



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 05:26 PM UTC

PDB ID : 7VB3 / pdb_00007vb3
Title : Crystal structure of hydroxynitrile lyase from *Linum usitatissimum*
Authors : Zheng, D.; Nakabayashi, M.; Asano, Y.
Deposited on : 2021-08-30
Resolution : 1.48 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

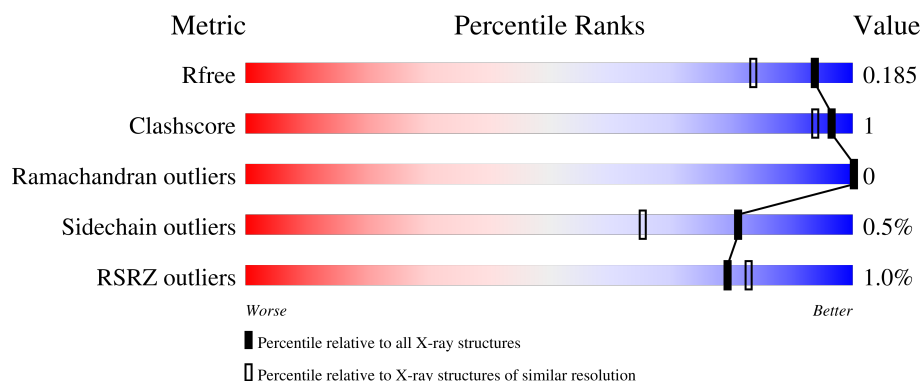
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.48 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	6779 (1.50-1.46)
Clashscore	190562	7025 (1.50-1.46)
Ramachandran outliers	187476	6917 (1.50-1.46)
Sidechain outliers	187428	6914 (1.50-1.46)
RSRZ outliers	180081	6781 (1.50-1.46)

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	443	% 87% 5% 8%
1	B	443	 91% • 7%
1	C	443	2% 87% 5% 8%
1	D	443	% 90% • 8%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 14555 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 Aliphatic (R)-hydroxynitrile lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	408	Total	C	N	O	S	0	6	0
			3142	1993	525	599	25			
1	B	414	Total	C	N	O	S	0	5	0
			3176	2011	528	612	25			
1	C	409	Total	C	N	O	S	0	1	0
			3115	1971	520	599	25			
1	D	408	Total	C	N	O	S	0	3	0
			3122	1980	521	597	24			

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	initiating methionine	UNP P93243
A	-19	GLY	-	expression tag	UNP P93243
A	-18	SER	-	expression tag	UNP P93243
A	-17	SER	-	expression tag	UNP P93243
A	-16	HIS	-	expression tag	UNP P93243
A	-15	HIS	-	expression tag	UNP P93243
A	-14	HIS	-	expression tag	UNP P93243
A	-13	HIS	-	expression tag	UNP P93243
A	-12	HIS	-	expression tag	UNP P93243
A	-11	HIS	-	expression tag	UNP P93243
A	-10	SER	-	expression tag	UNP P93243
A	-9	SER	-	expression tag	UNP P93243
A	-8	GLY	-	expression tag	UNP P93243
A	-7	LEU	-	expression tag	UNP P93243
A	-6	VAL	-	expression tag	UNP P93243
A	-5	PRO	-	expression tag	UNP P93243
A	-4	ARG	-	expression tag	UNP P93243
A	-3	GLY	-	expression tag	UNP P93243
A	-2	SER	-	expression tag	UNP P93243
A	-1	HIS	-	expression tag	UNP P93243
A	0	MET	-	expression tag	UNP P93243

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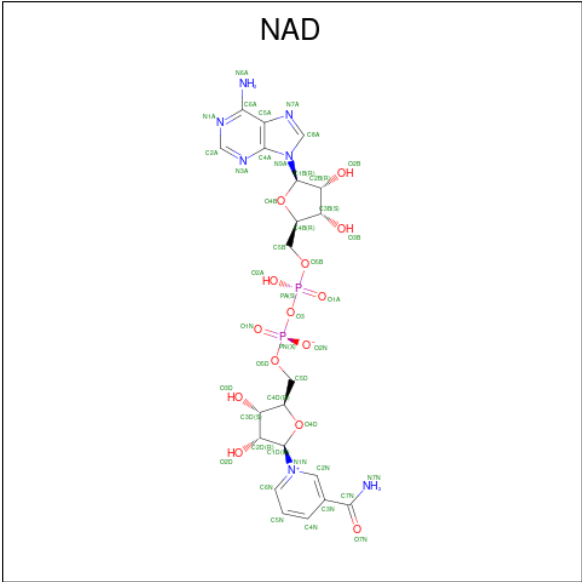
Chain	Residue	Modelled	Actual	Comment	Reference
A	117	THR	VAL	conflict	UNP P93243
B	-20	MET	-	initiating methionine	UNP P93243
B	-19	GLY	-	expression tag	UNP P93243
B	-18	SER	-	expression tag	UNP P93243
B	-17	SER	-	expression tag	UNP P93243
B	-16	HIS	-	expression tag	UNP P93243
B	-15	HIS	-	expression tag	UNP P93243
B	-14	HIS	-	expression tag	UNP P93243
B	-13	HIS	-	expression tag	UNP P93243
B	-12	HIS	-	expression tag	UNP P93243
B	-11	HIS	-	expression tag	UNP P93243
B	-10	SER	-	expression tag	UNP P93243
B	-9	SER	-	expression tag	UNP P93243
B	-8	GLY	-	expression tag	UNP P93243
B	-7	LEU	-	expression tag	UNP P93243
B	-6	VAL	-	expression tag	UNP P93243
B	-5	PRO	-	expression tag	UNP P93243
B	-4	ARG	-	expression tag	UNP P93243
B	-3	GLY	-	expression tag	UNP P93243
B	-2	SER	-	expression tag	UNP P93243
B	-1	HIS	-	expression tag	UNP P93243
B	0	MET	-	expression tag	UNP P93243
B	117	THR	VAL	conflict	UNP P93243
C	-20	MET	-	initiating methionine	UNP P93243
C	-19	GLY	-	expression tag	UNP P93243
C	-18	SER	-	expression tag	UNP P93243
C	-17	SER	-	expression tag	UNP P93243
C	-16	HIS	-	expression tag	UNP P93243
C	-15	HIS	-	expression tag	UNP P93243
C	-14	HIS	-	expression tag	UNP P93243
C	-13	HIS	-	expression tag	UNP P93243
C	-12	HIS	-	expression tag	UNP P93243
C	-11	HIS	-	expression tag	UNP P93243
C	-10	SER	-	expression tag	UNP P93243
C	-9	SER	-	expression tag	UNP P93243
C	-8	GLY	-	expression tag	UNP P93243
C	-7	LEU	-	expression tag	UNP P93243
C	-6	VAL	-	expression tag	UNP P93243
C	-5	PRO	-	expression tag	UNP P93243
C	-4	ARG	-	expression tag	UNP P93243
C	-3	GLY	-	expression tag	UNP P93243
C	-2	SER	-	expression tag	UNP P93243

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	HIS	-	expression tag	UNP P93243
C	0	MET	-	expression tag	UNP P93243
C	117	THR	VAL	conflict	UNP P93243
D	-20	MET	-	initiating methionine	UNP P93243
D	-19	GLY	-	expression tag	UNP P93243
D	-18	SER	-	expression tag	UNP P93243
D	-17	SER	-	expression tag	UNP P93243
D	-16	HIS	-	expression tag	UNP P93243
D	-15	HIS	-	expression tag	UNP P93243
D	-14	HIS	-	expression tag	UNP P93243
D	-13	HIS	-	expression tag	UNP P93243
D	-12	HIS	-	expression tag	UNP P93243
D	-11	HIS	-	expression tag	UNP P93243
D	-10	SER	-	expression tag	UNP P93243
D	-9	SER	-	expression tag	UNP P93243
D	-8	GLY	-	expression tag	UNP P93243
D	-7	LEU	-	expression tag	UNP P93243
D	-6	VAL	-	expression tag	UNP P93243
D	-5	PRO	-	expression tag	UNP P93243
D	-4	ARG	-	expression tag	UNP P93243
D	-3	GLY	-	expression tag	UNP P93243
D	-2	SER	-	expression tag	UNP P93243
D	-1	HIS	-	expression tag	UNP P93243
D	0	MET	-	expression tag	UNP P93243
D	117	THR	VAL	conflict	UNP P93243

- Molecule 2 is NICOTINAMIDE-ADENINE-DINUCLEOTIDE (CCD ID: NAD) (formula: $C_{21}H_{27}N_7O_{14}P_2$) (labeled as "Ligand of Interest" by depositor).

