3	QBQ	K	302	32/32	0.98	0.05	28,34,38,40	0
3	QBQ	C	302	32/32	0.98	0.06	36,40,45,46	0
2	ZN	A	301	1/1	0.98	0.03	51,51,51,51	0
2	ZN	D	301	1/1	0.98	0.04	51,51,51,51	0
3	QBQ	O	302	32/32	0.98	0.05	38,39,43,44	0
2	ZN	Q	301	1/1	0.98	0.06	61,61,61,61	0
3	QBQ	Q	302	32/32	0.98	0.05	39,42,46,47	0
2	ZN	T	301	1/1	0.98	0.03	64,64,64,64	0
2	ZN	F	301	1/1	0.98	0.03	45,45,45,45	0
3	QBQ	B	302	32/32	0.98	0.05	33,37,39,39	0
2	ZN	O	301	1/1	0.99	0.04	62,62,62,62	0
2	ZN	E	301	1/1	0.99	0.02	53,53,53,53	0
2	ZN	C	301	1/1	0.99	0.04	55,55,55,55	0
2	ZN	R	301	1/1	0.99	0.03	52,52,52,52	0
2	ZN	S	301	1/1	0.99	0.03	58,58,58,58	0

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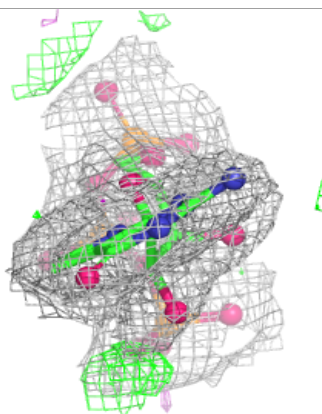
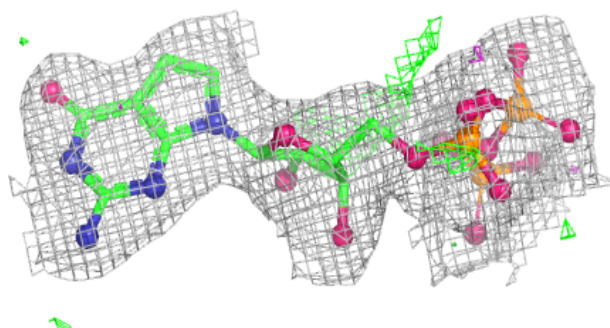
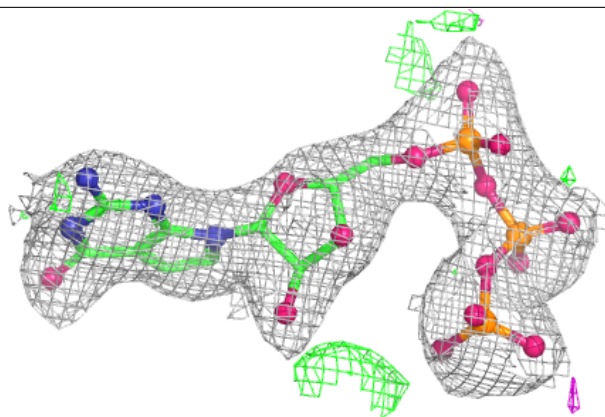
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	G	301	1/1	0.99	0.03	52,52,52,52	0
2	ZN	H	301	1/1	0.99	0.02	56,56,56,56	0
2	ZN	I	301	1/1	0.99	0.03	53,53,53,53	0
2	ZN	J	301	1/1	0.99	0.02	54,54,54,54	0
2	ZN	K	301	1/1	0.99	0.03	62,62,62,62	0
2	ZN	B	301	1/1	0.99	0.02	54,54,54,54	0
2	ZN	M	301	1/1	0.99	0.05	55,55,55,55	0
2	ZN	N	301	1/1	0.99	0.03	64,64,64,64	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

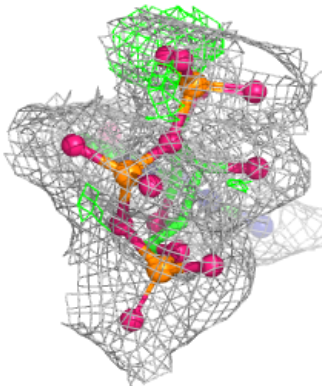
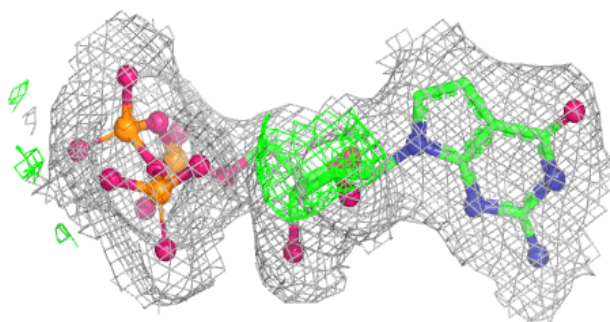
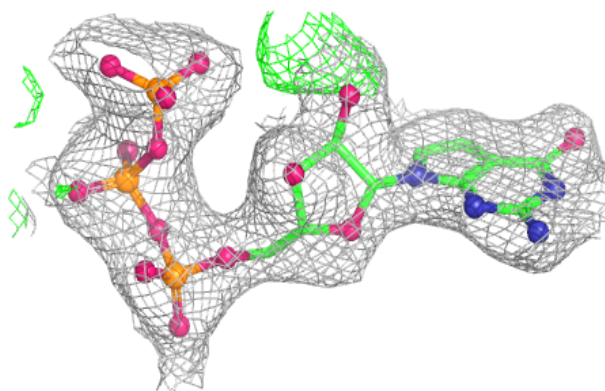


Electron density around QBQ G 302:

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

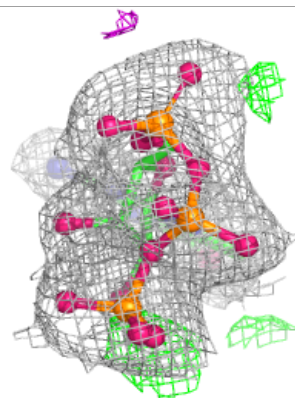
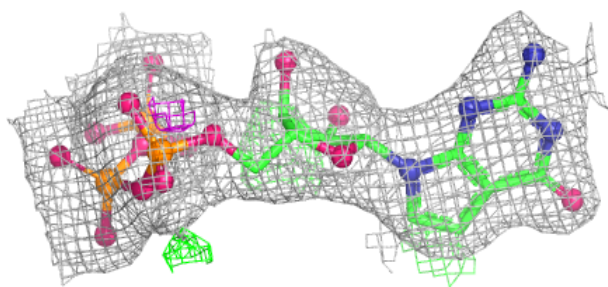
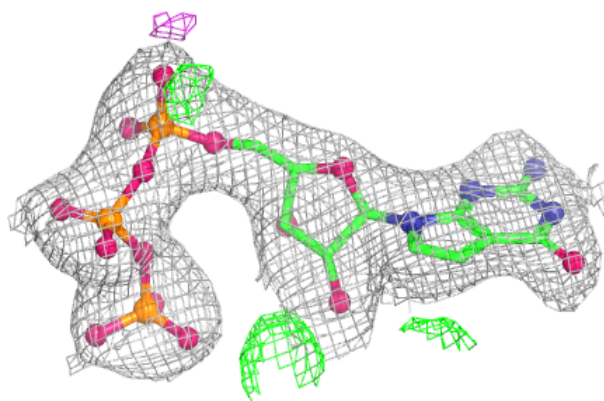
**Electron density around QBQ I 302:**

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

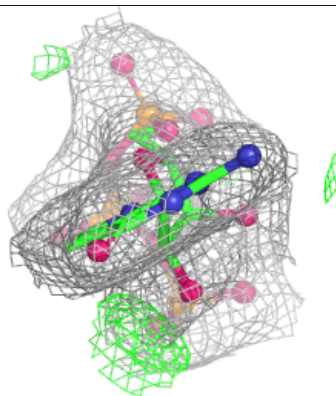
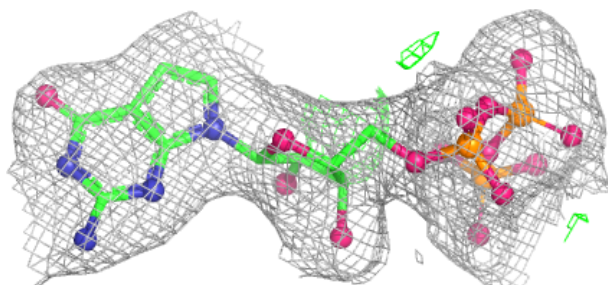
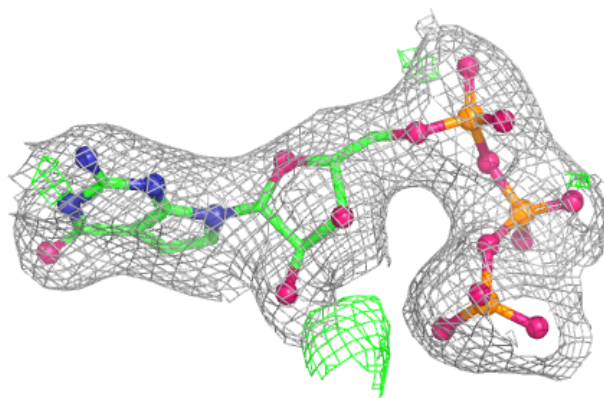


Electron density around QBQ P 302:

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

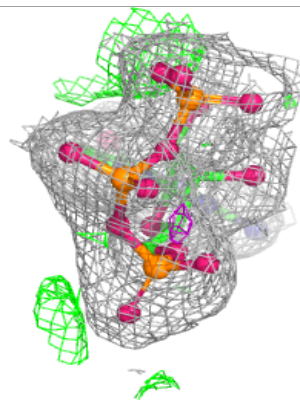
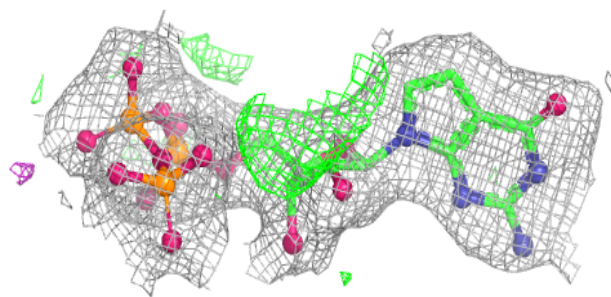
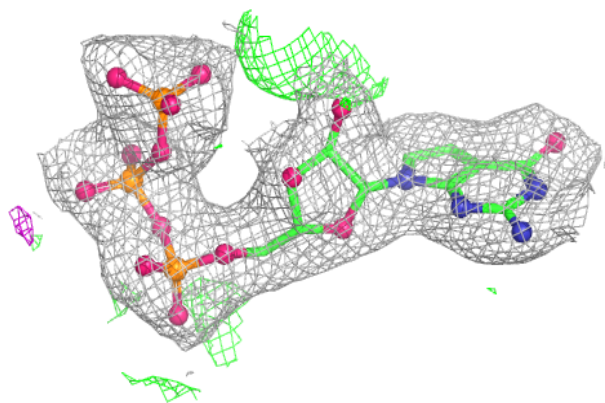
**Electron density around QBQ S 302:**

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



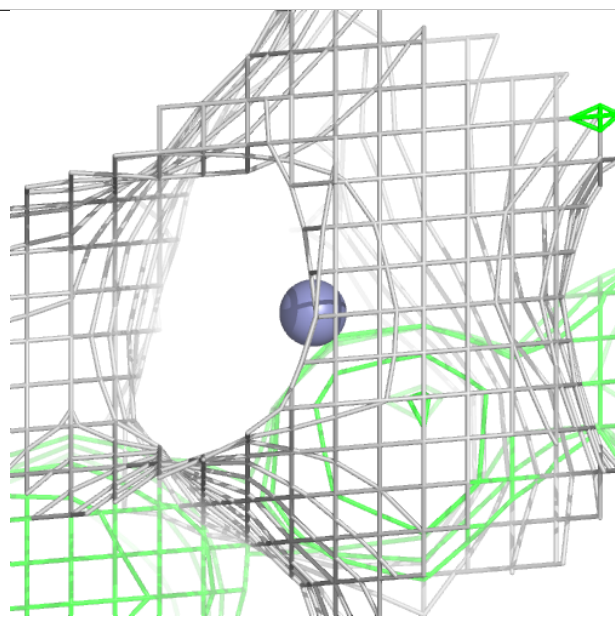
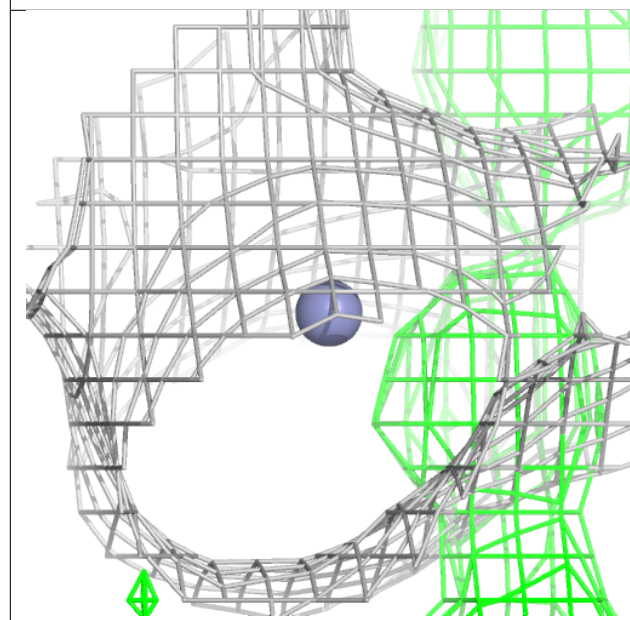
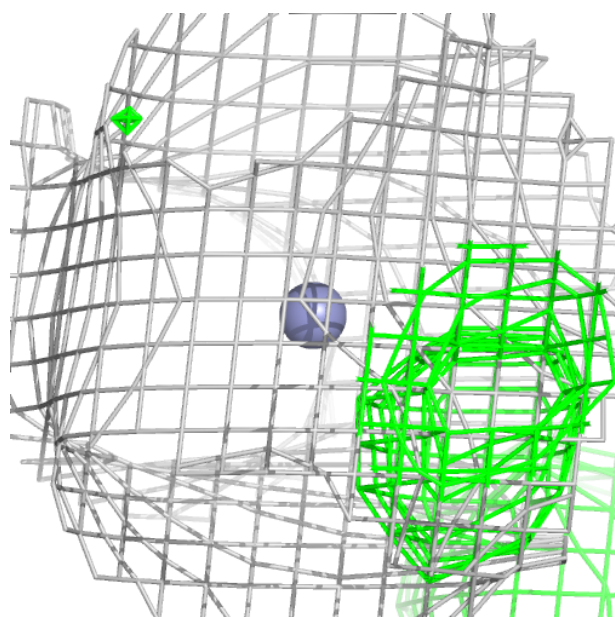
Electron density around QBQ F 302:

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



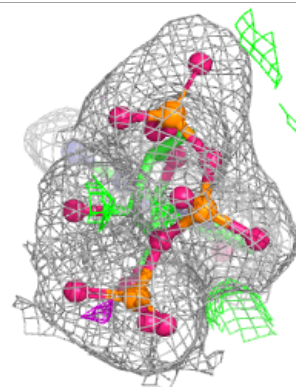
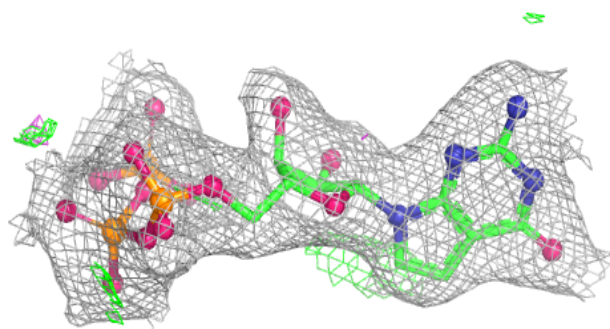
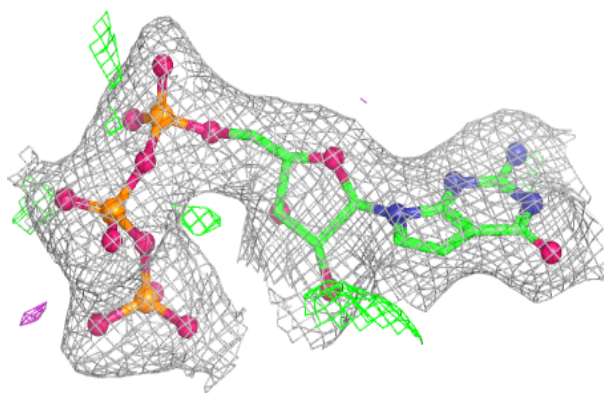
Electron density around ZN P 301:

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



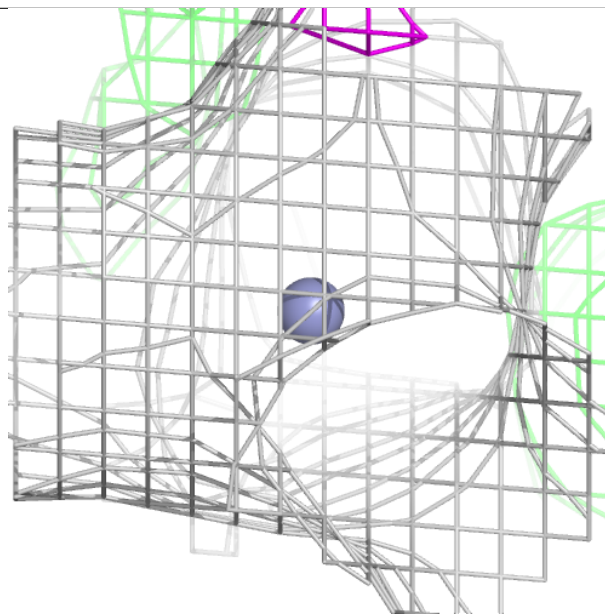
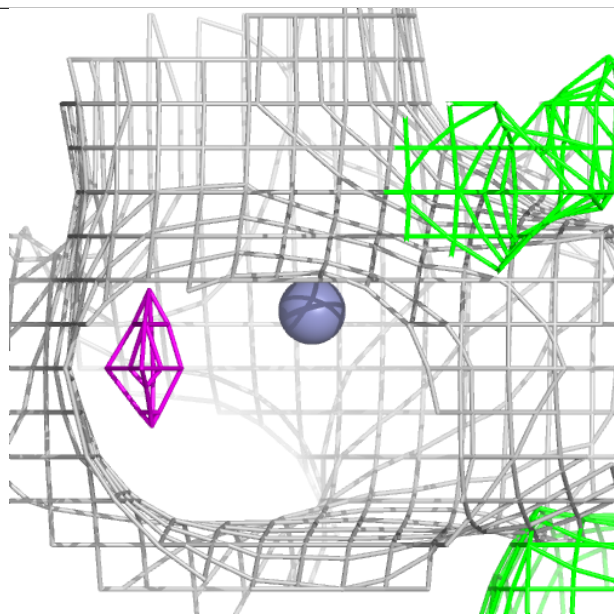
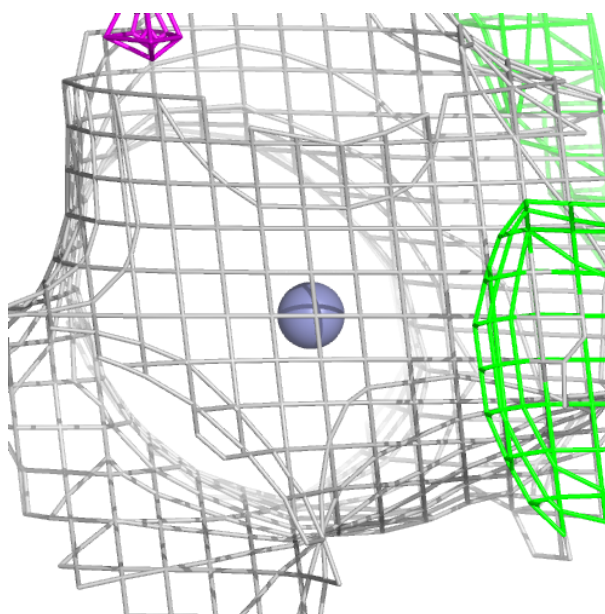
Electron density around QBQ H 302:

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



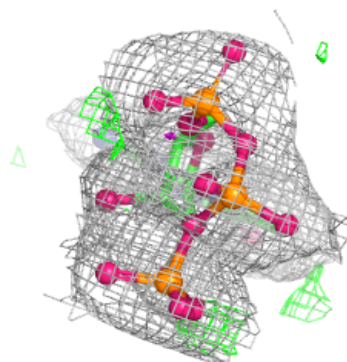
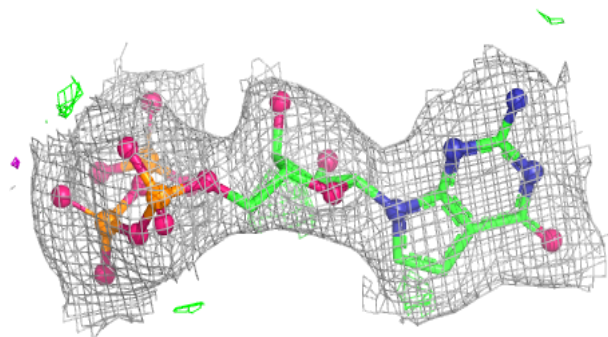
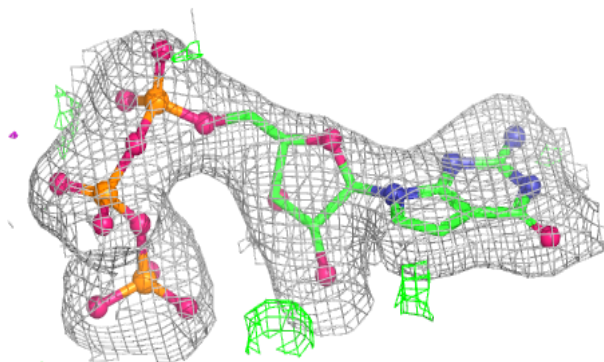
Electron density around ZN L 301:

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

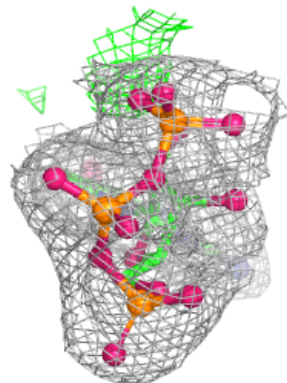
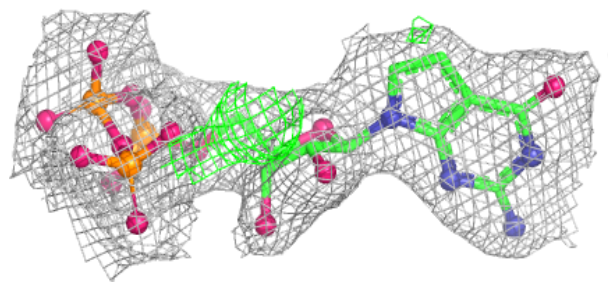
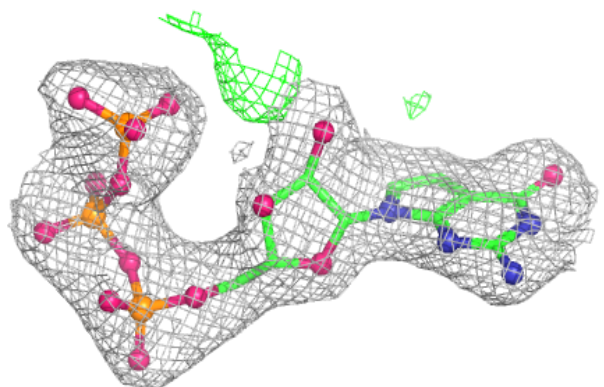


Electron density around QBQ L 302:

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

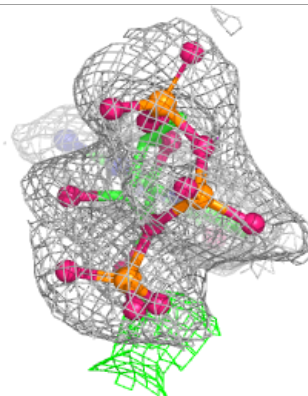
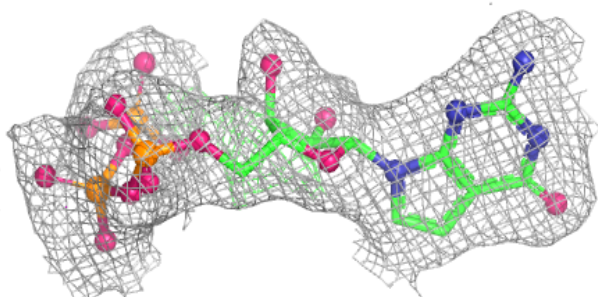
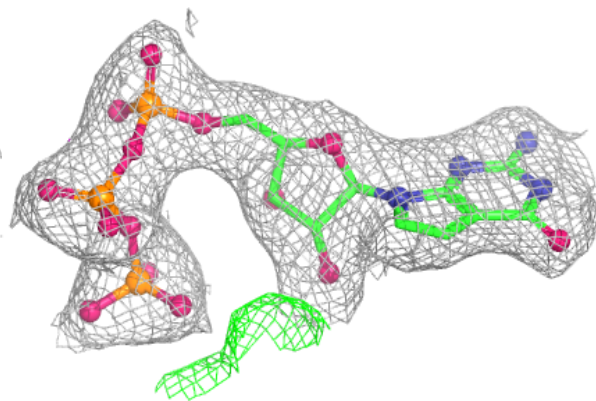
**Electron density around QBQ M 302:**

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

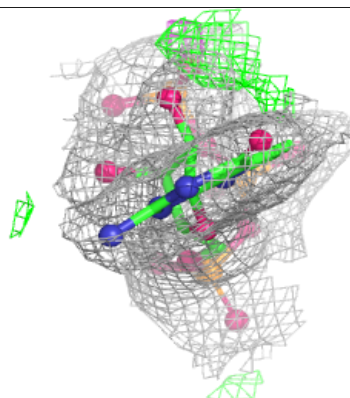
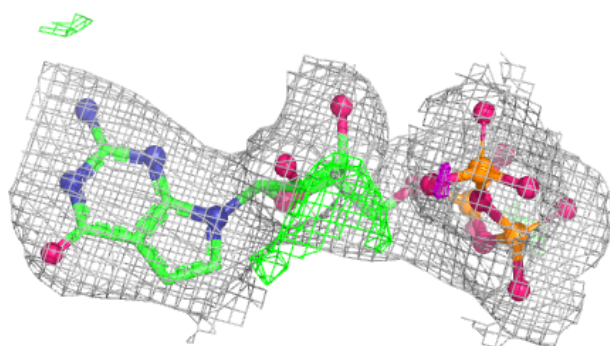
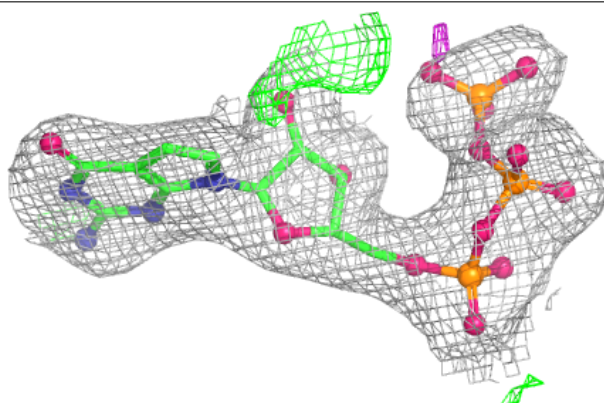