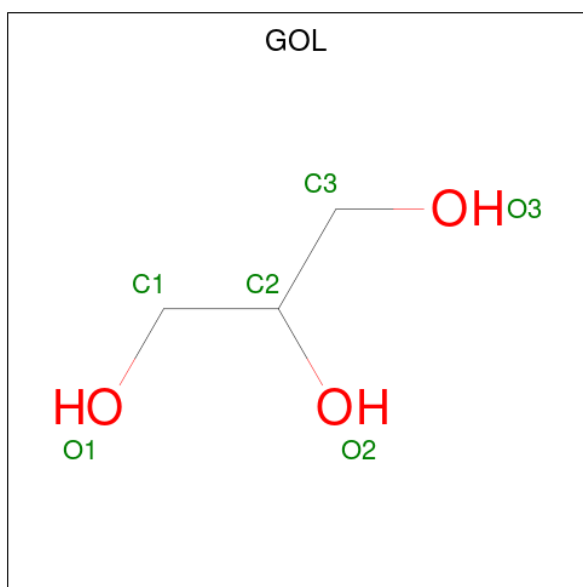


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	B	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	C	1	Total	C	N	O	P	0	0
			44	21	7	14	2		
2	D	1	Total	C	N	O	P	0	0
			44	21	7	14	2		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		
3	B	2	Total	Zn	0	0
			2	2		
3	C	2	Total	Zn	0	0
			2	2		
3	D	2	Total	Zn	0	0
			2	2		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

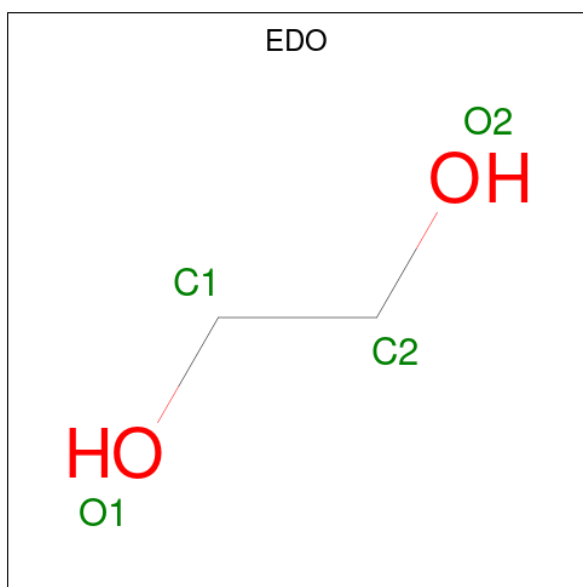
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Mg	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Mg	0	0
			1	1		
5	C	1	Total	Mg	0	0
			1	1		
5	D	1	Total	Mg	0	0
			1	1		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



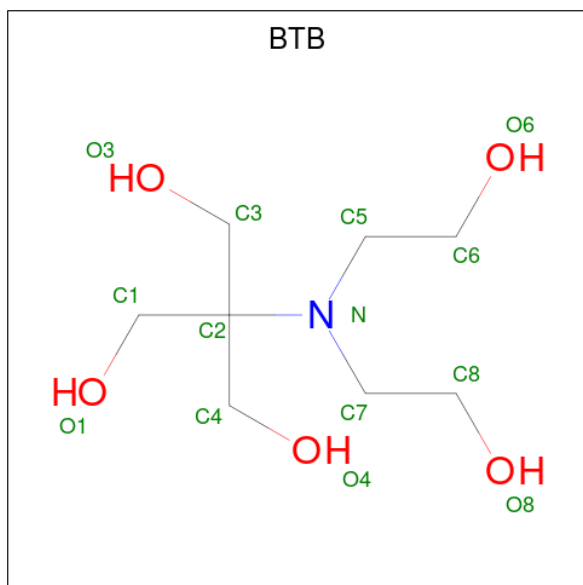
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	A	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	B	1	Total	C	O	0	0
			4	2	2		
6	C	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (CCD ID: BTB) (formula: C₈H₁₉NO₅).

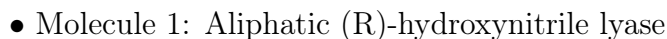
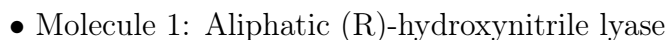
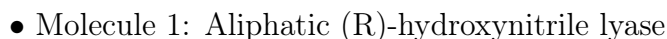


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	474	Total	O	0	0
			474	474		
8	B	478	Total	O	0	0
			478	478		
8	C	346	Total	O	0	0
			346	346		
8	D	370	Total	O	0	0
			370	370		

- Molecule 1: Aliphatic (R)-hydroxynitrile lyase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	94.12Å 52.18Å 168.51Å 90.00° 95.01° 90.00°	Depositor
Resolution (Å)	84.07 – 1.48 84.07 – 1.48	Depositor EDS
% Data completeness (in resolution range)	99.5 (84.07-1.48) 99.5 (84.07-1.48)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 1.48Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.156 , 0.183 0.157 , 0.185	Depositor DCC
R_{free} test set	13630 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	14.5	Xtriage
Anisotropy	0.081	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 33.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	14555	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.12% 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: MG, GOL, BTB, NAD, SNC, ZN, EDO

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	1.10	5/3206 (0.2%)	1.14	3/4330 (0.1%)
1	B	1.07	0/3237	1.18	6/4373 (0.1%)
1	C	1.08	4/3163 (0.1%)	1.25	6/4275 (0.1%)
1	D	1.04	0/3177	1.18	3/4294 (0.1%)
All	All	1.07	9/12783 (0.1%)	1.19	18/17272 (0.1%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	228	VAL	C-O	5.94	1.30	1.24
1	C	85	HIS	CE1-NE2	5.92	1.38	1.32
1	A	88	VAL	C-O	5.77	1.30	1.23
1	A	248	ASP	C-O	5.70	1.30	1.23
1	A	226	GLY	C-O	5.36	1.29	1.23
1	A	281	VAL	C-O	5.25	1.30	1.24
1	C	238	GLU	C-O	5.21	1.30	1.24
1	A	21	ALA	C-O	5.21	1.30	1.23
1	C	294	SER	CA-CB	-5.20	1.45	1.53

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	335	PHE	CA-CB-CG	-6.67	107.13	113.80
1	B	282	ARG	NE-CZ-NH2	5.87	124.48	119.20
1	B	100	ASN	CB-CA-C	-5.69	101.76	111.26
1	B	335	PHE	CA-CB-CG	-5.65	108.15	113.80
1	A	190	PRO	N-CA-CB	5.64	106.35	103.19
1	B	326	TYR	O-C-N	5.57	129.50	123.10
1	C	151	ASP	CA-CB-CG	5.43	118.03	112.60
1	D	335	PHE	CA-CB-CG	-5.31	108.49	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	191	PRO	N-CA-C	5.25	117.10	110.70
1	B	147	PHE	CA-CB-CG	5.22	119.02	113.80
1	C	335	PHE	CA-CB-CG	-5.21	108.59	113.80
1	D	147	PHE	CA-CB-CG	5.16	118.96	113.80
1	A	183	ASN	CA-C-O	-5.11	112.97	119.31
1	C	50	ASN	CA-CB-CG	5.08	117.68	112.60
1	D	191	PRO	N-CA-C	5.04	116.85	110.70
1	C	233	VAL	N-CA-C	-5.03	105.59	110.42
1	C	338	PHE	CA-CB-CG	-5.02	108.78	113.80
1	B	405	GLU	CB-CG-CD	5.00	121.10	112.60

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3142	0	3176	9	0
1	B	3176	0	3201	4	0
1	C	3115	0	3130	11	0
1	D	3122	0	3146	7	0
2	A	44	0	26	0	0
2	B	44	0	26	0	0
2	C	44	0	26	0	0
2	D	44	0	26	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	A	24	0	32	1	0
4	B	24	0	32	0	0
4	C	6	0	8	0	0
4	D	12	0	16	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	D	1	0	0	0	0
6	A	12	0	18	1	0
6	B	16	0	24	2	0
6	C	4	0	6	0	0
6	D	4	0	6	0	0
7	A	28	0	38	1	0
7	B	14	0	19	0	0
8	A	474	0	0	1	0
8	B	478	0	0	1	0
8	C	346	0	0	1	0
8	D	370	0	0	1	0
All	All	14555	0	12956	26	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 1.

All (26) 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:C:344:VAL:HG12	1:D:338:PHE:HZ	1.51	0.75
6:B:511:EDO:H12	8:B:855:HOH:O	1.90	0.72
1:C:344:VAL:HG12	1:D:338:PHE:CZ	2.34	0.59
1:C:327:GLU:HA	1:C:327:GLU:OE1	2.03	0.57
1:A:338:PHE:HZ	1:B:344:VAL:HG12	1.69	0.56
6:A:511:EDO:H12	8:A:609:HOH:O	2.09	0.52
1:C:329:ASP:OD1	1:D:335:PHE:HB3	2.09	0.52
1:C:91:ILE:HD12	1:C:105:ASP:HB2	1.93	0.50
1:D:61:SER:HB2	1:D:414:LYS:HD2	1.94	0.49
1:C:329:ASP:O	1:C:329:ASP:CG	2.51	0.49
1:A:344:VAL:HG12	1:B:338:PHE:HZ	1.77	0.49
1:C:323:GLU:HG2	1:D:339:LEU:HD22	1.97	0.47
1:A:380:TYR:CE2	1:A:382:ALA:HA	2.51	0.46
1:D:82:ILE:HD11	1:D:159:LEU:HD13	1.98	0.46
7:A:513:BTB:H41	7:A:513:BTB:H72	1.77	0.45
1:A:338:PHE:CZ	1:B:344:VAL:HG12	2.50	0.45
1:C:164:GLY:HA3	8:C:910:HOH:O	2.18	0.43
1:A:237:LYS:HD3	4:A:505:GOL:H12	2.00	0.43
1:C:271:LEU:O	1:C:272:PRO:C	2.59	0.43
1:A:368:PRO:HD2	1:C:29:LEU:HD12	2.00	0.43
1:C:392:ASP:HB3	1:C:395:GLU:HG2	2.00	0.43
1:A:91:ILE:HD12	1:A:105:ASP:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:379:GLY:O	6:B:512:EDO:H22	2.20	0.42
1:D:157:LYS:HE2	8:D:906:HOH:O	2.20	0.41
1:A:248:ASP:O	1:A:267:ASN:HA	2.20	0.41
1:A:61:SER:HB2	1:A:414:LYS:HD2	2.02	0.41

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	411/443 (93%)	402 (98%)	9 (2%)	0	100	100
1	B	416/443 (94%)	407 (98%)	9 (2%)	0	100	100
1	C	407/443 (92%)	397 (98%)	10 (2%)	0	100	100
1	D	408/443 (92%)	397 (97%)	11 (3%)	0	100	100
All	All	1642/1772 (93%)	1603 (98%)	39 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	343/367 (94%)	340 (99%)	3 (1%)	70	48
1	B	347/367 (95%)	345 (99%)	2 (1%)	78	60

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	339/367 (92%)	338 (100%)	1 (0%)	86	75
1	D	340/367 (93%)	339 (100%)	1 (0%)	86	75
All	All	1369/1468 (93%)	1362 (100%)	7 (0%)	81	65

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	81	LYS
1	A	182	LEU
1	A	376	LYS
1	B	81	LYS
1	B	172	MET
1	C	81	LYS
1	D	81	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	130	GLN
1	B	44	GLN
1	B	135	ASN
1	C	328	ASN
1	D	44	GLN
1	D	130	GLN
1	D	328	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	SNC	B	265	1	4,7,8	0.86	0	2,7,9	1.04	0
1	SNC	A	265	1	4,7,8	0.62	0	2,7,9	0.48	0
1	SNC	D	265	1	4,7,8	0.82	0	2,7,9	0.80	0
1	SNC	C	265	1	4,7,8	0.68	0	2,7,9	1.03	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	SNC	B	265	1	-	0/0/6/8	-
1	SNC	A	265	1	-	0/0/6/8	-
1	SNC	D	265	1	-	0/0/6/8	-
1	SNC	C	265	1	-	0/0/6/8	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 39 ligands modelled in this entry, 12 are monoatomic - leaving 27 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	$\# Z > 2$
7	BTB	B	513	-	13,13,13	1.47	3 (23%)	7,16,16	0.53	0
4	GOL	A	504	-	5,5,5	0.27	0	5,5,5	0.25	0
2	NAD	A	501	-	46,48,48	0.58	0	64,73,73	1.02	6 (9%)
6	EDO	A	509	-	3,3,3	0.53	0	2,2,2	0.35	0
4	GOL	B	506	-	5,5,5	0.13	0	5,5,5	0.37	0
2	NAD	C	501	-	46,48,48	0.77	1 (2%)	64,73,73	0.83	3 (4%)
4	GOL	B	507	-	5,5,5	0.20	0	5,5,5	0.44	0
4	GOL	D	505	-	5,5,5	0.09	0	5,5,5	0.27	0
7	BTB	A	513	-	13,13,13	1.43	3 (23%)	7,16,16	1.11	1 (14%)
4	GOL	D	504	-	5,5,5	0.18	0	5,5,5	0.13	0
6	EDO	D	507	-	3,3,3	0.24	0	2,2,2	0.37	0
4	GOL	A	507	-	5,5,5	0.17	0	5,5,5	0.45	0
4	GOL	B	505	-	5,5,5	0.15	0	5,5,5	0.44	0
6	EDO	B	511	-	3,3,3	0.37	0	2,2,2	0.09	0
4	GOL	A	505	-	5,5,5	0.47	0	5,5,5	0.75	0
7	BTB	A	512	-	13,13,13	1.43	3 (23%)	7,16,16	0.52	0
6	EDO	B	509	-	3,3,3	0.14	0	2,2,2	0.49	0
4	GOL	A	506	-	5,5,5	0.11	0	5,5,5	0.35	0
6	EDO	A	511	-	3,3,3	0.27	0	2,2,2	0.36	0
6	EDO	B	510	-	3,3,3	0.11	0	2,2,2	0.22	0
4	GOL	C	504	-	5,5,5	0.19	0	5,5,5	0.31	0
6	EDO	B	512	-	3,3,3	0.07	0	2,2,2	0.36	0
4	GOL	B	504	-	5,5,5	0.30	0	5,5,5	0.33	0
6	EDO	C	506	-	3,3,3	0.26	0	2,2,2	0.30	0
2	NAD	D	501	-	46,48,48	1.01	2 (4%)	64,73,73	0.82	1 (1%)
2	NAD	B	501	-	46,48,48	0.82	1 (2%)	64,73,73	0.72	1 (1%)
6	EDO	A	510	-	3,3,3	0.24	0	2,2,2	0.09	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	BTB	B	513	-	-	3/21/21/21	-
4	GOL	A	504	-	-	0/4/4/4	-
2	NAD	A	501	-	-	4/30/62/62	0/5/5/5
6	EDO	A	509	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	506	-	-	0/4/4/4	-
2	NAD	C	501	-	-	4/30/62/62	0/5/5/5
4	GOL	B	507	-	-	2/4/4/4	-
4	GOL	D	505	-	-	0/4/4/4	-
7	BTB	A	513	-	-	4/21/21/21	-
4	GOL	D	504	-	-	0/4/4/4	-
6	EDO	D	507	-	-	0/1/1/1	-
4	GOL	A	507	-	-	4/4/4/4	-
4	GOL	B	505	-	-	0/4/4/4	-
6	EDO	B	511	-	-	0/1/1/1	-
4	GOL	A	505	-	-	4/4/4/4	-
7	BTB	A	512	-	-	5/21/21/21	-
6	EDO	B	509	-	-	0/1/1/1	-
4	GOL	A	506	-	-	1/4/4/4	-
6	EDO	A	511	-	-	0/1/1/1	-
6	EDO	B	510	-	-	1/1/1/1	-
4	GOL	C	504	-	-	0/4/4/4	-
6	EDO	B	512	-	-	0/1/1/1	-
4	GOL	B	504	-	-	0/4/4/4	-
6	EDO	C	506	-	-	0/1/1/1	-
2	NAD	D	501	-	-	3/30/62/62	0/5/5/5
2	NAD	B	501	-	-	3/30/62/62	0/5/5/5
6	EDO	A	510	-	-	1/1/1/1	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	NAD	PA-O3	4.63	1.64	1.59
7	A	513	BTB	C2-N	3.46	1.55	1.48
7	A	512	BTB	C2-N	3.37	1.55	1.48
7	B	513	BTB	C5-N	3.06	1.52	1.48
2	D	501	NAD	O4D-C1D	2.90	1.44	1.40
2	C	501	NAD	O4D-C1D	2.82	1.44	1.40
7	B	513	BTB	C7-N	2.82	1.51	1.48
7	B	513	BTB	C2-N	2.79	1.54	1.48
2	B	501	NAD	PA-O3	2.58	1.62	1.59
7	A	512	BTB	C5-N	2.52	1.51	1.48
7	A	513	BTB	C7-N	2.49	1.51	1.48
7	A	512	BTB	C7-N	2.35	1.51	1.48
7	A	513	BTB	C5-N	2.07	1.50	1.48

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	NAD	C6N-N1N-C2N	-3.04	119.29	121.88
2	A	501	NAD	C5N-C4N-C3N	-2.80	117.61	120.36
2	C	501	NAD	C6N-N1N-C2N	-2.62	119.65	121.88
2	C	501	NAD	O4B-C1B-N9A	2.59	113.07	108.09
7	A	513	BTB	O3-C3-C2	2.45	117.16	111.40
2	A	501	NAD	C6N-N1N-C1D	2.36	124.36	119.73
2	A	501	NAD	O7N-C7N-N7N	2.30	125.94	122.62
2	C	501	NAD	O2N-PN-O3	2.29	113.46	107.27
2	D	501	NAD	O2D-C2D-C3D	-2.23	104.66	111.82
2	B	501	NAD	O2N-PN-O1N	2.22	122.79	112.44
2	A	501	NAD	O4B-C1B-N9A	2.14	112.20	108.09
2	A	501	NAD	O7N-C7N-C3N	-2.05	117.09	119.60

There are no chirality outliers.