Electron density around QBQ N 302:

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

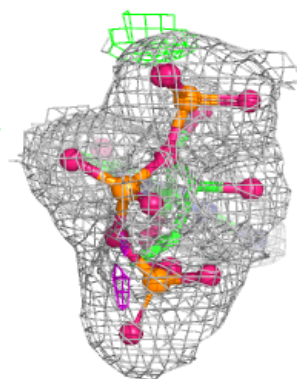
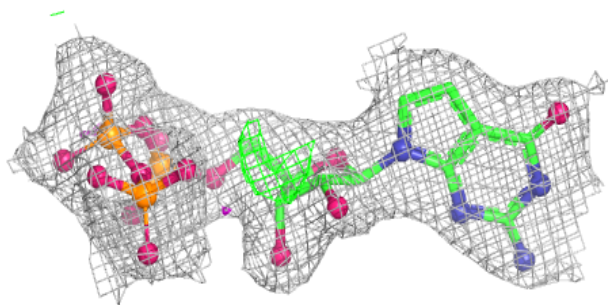
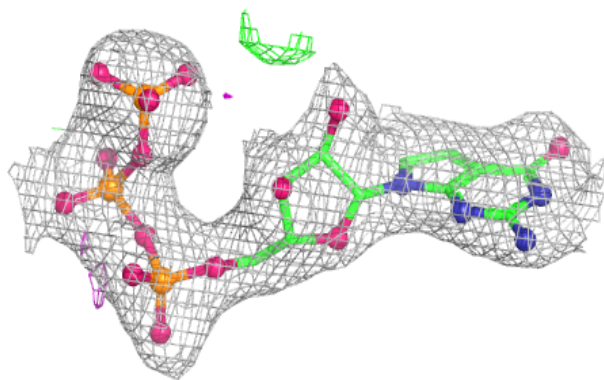
**Electron density around QBQ D 302:**

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

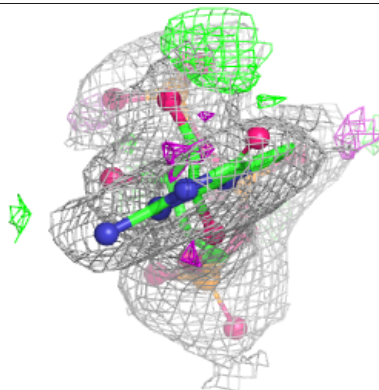
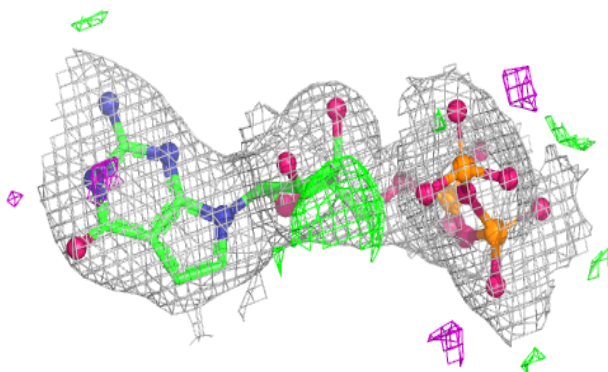
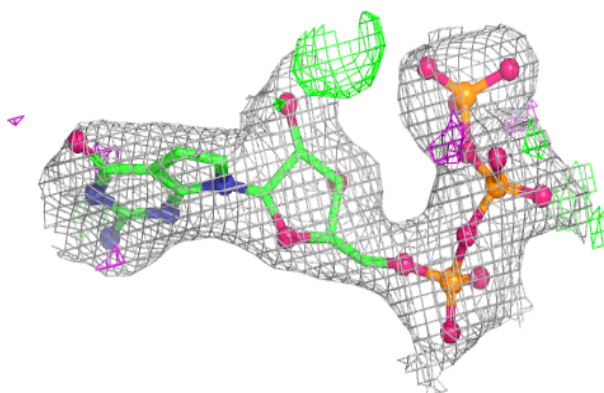


Electron density around QBQ R 302:

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

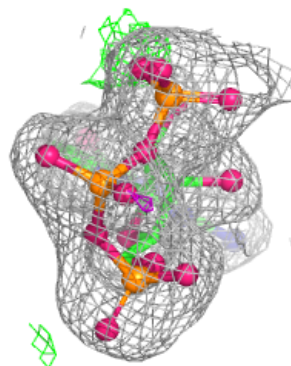
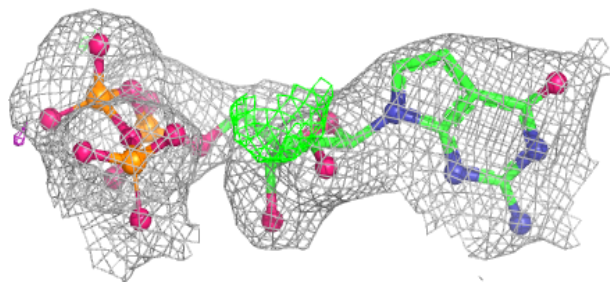
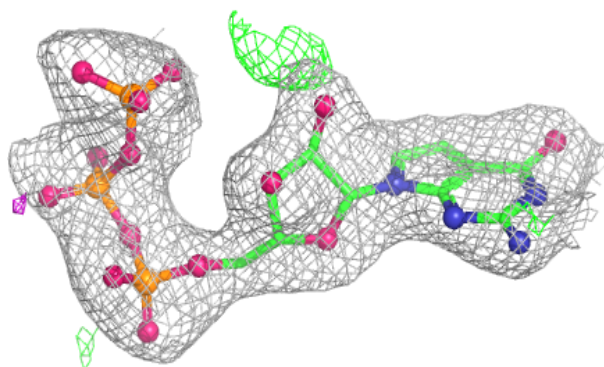
**Electron density around QBQ E 302:**

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

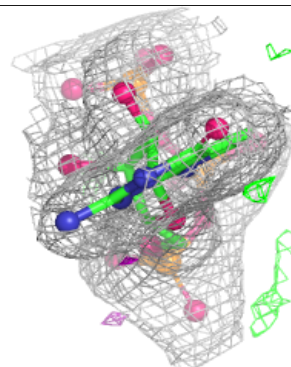
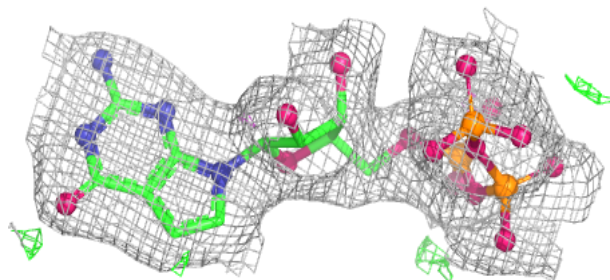
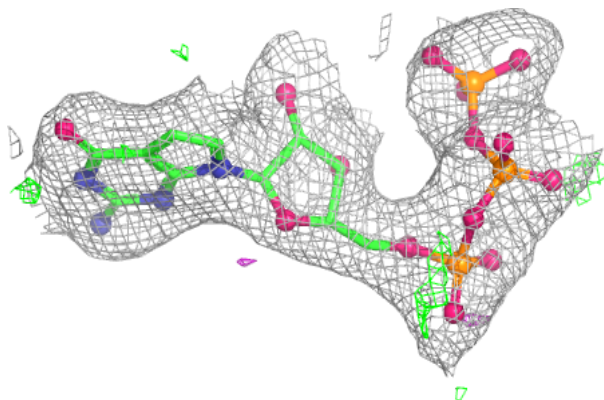


Electron density around QBQ T 302:

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

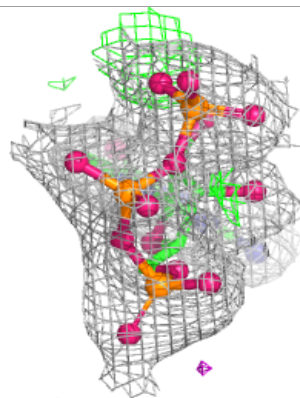
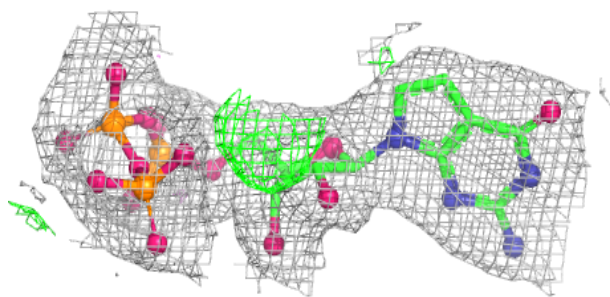
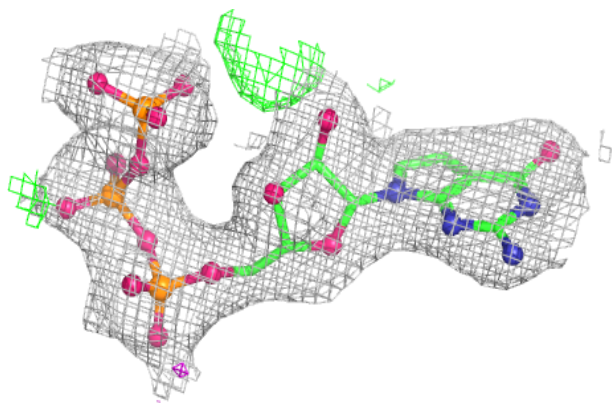
**Electron density around QBQ J 302:**

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

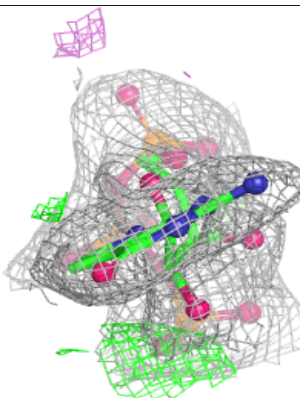
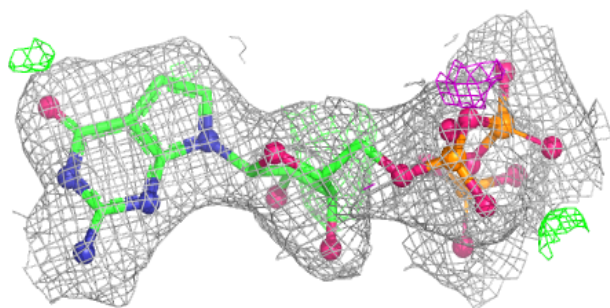
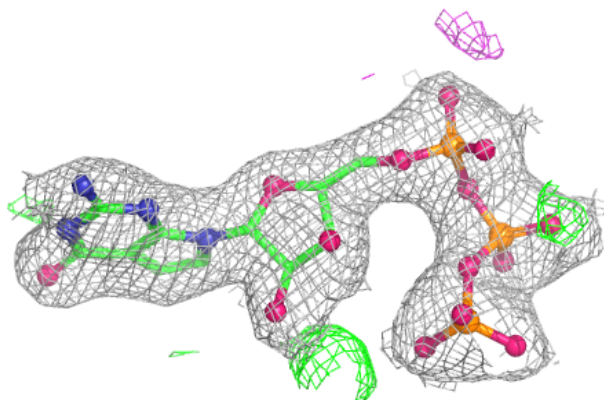


Electron density around QBQ K 302:

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

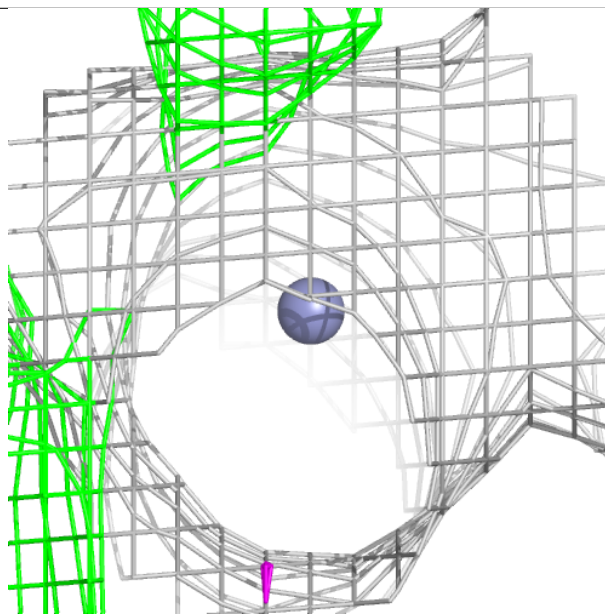
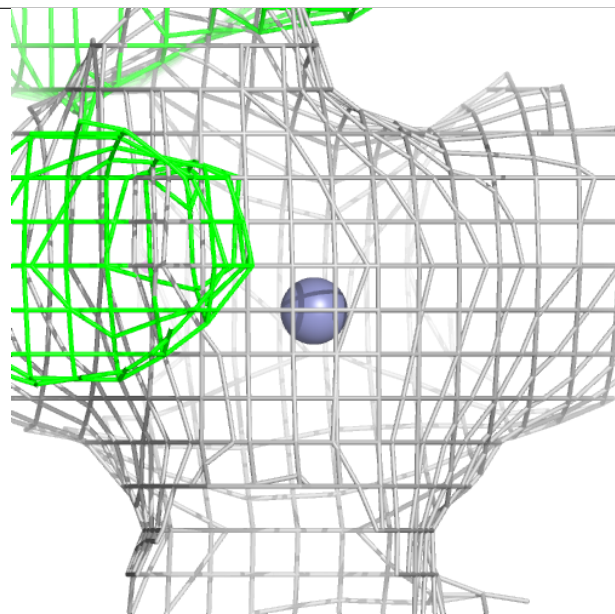
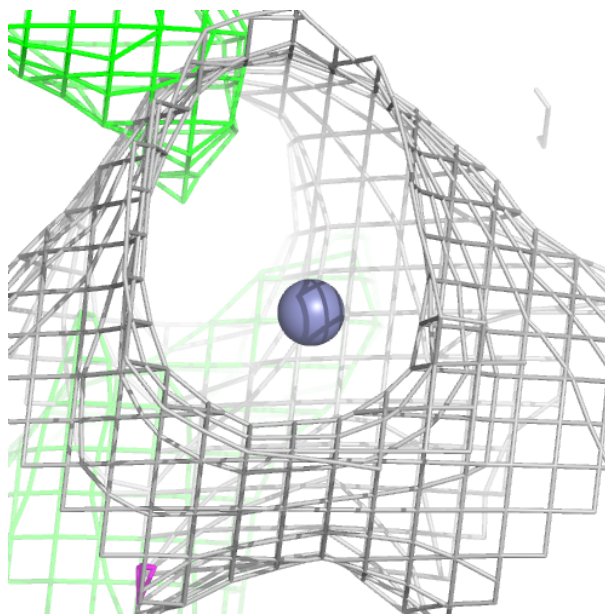
**Electron density around QBQ C 302:**