All (39) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	NAD	O4D-C1D-N1N-C2N
2	A	501	NAD	O4D-C1D-N1N-C6N
2	A	501	NAD	C2D-C1D-N1N-C2N
2	A	501	NAD	C2D-C1D-N1N-C6N
2	B	501	NAD	O4D-C1D-N1N-C2N
2	B	501	NAD	O4D-C1D-N1N-C6N
2	B	501	NAD	C2D-C1D-N1N-C2N
2	C	501	NAD	O4D-C1D-N1N-C2N
2	C	501	NAD	O4D-C1D-N1N-C6N
2	C	501	NAD	C2D-C1D-N1N-C2N
2	D	501	NAD	O4D-C1D-N1N-C2N
2	D	501	NAD	O4D-C1D-N1N-C6N
2	D	501	NAD	C2D-C1D-N1N-C2N
4	A	505	GOL	C1-C2-C3-O3
4	A	505	GOL	O2-C2-C3-O3
4	B	507	GOL	C1-C2-C3-O3
7	A	512	BTB	C1-C2-C4-O4
7	A	512	BTB	C3-C2-C4-O4
7	A	512	BTB	N-C2-C4-O4
7	A	513	BTB	O1-C1-C2-C3
7	A	513	BTB	O1-C1-C2-C4
7	A	513	BTB	O1-C1-C2-N
7	B	513	BTB	C1-C2-C3-O3
7	B	513	BTB	C4-C2-C3-O3

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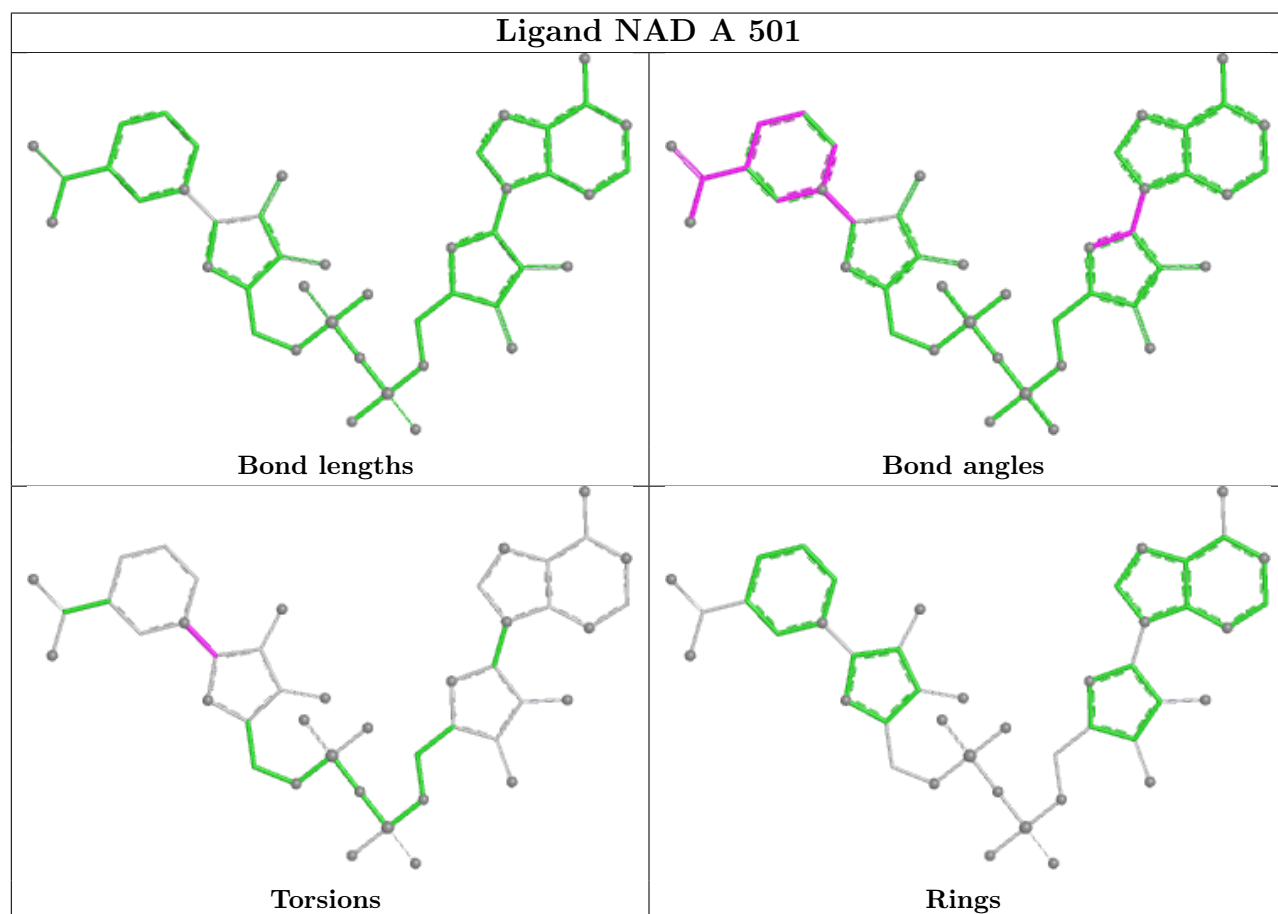
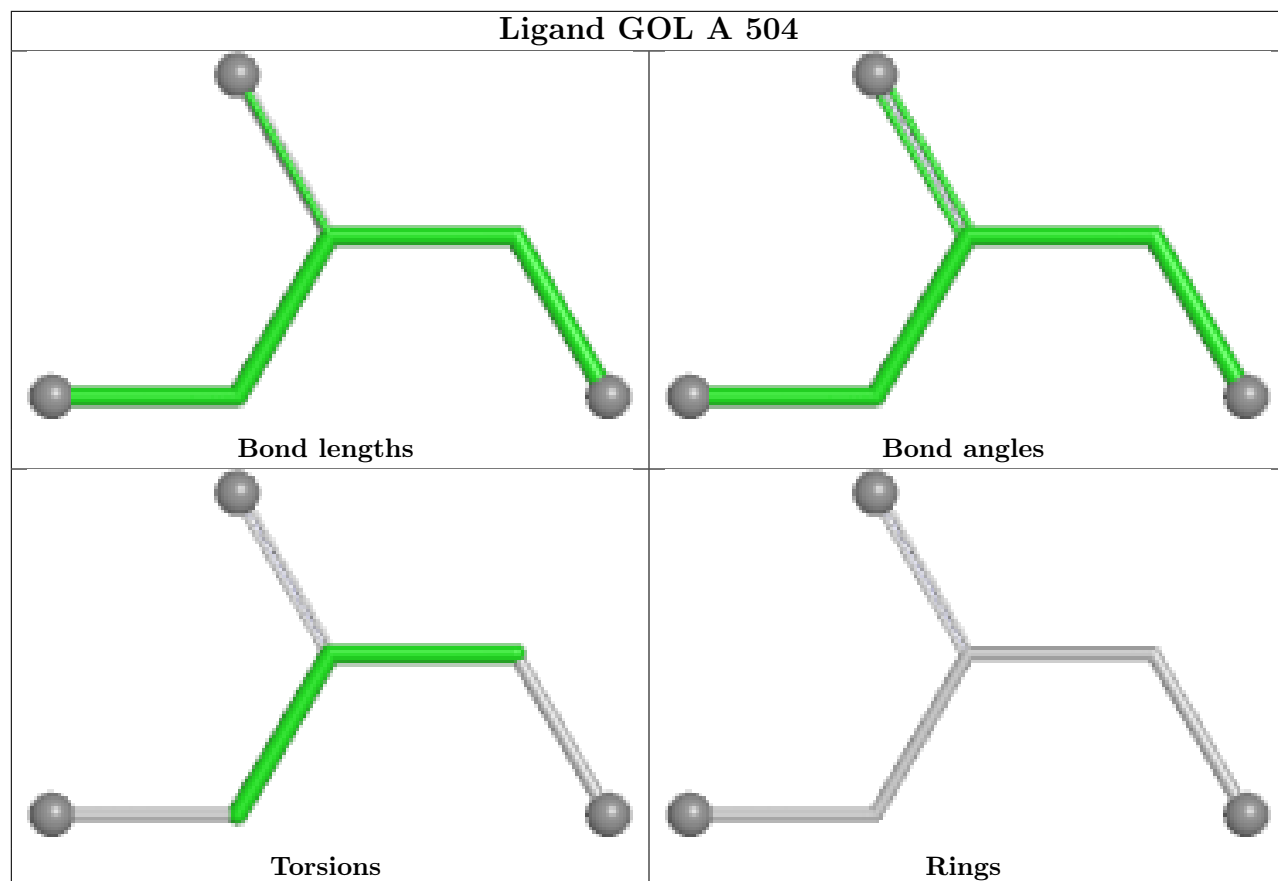
Mol	Chain	Res	Type	Atoms
7	B	513	BTB	N-C2-C3-O3
4	A	505	GOL	O1-C1-C2-C3
4	A	507	GOL	O1-C1-C2-C3
4	A	507	GOL	C1-C2-C3-O3
4	B	507	GOL	O2-C2-C3-O3
7	A	513	BTB	N-C5-C6-O6
4	A	505	GOL	O1-C1-C2-O2
4	A	507	GOL	O1-C1-C2-O2
4	A	506	GOL	O1-C1-C2-O2
2	C	501	NAD	C2D-C1D-N1N-C6N
6	A	510	EDO	O1-C1-C2-O2
4	A	507	GOL	O2-C2-C3-O3
6	B	510	EDO	O1-C1-C2-O2
7	A	512	BTB	N-C5-C6-O6
7	A	512	BTB	O1-C1-C2-C4

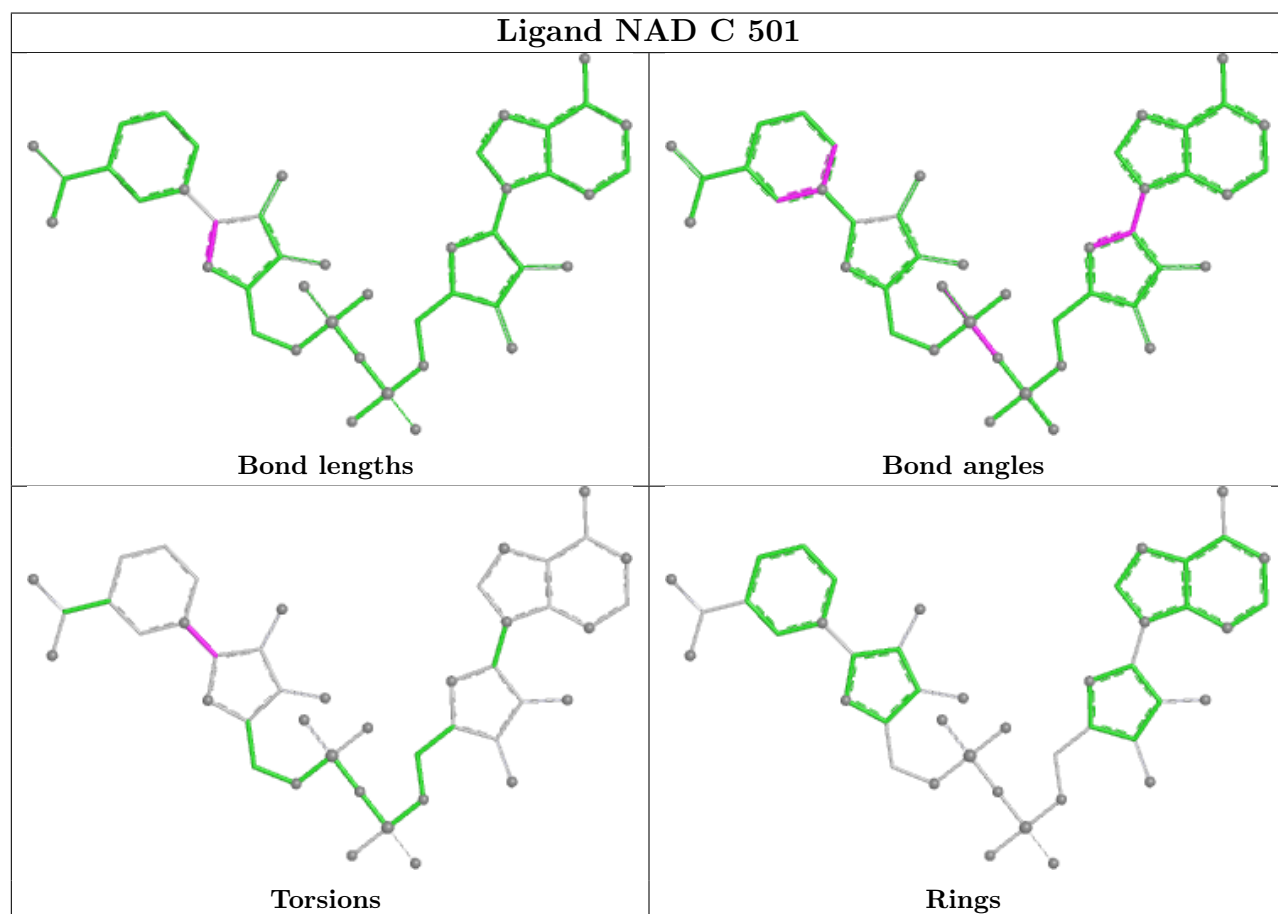
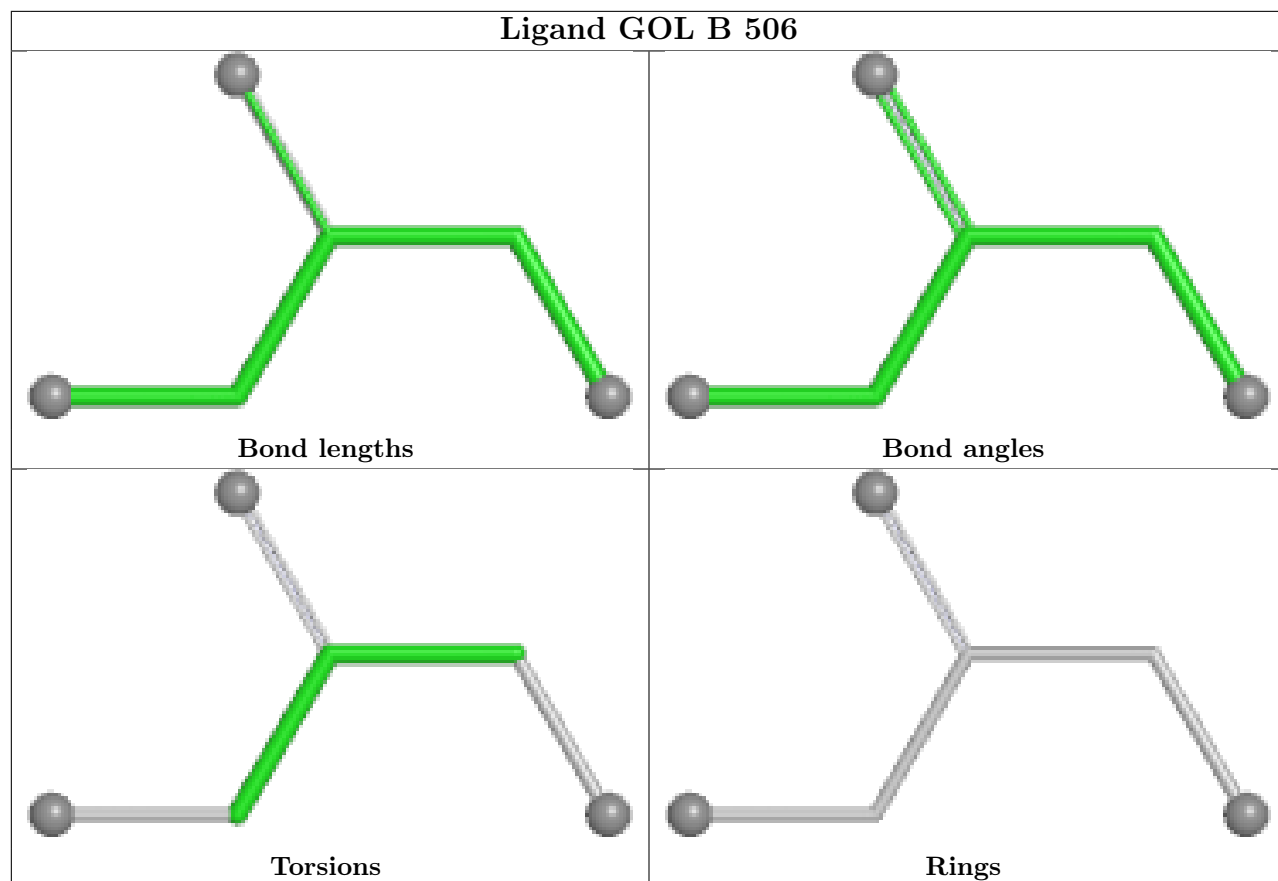
There are no ring outliers.