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



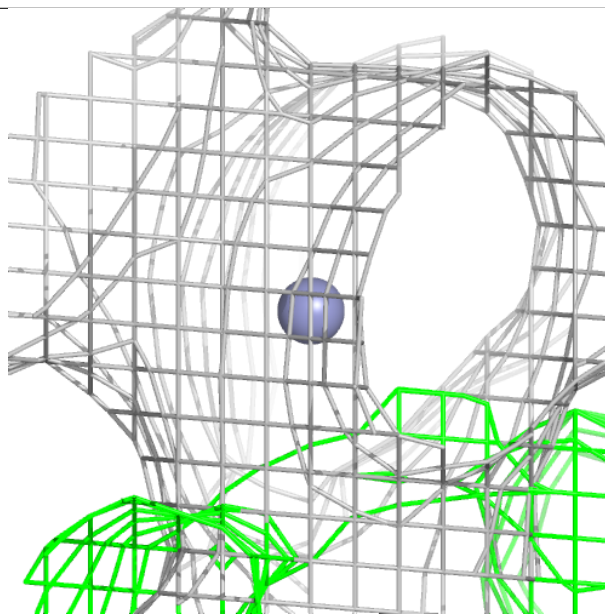
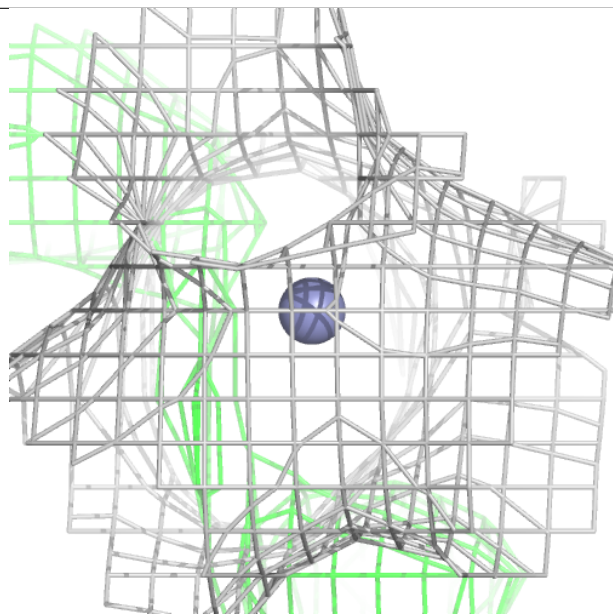
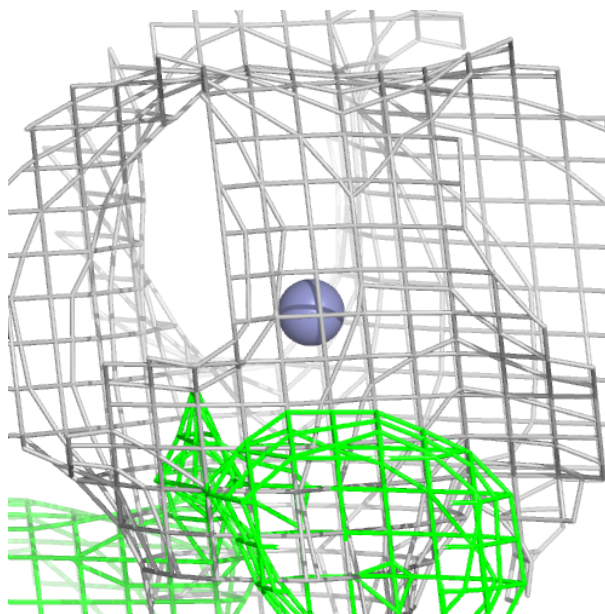
Electron density around ZN A 301:

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



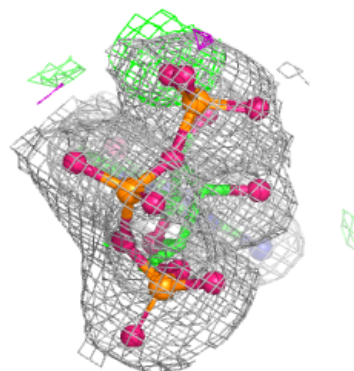
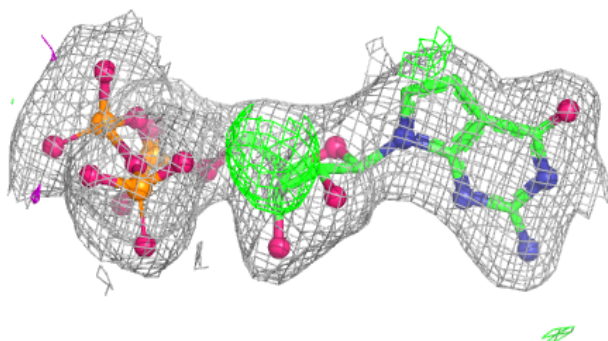
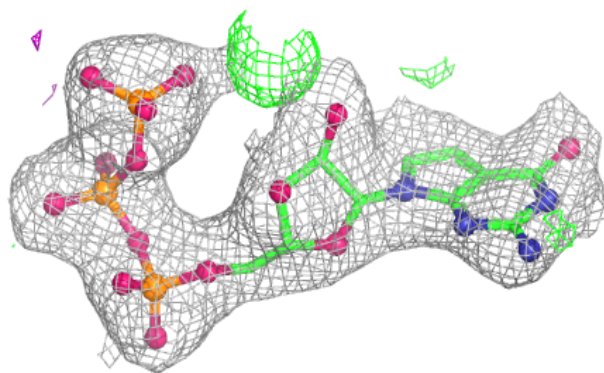
Electron density around ZN D 301:

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



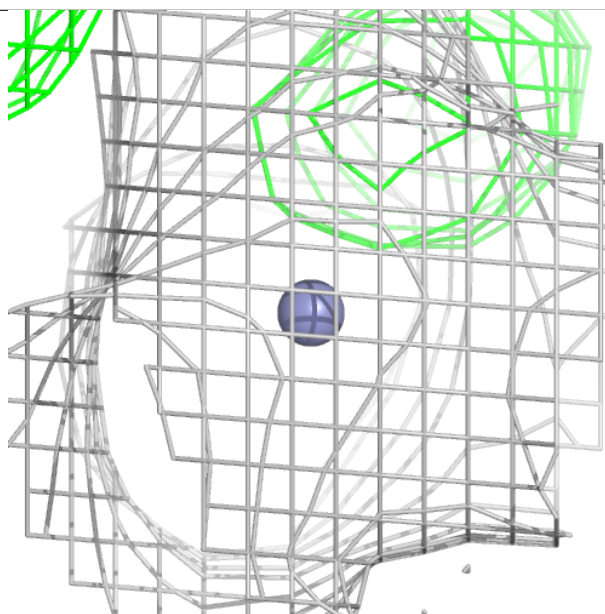
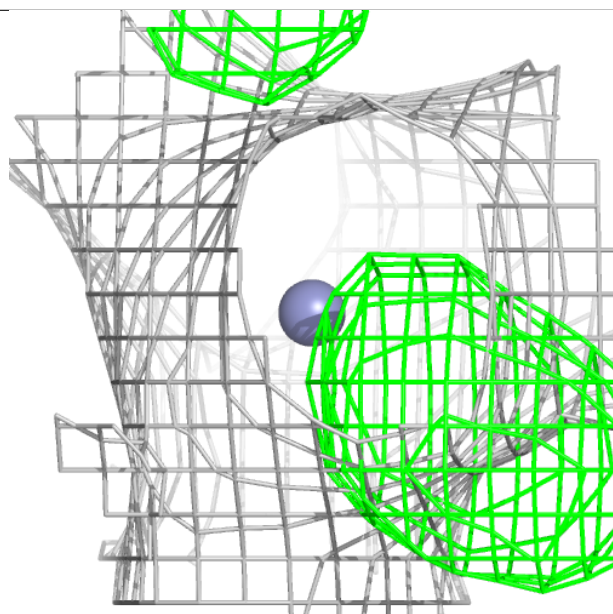
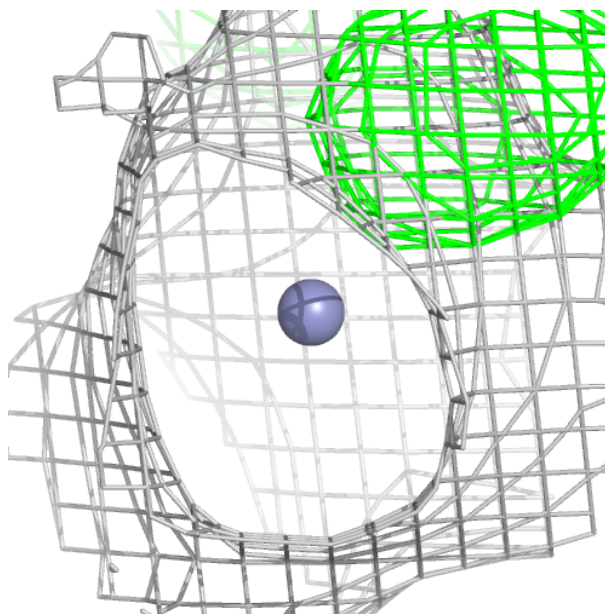
Electron density around QBQ O 302:

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



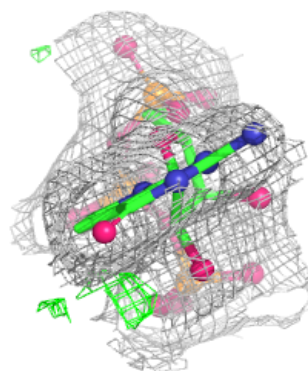
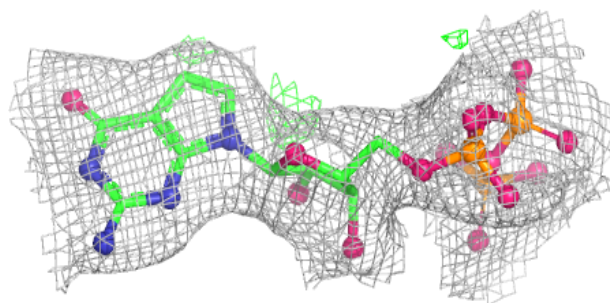
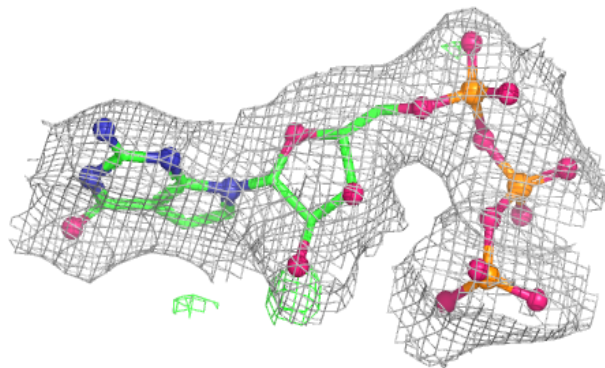
Electron density around ZN Q 301:

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



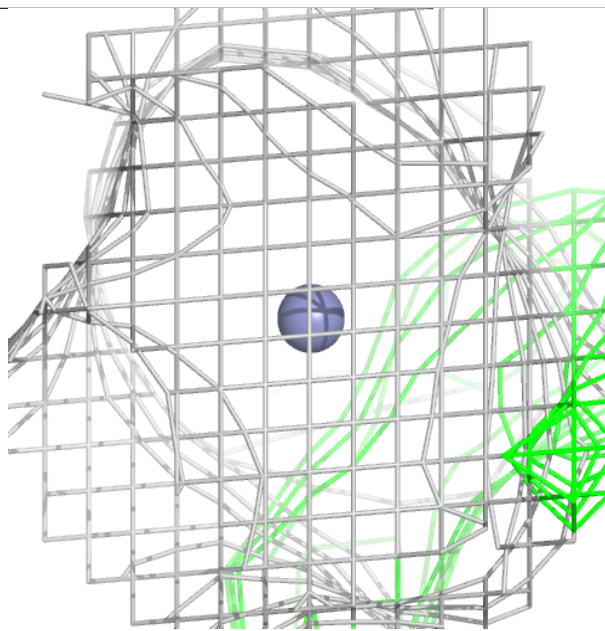
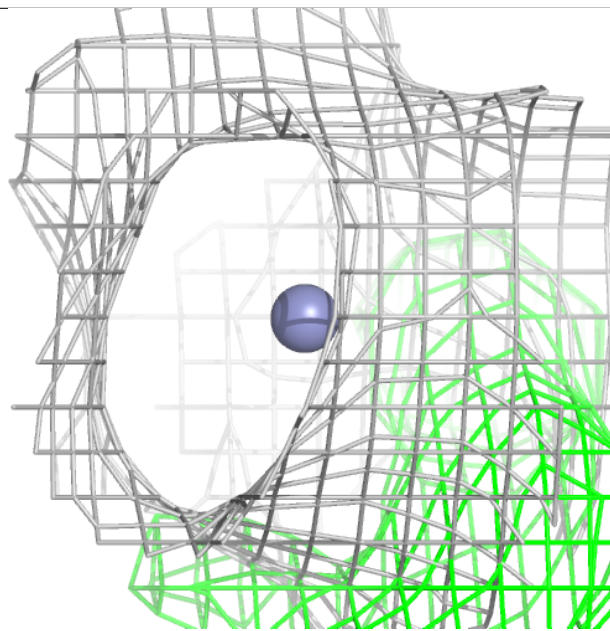
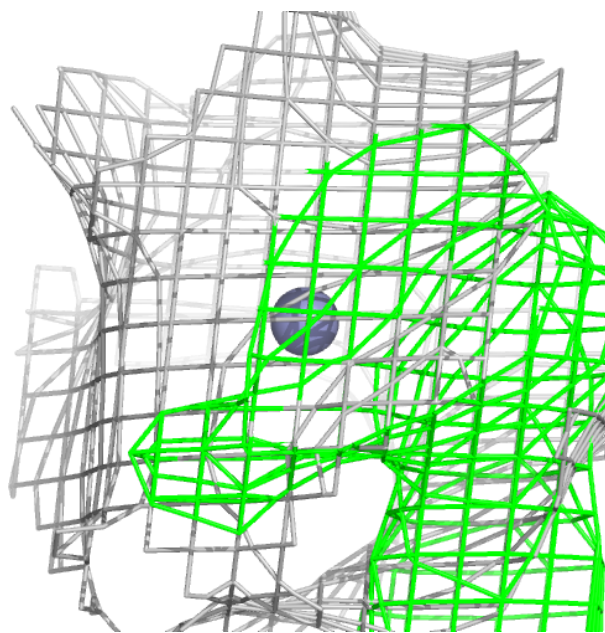
Electron density around QBQ Q 302:

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



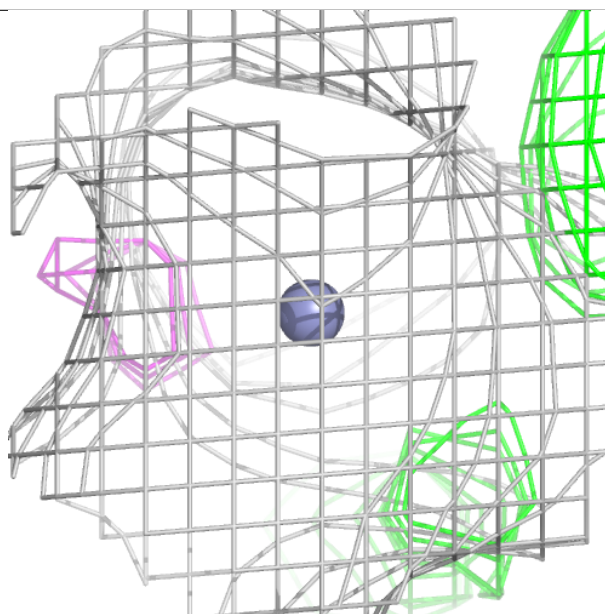
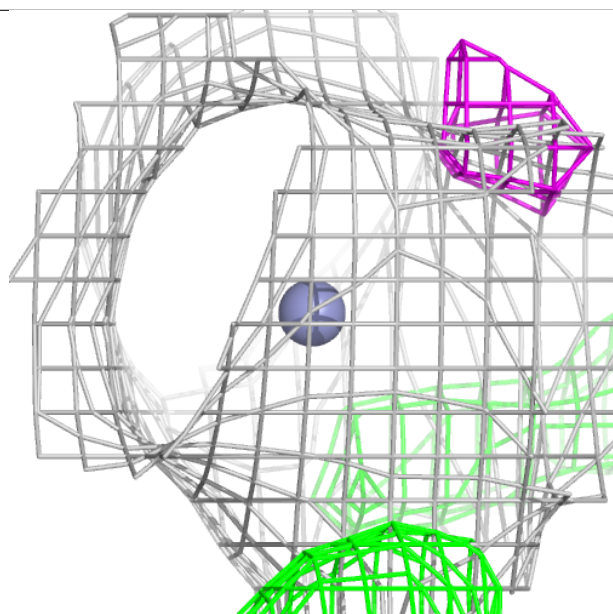
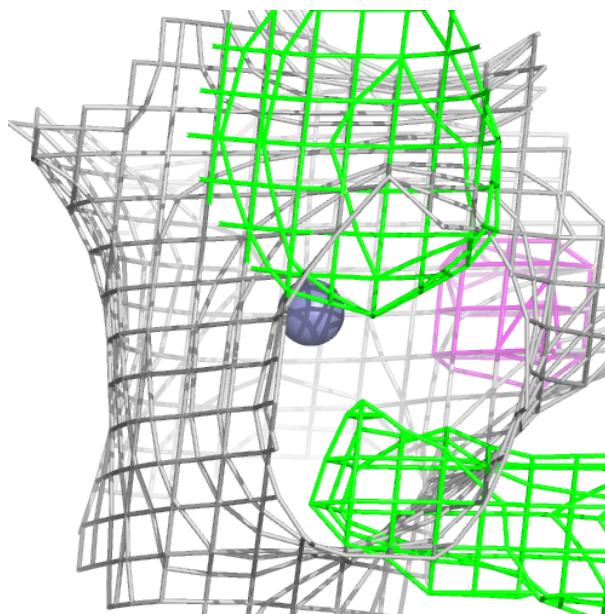
Electron density around ZN T 301:

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



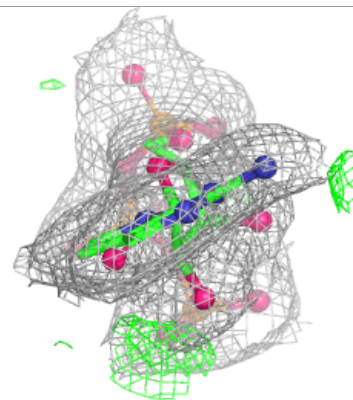
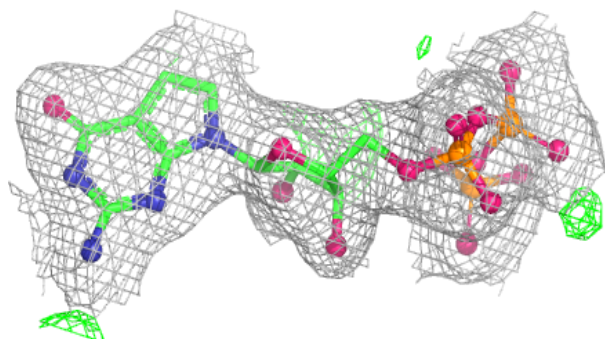
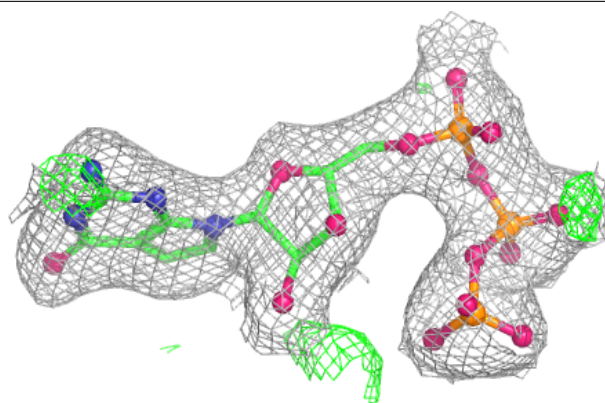
Electron density around ZN F 301:

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



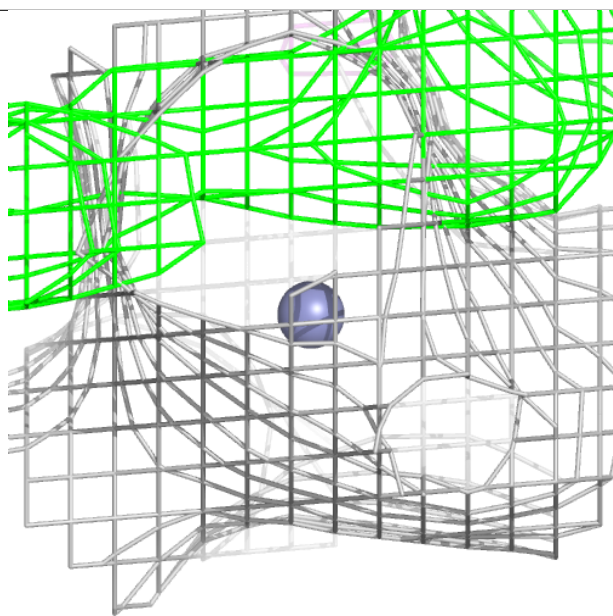
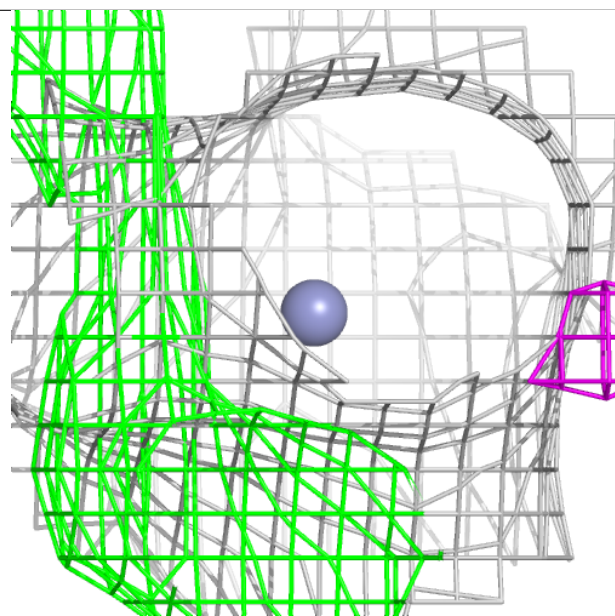
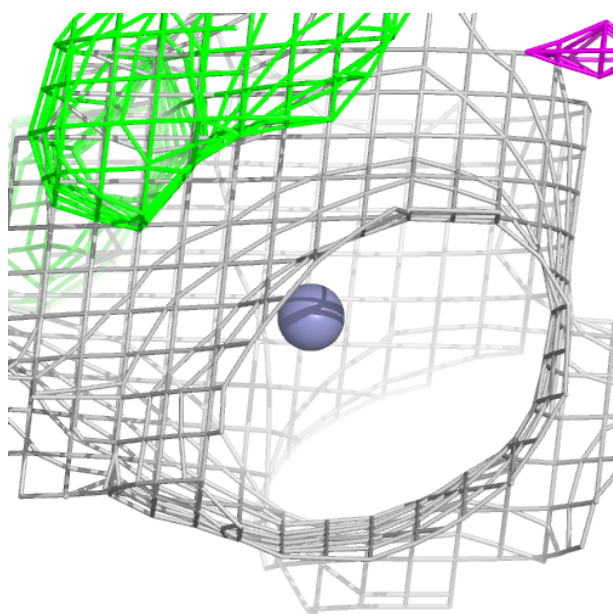
Electron density around QBQ B 302:

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



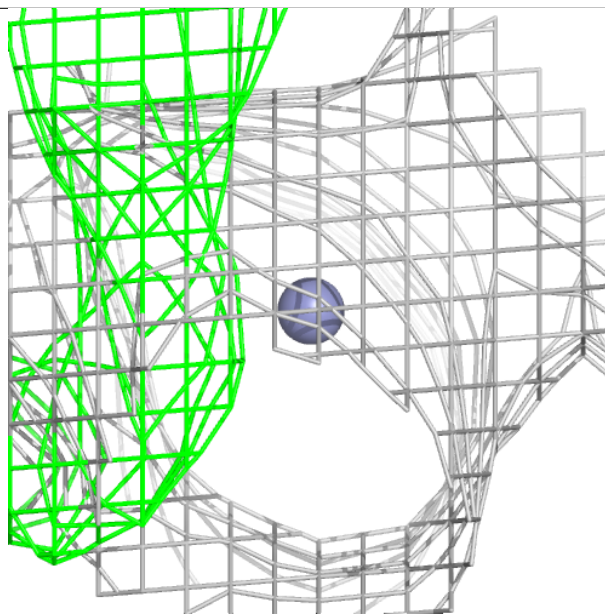
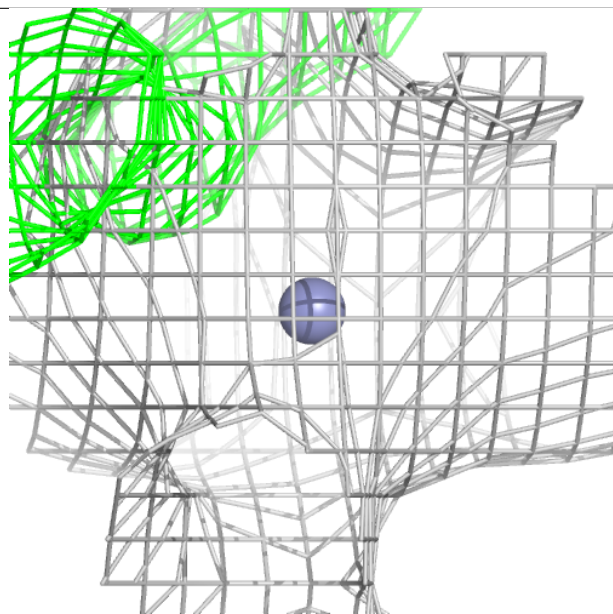
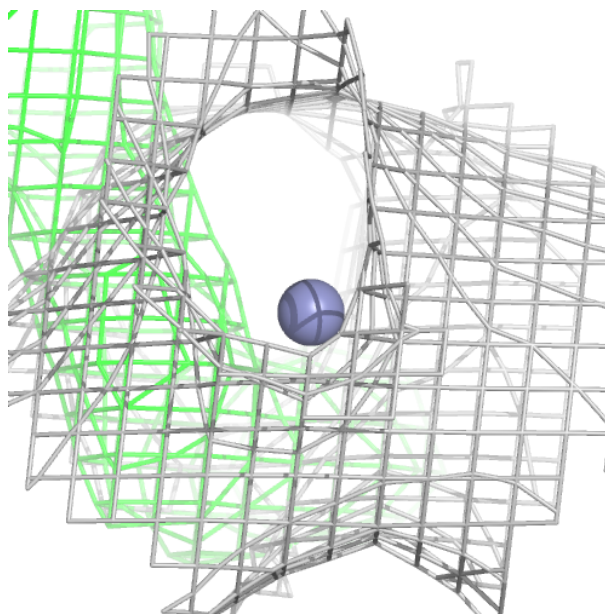
Electron density around ZN O 301:

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



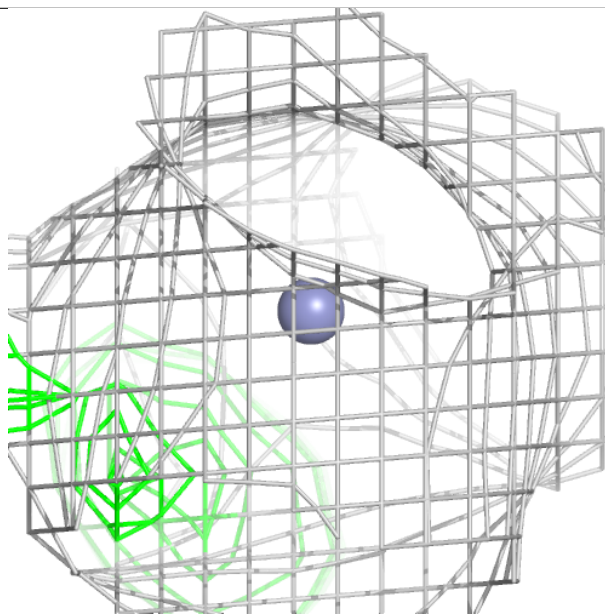
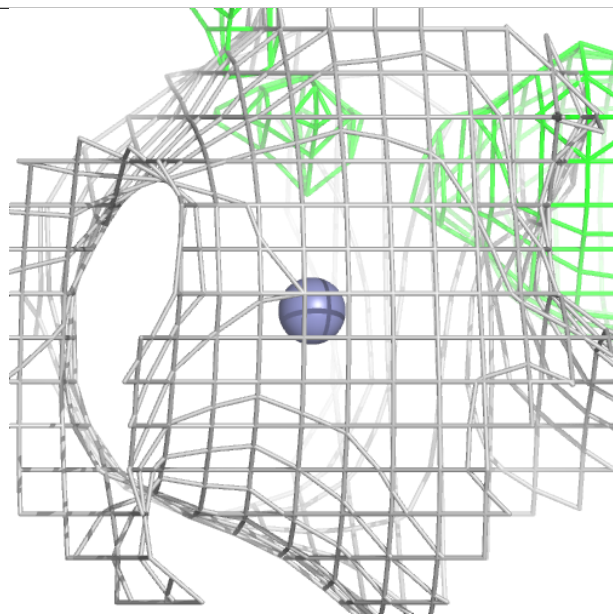
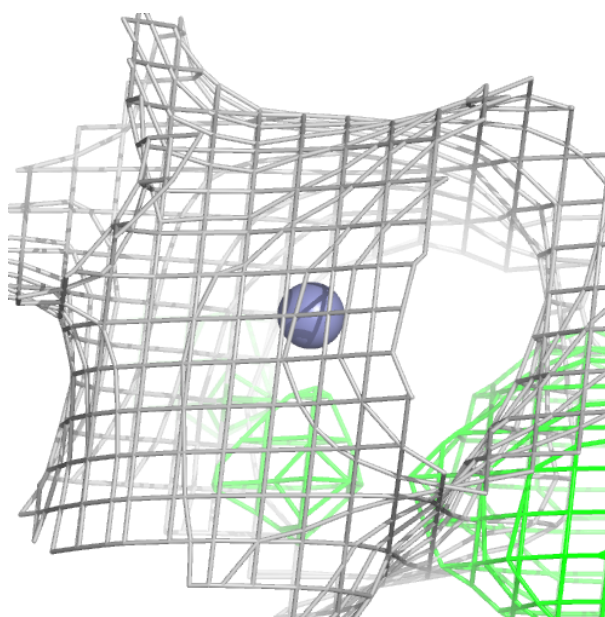
Electron density around ZN E 301:

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



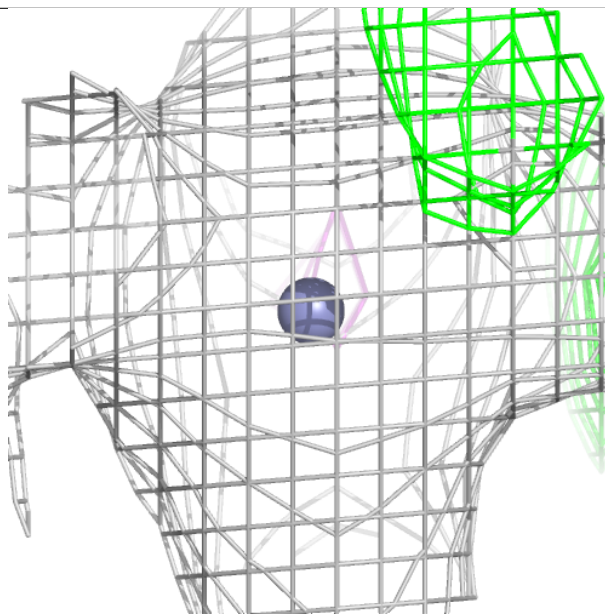
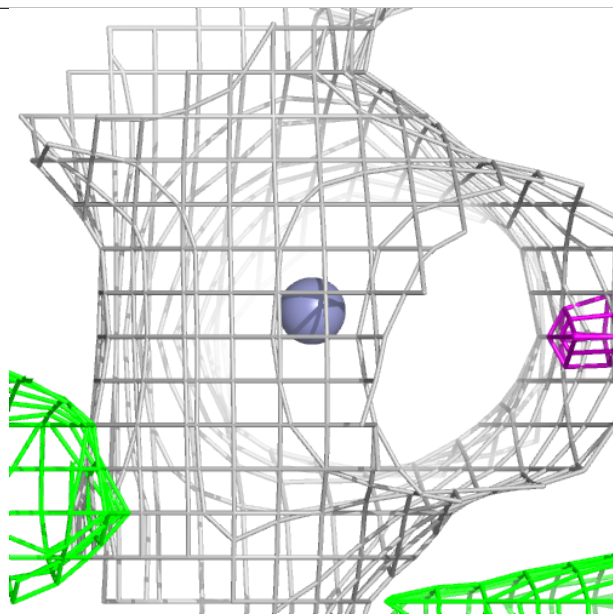
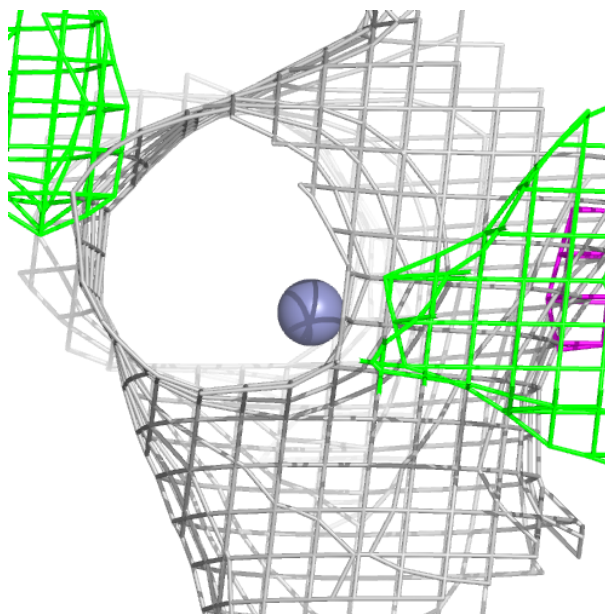
Electron density around ZN C 301:

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



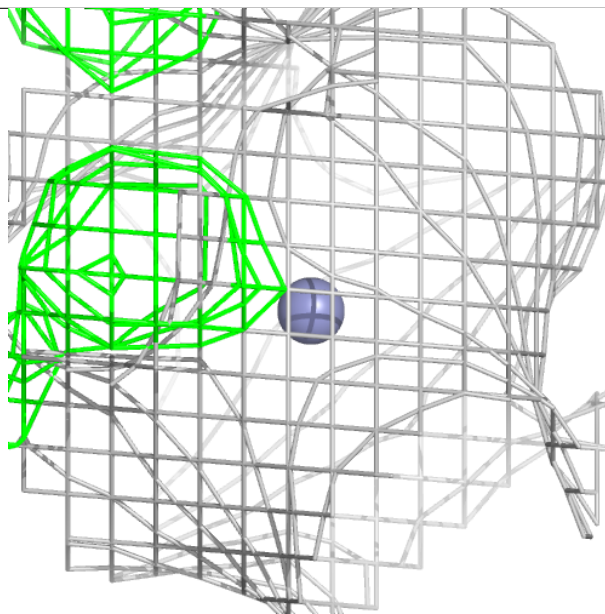
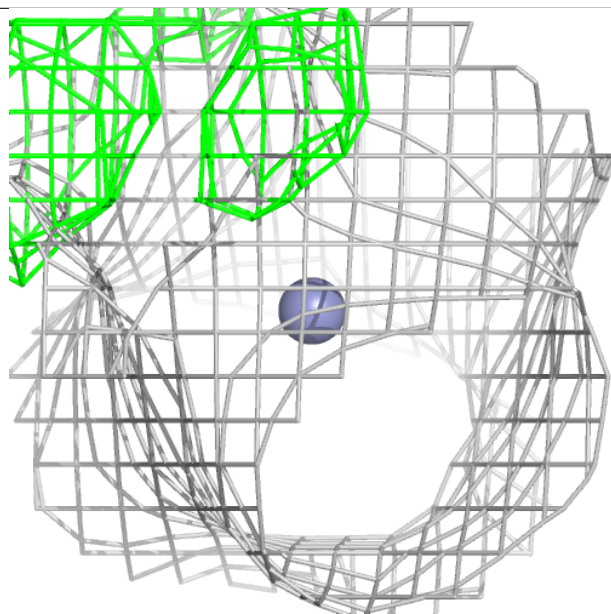
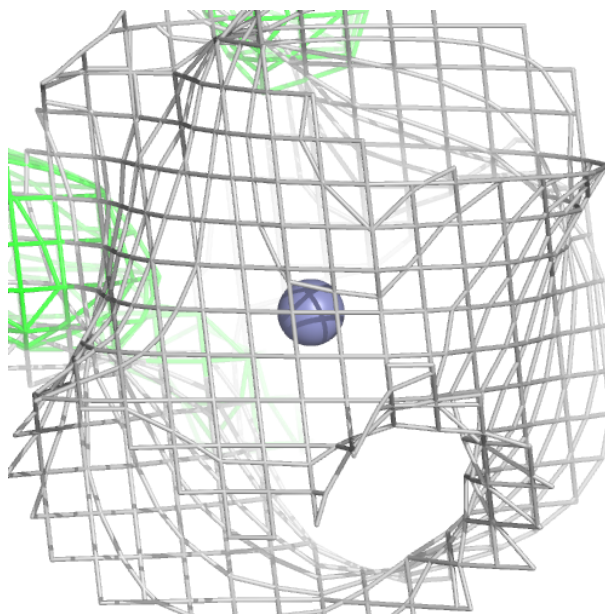
Electron density around ZN R 301:

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



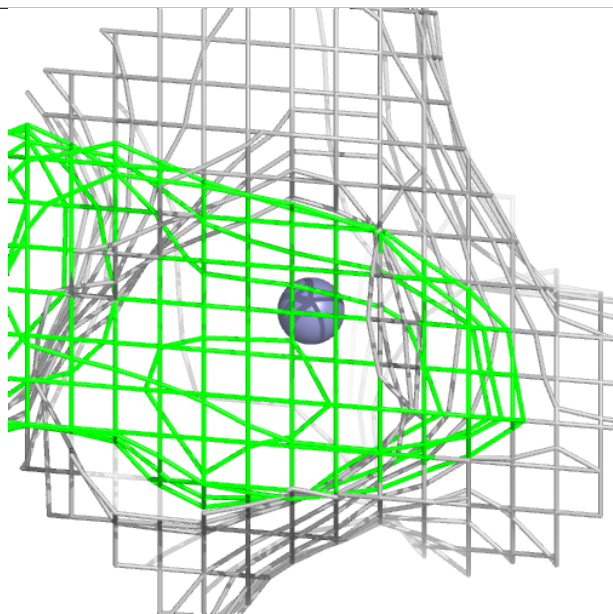
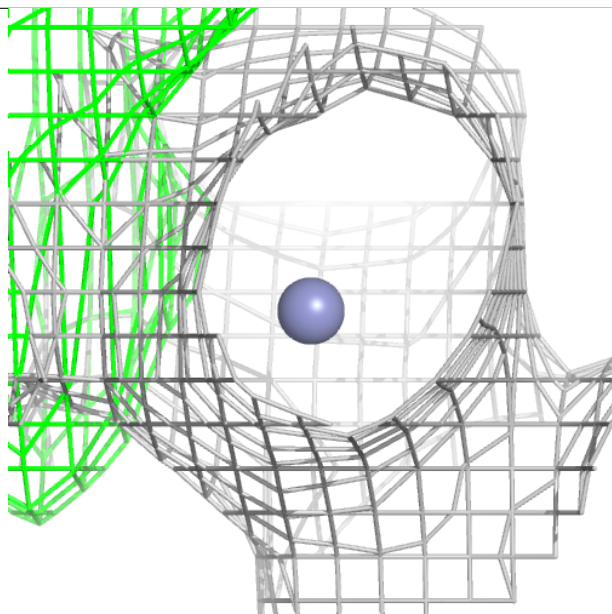
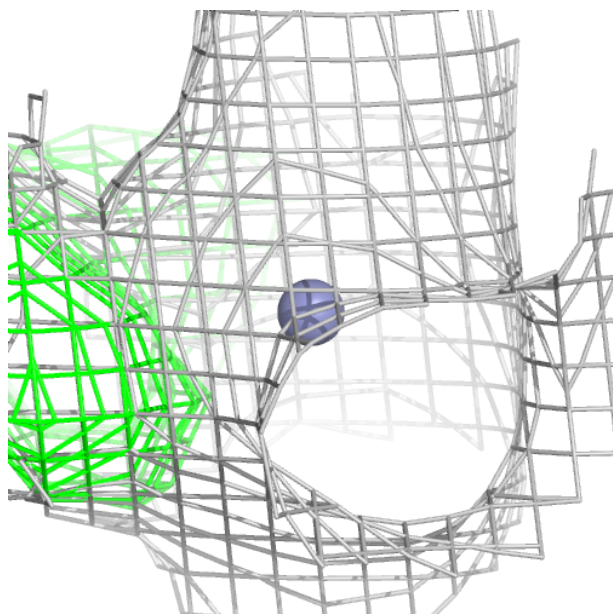
Electron density around ZN S 301:

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



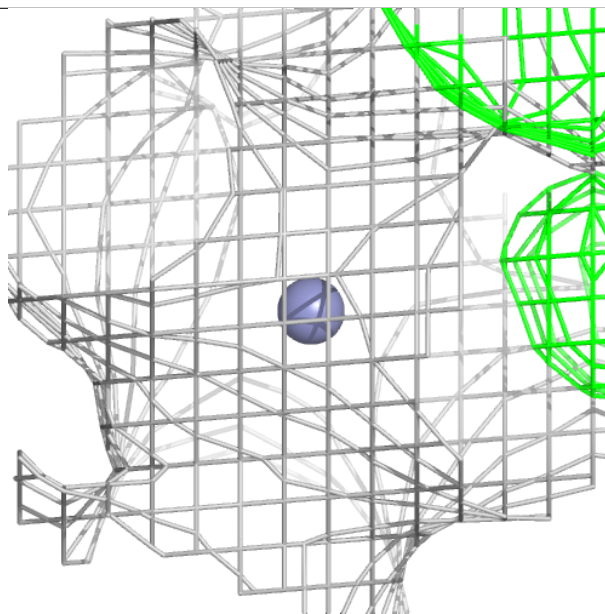
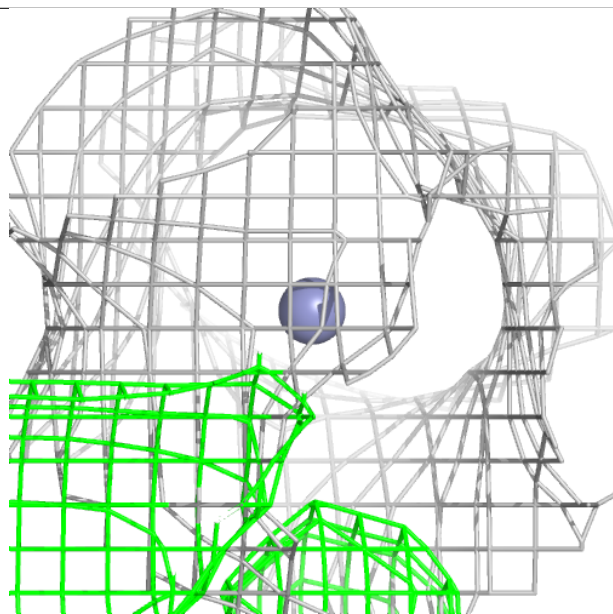
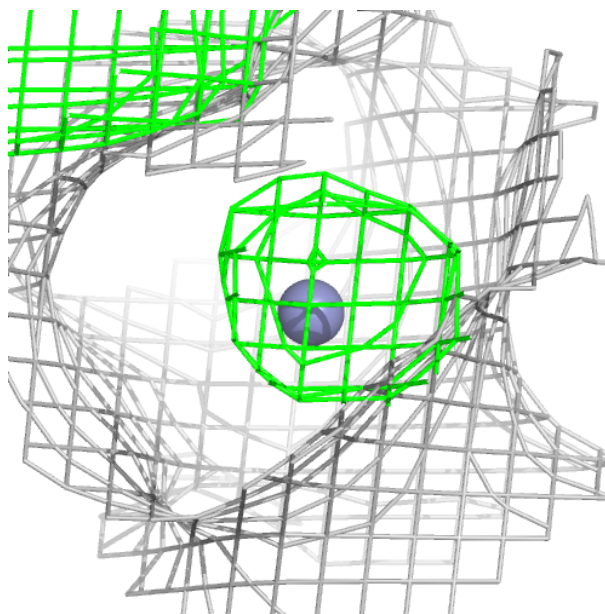
Electron density around ZN G 301:

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



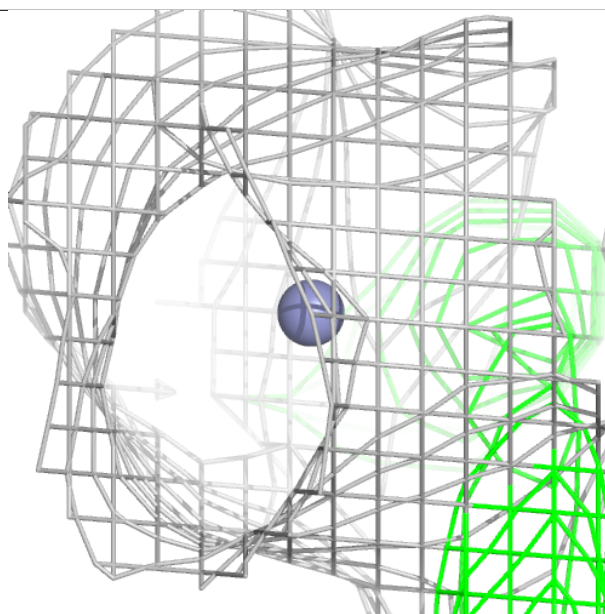
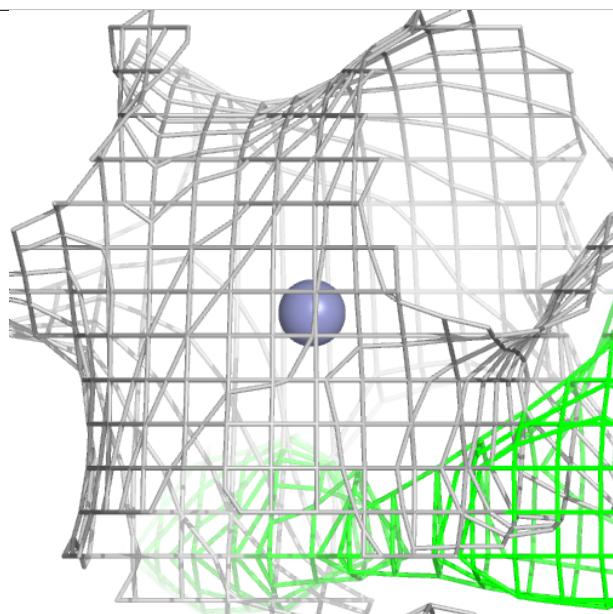
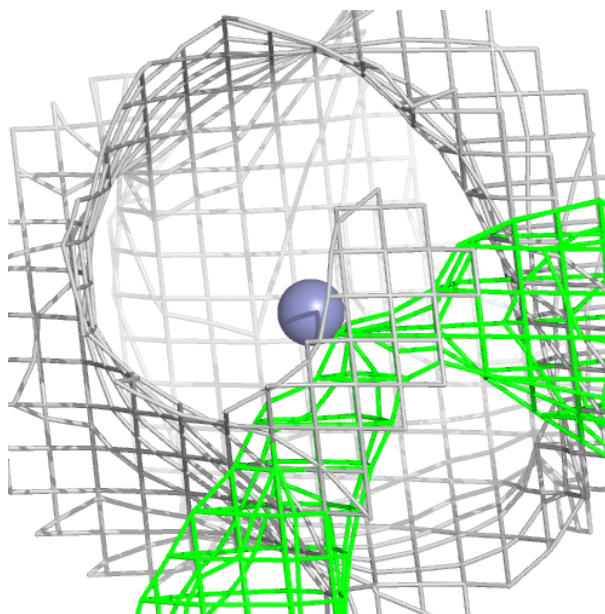
Electron density around ZN H 301:

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



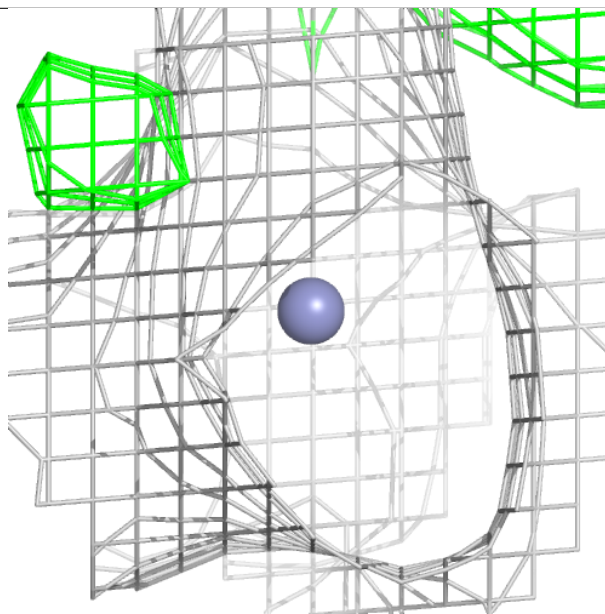
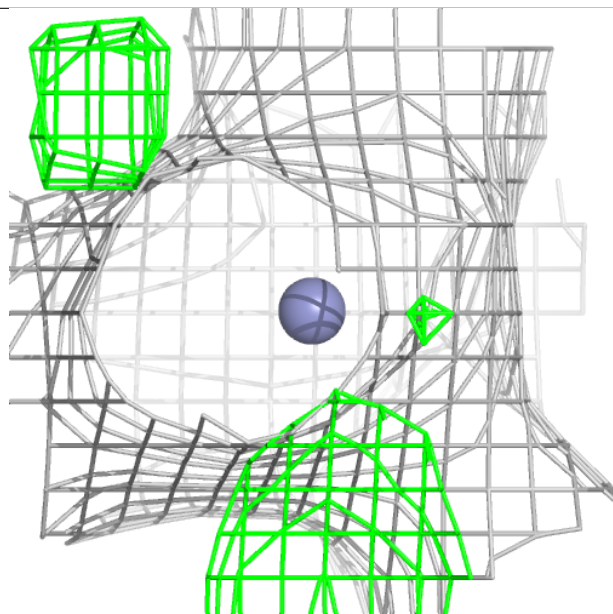
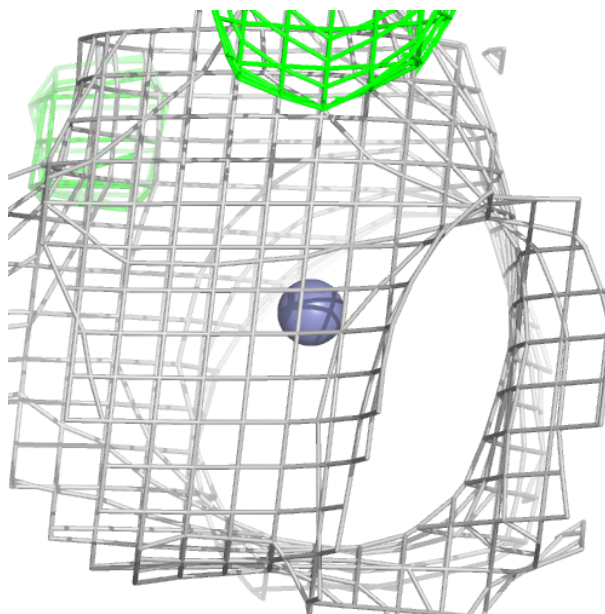
Electron density around ZN I 301:

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



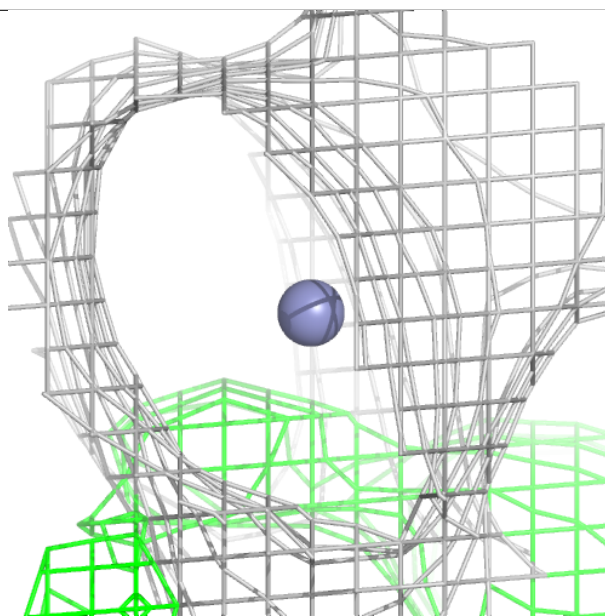
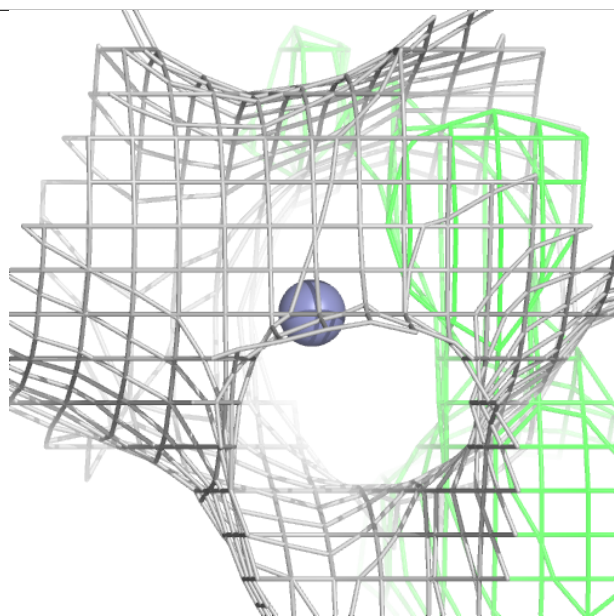
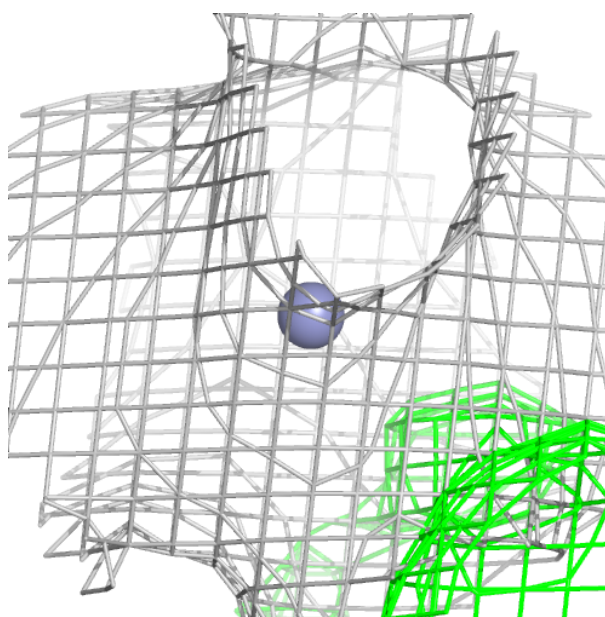
Electron density around ZN J 301:

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



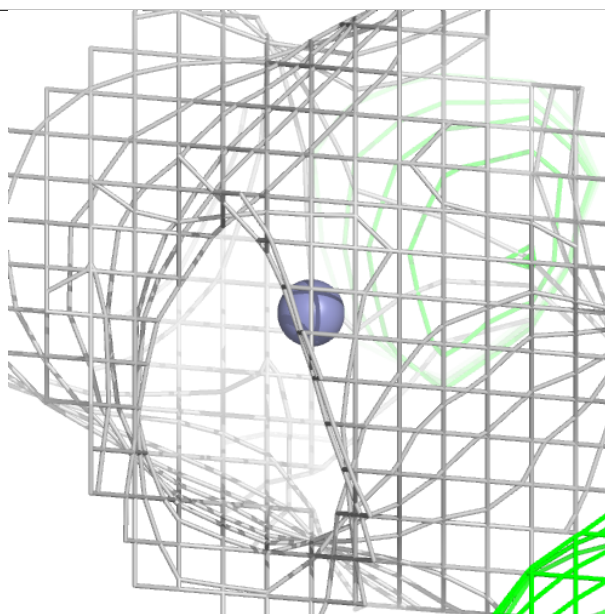
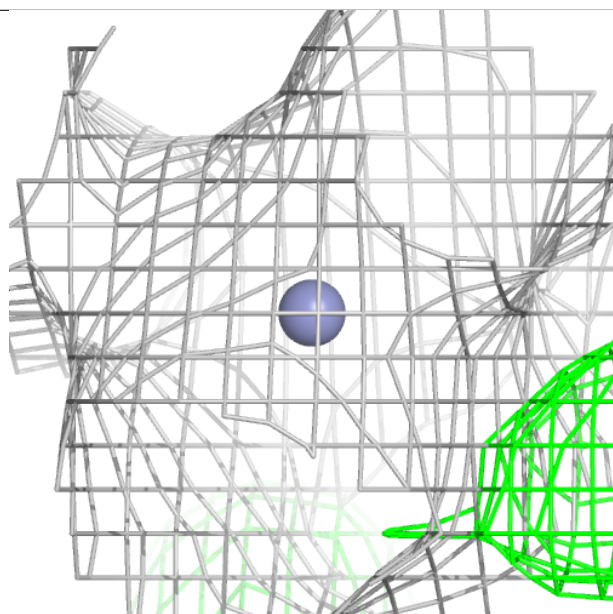
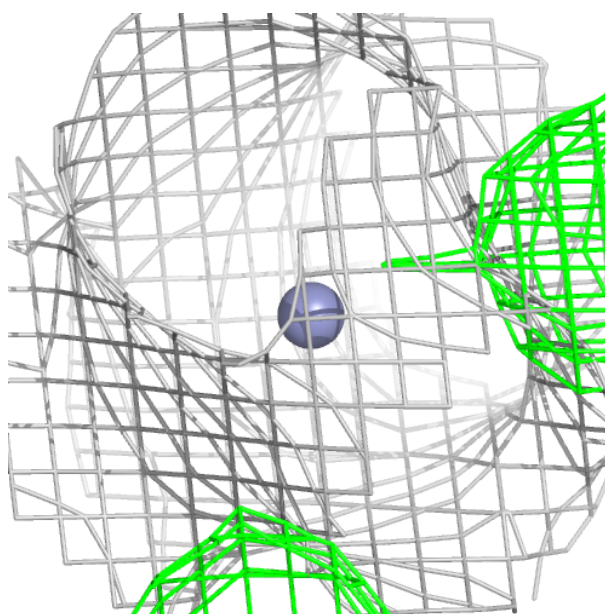
Electron density around ZN K 301:

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



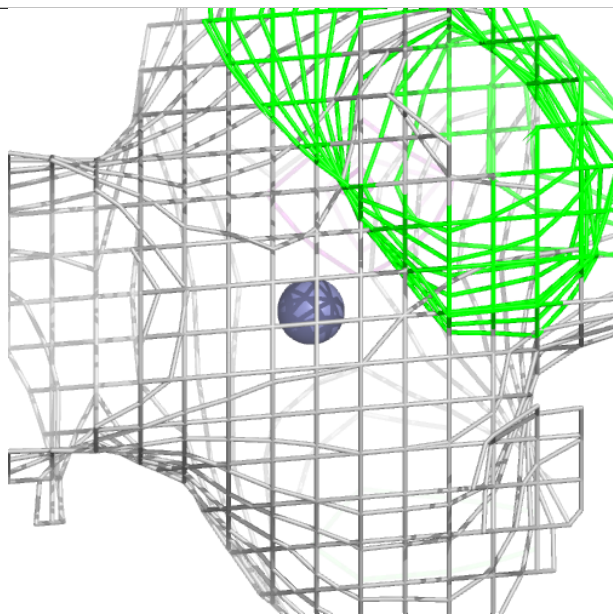
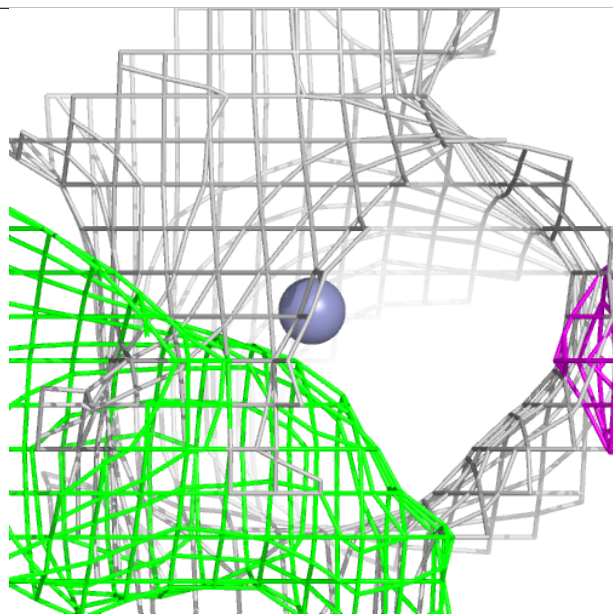
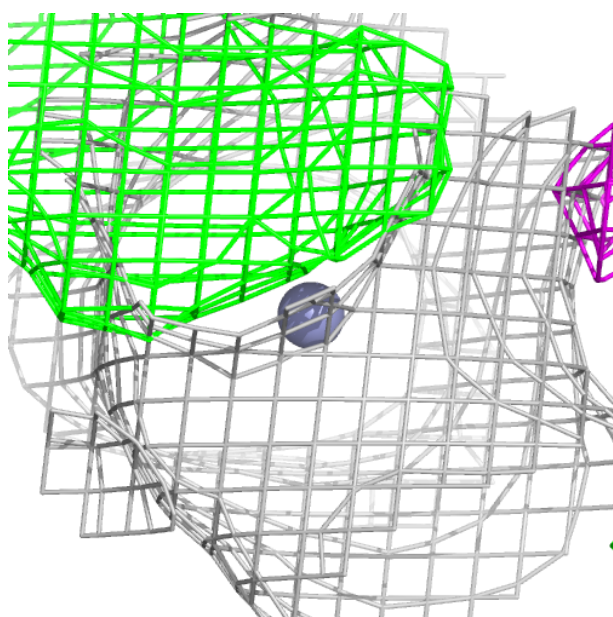
Electron density around ZN B 301:

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



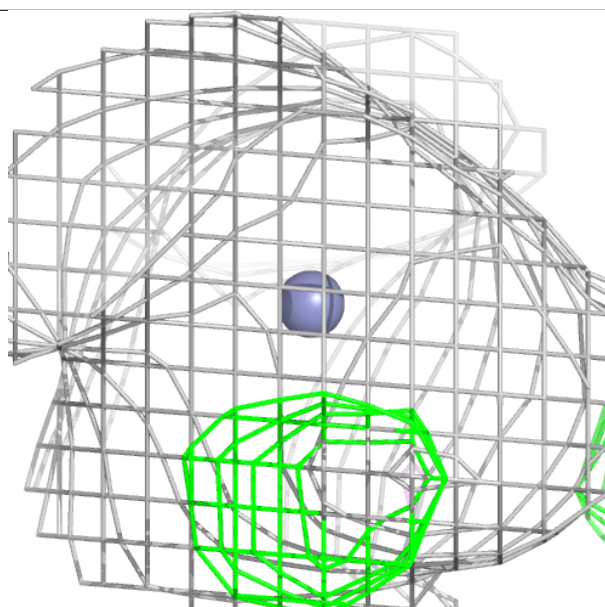
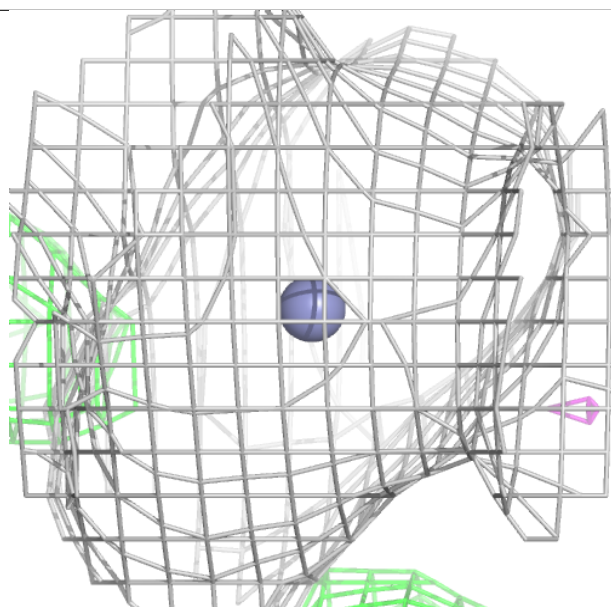
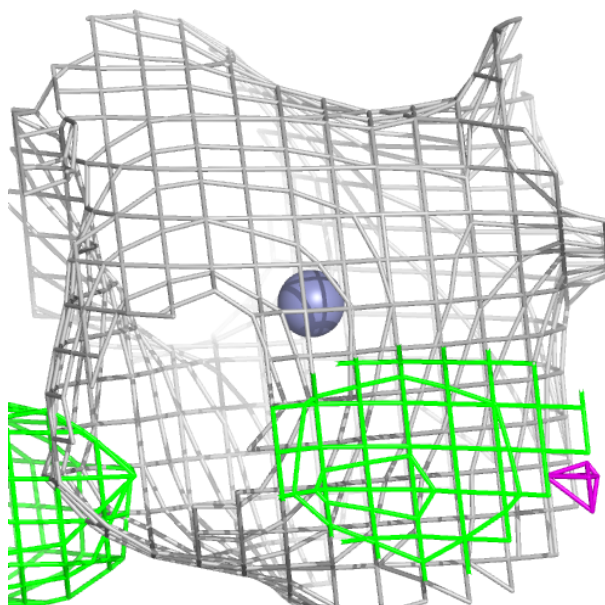
Electron density around ZN M 301:

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



Electron density around ZN N 301:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.