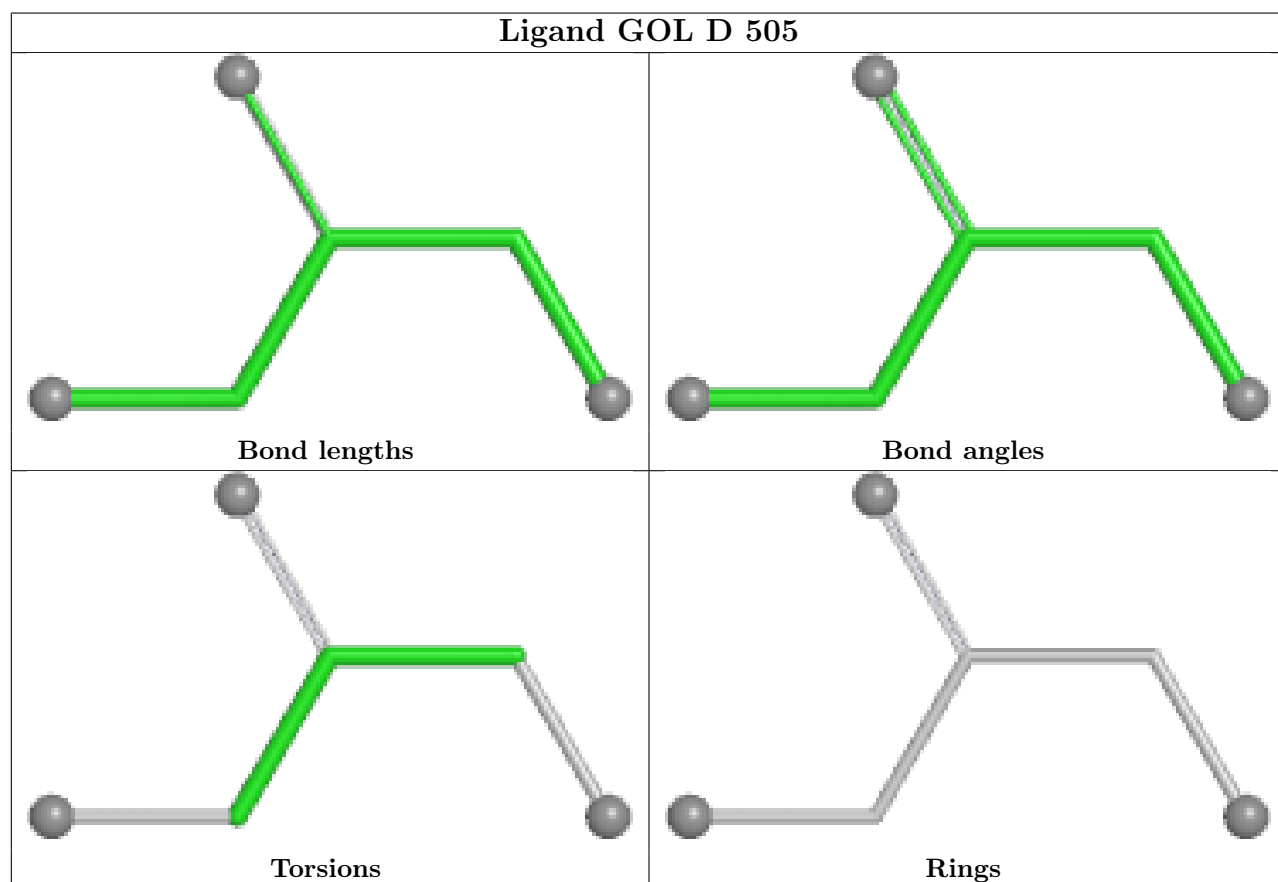
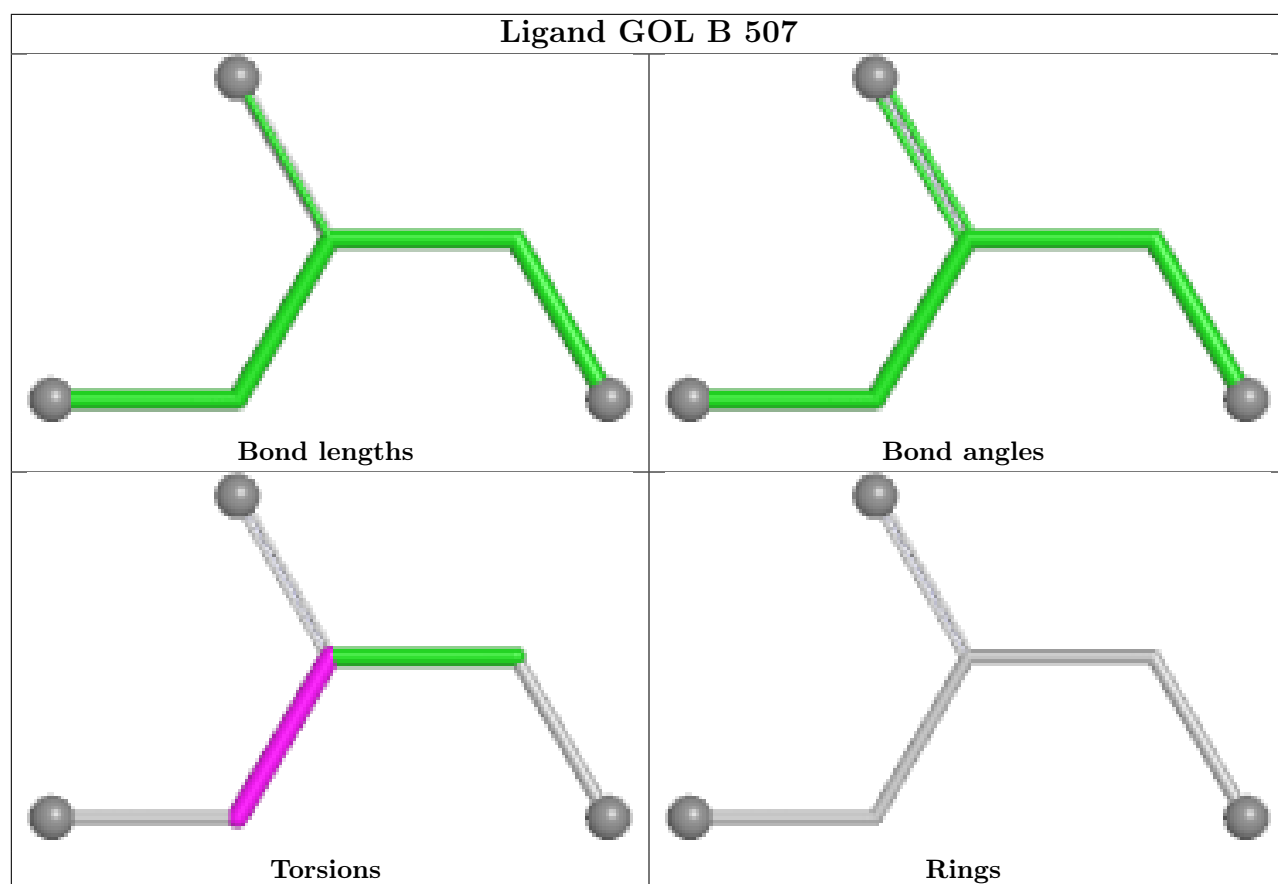
5 monomers are involved in 5 short contacts:

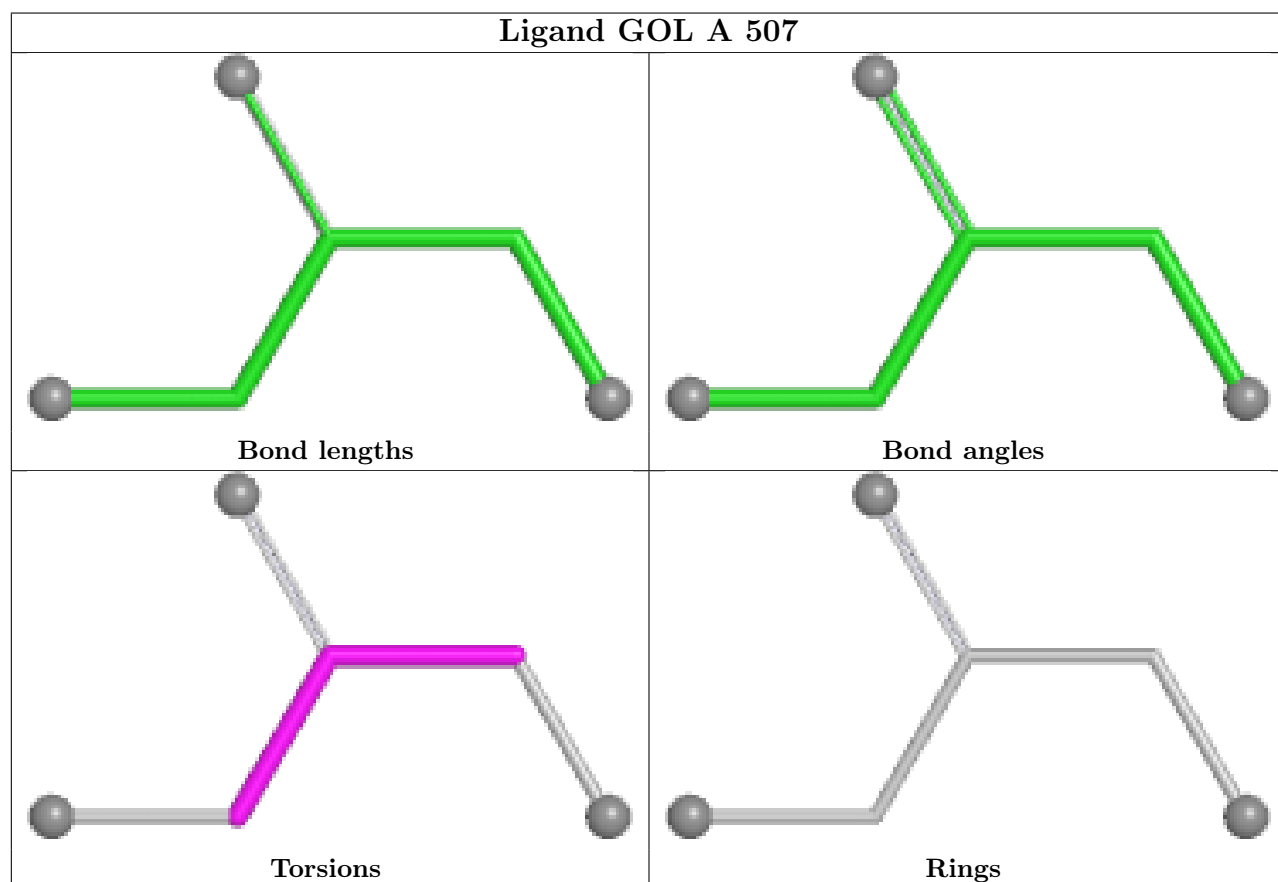
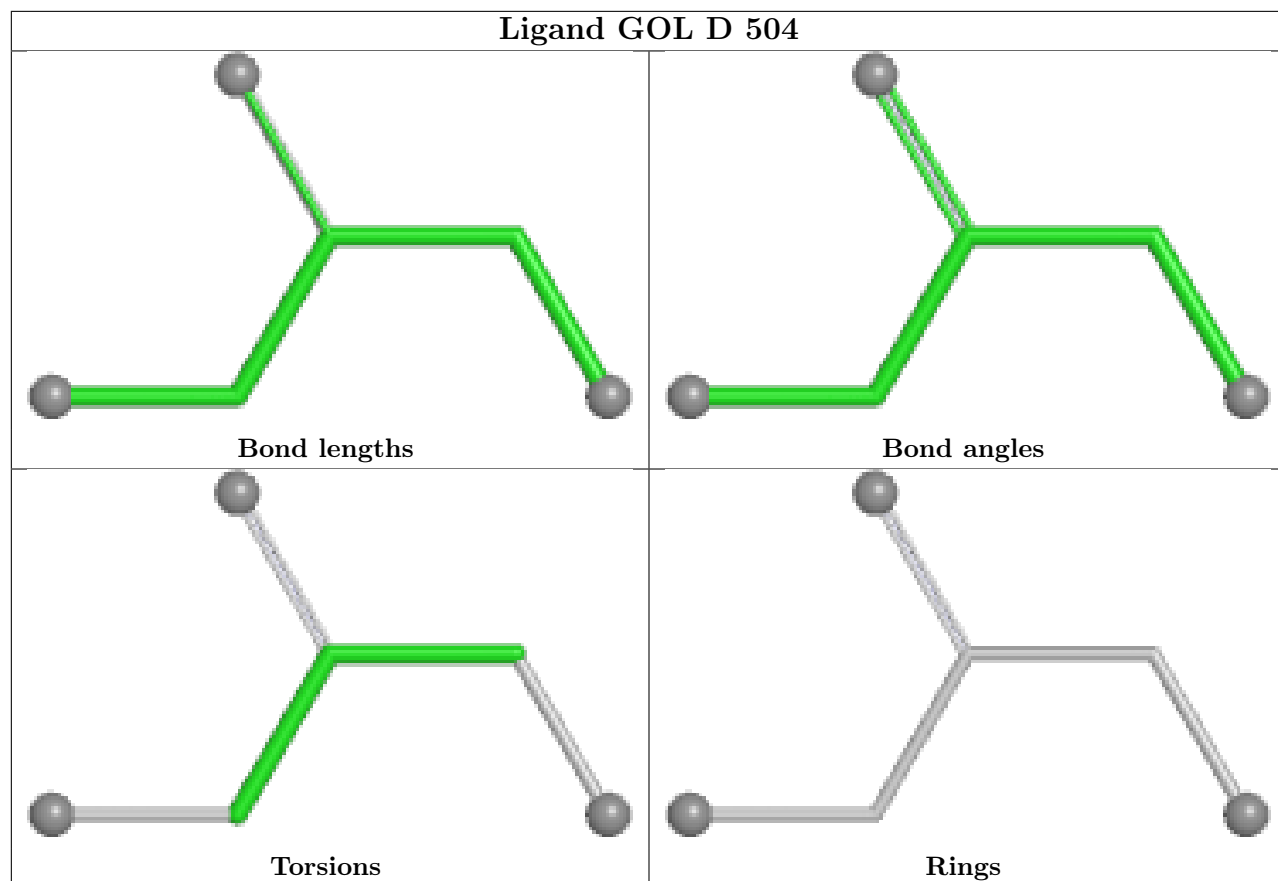
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	513	BTB	1	0
6	B	511	EDO	1	0
4	A	505	GOL	1	0
6	A	511	EDO	1	0
6	B	512	EDO	1	0

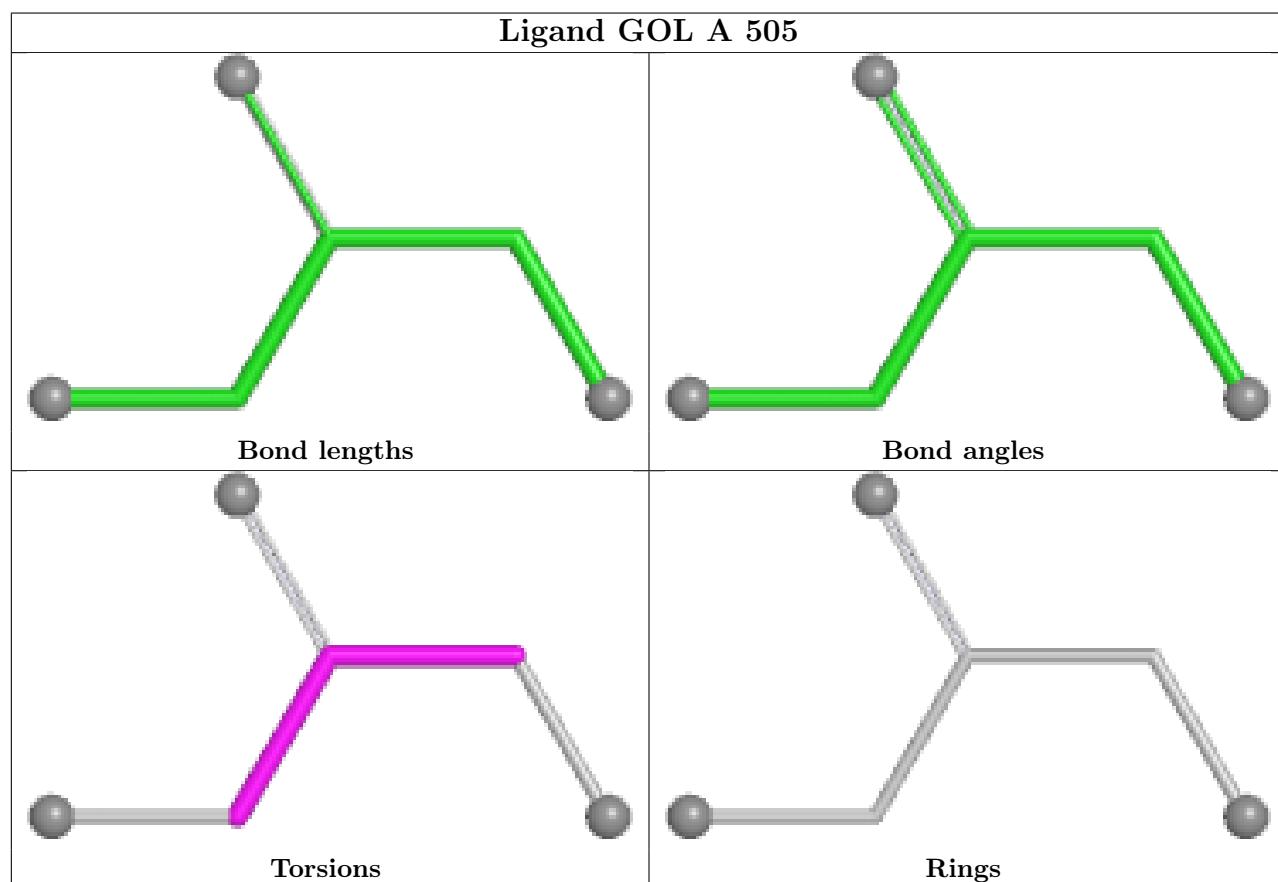
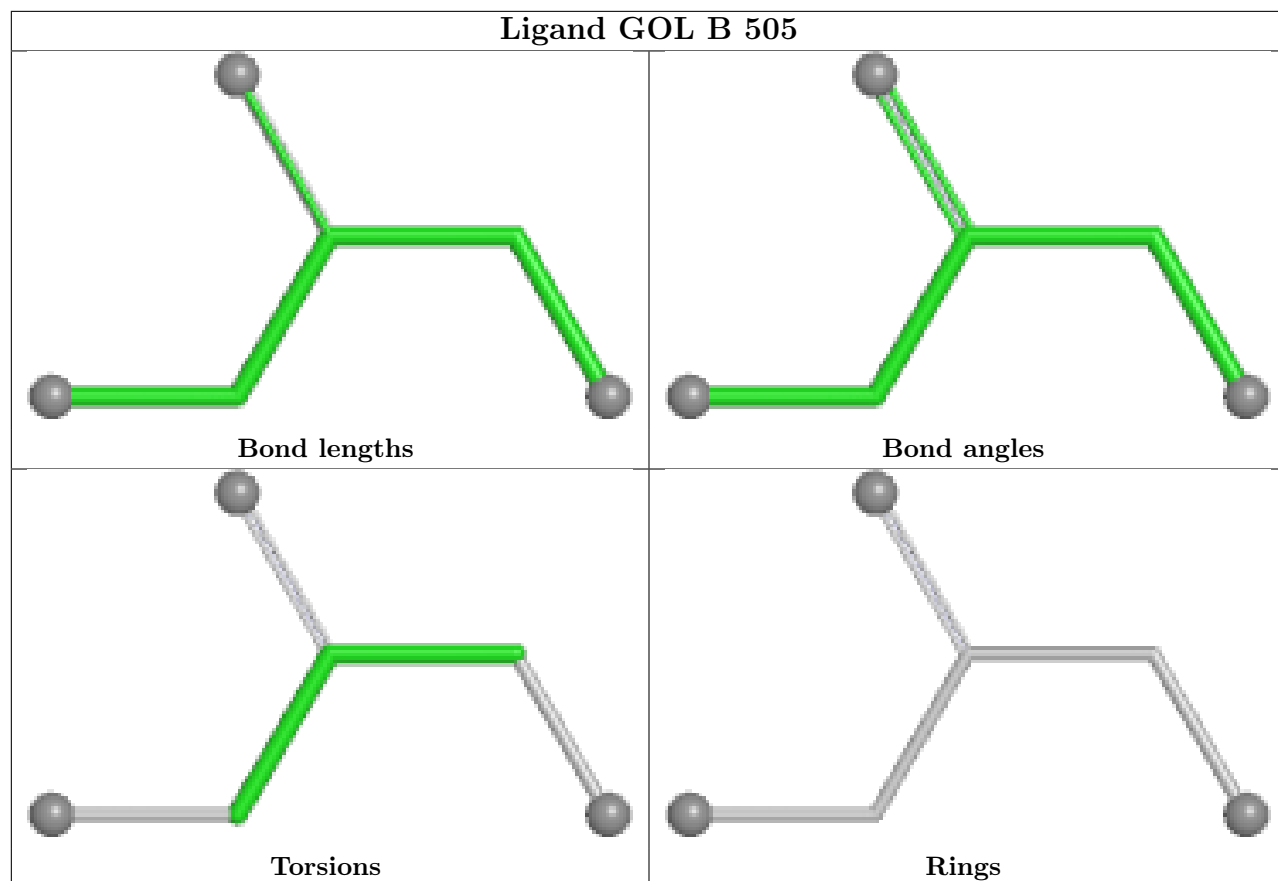
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

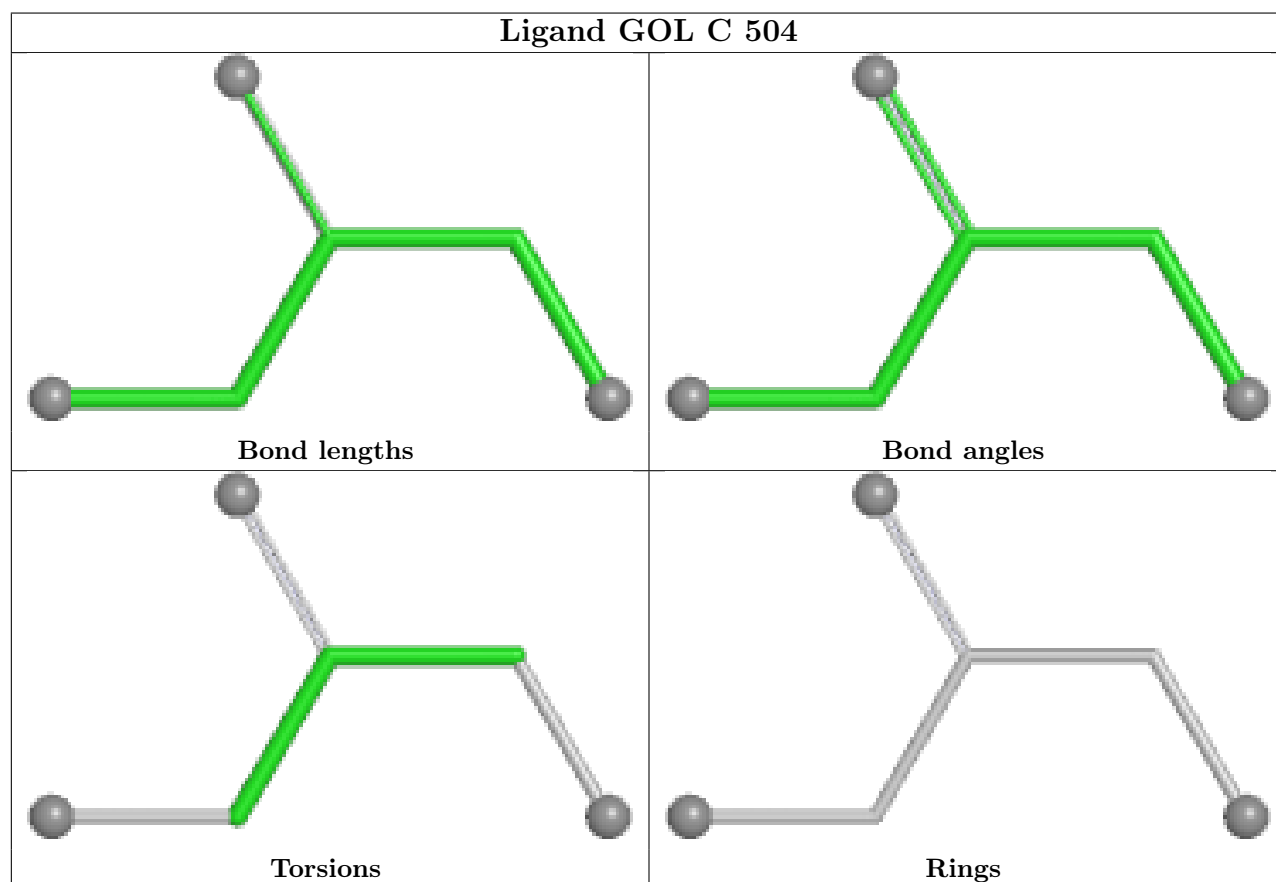
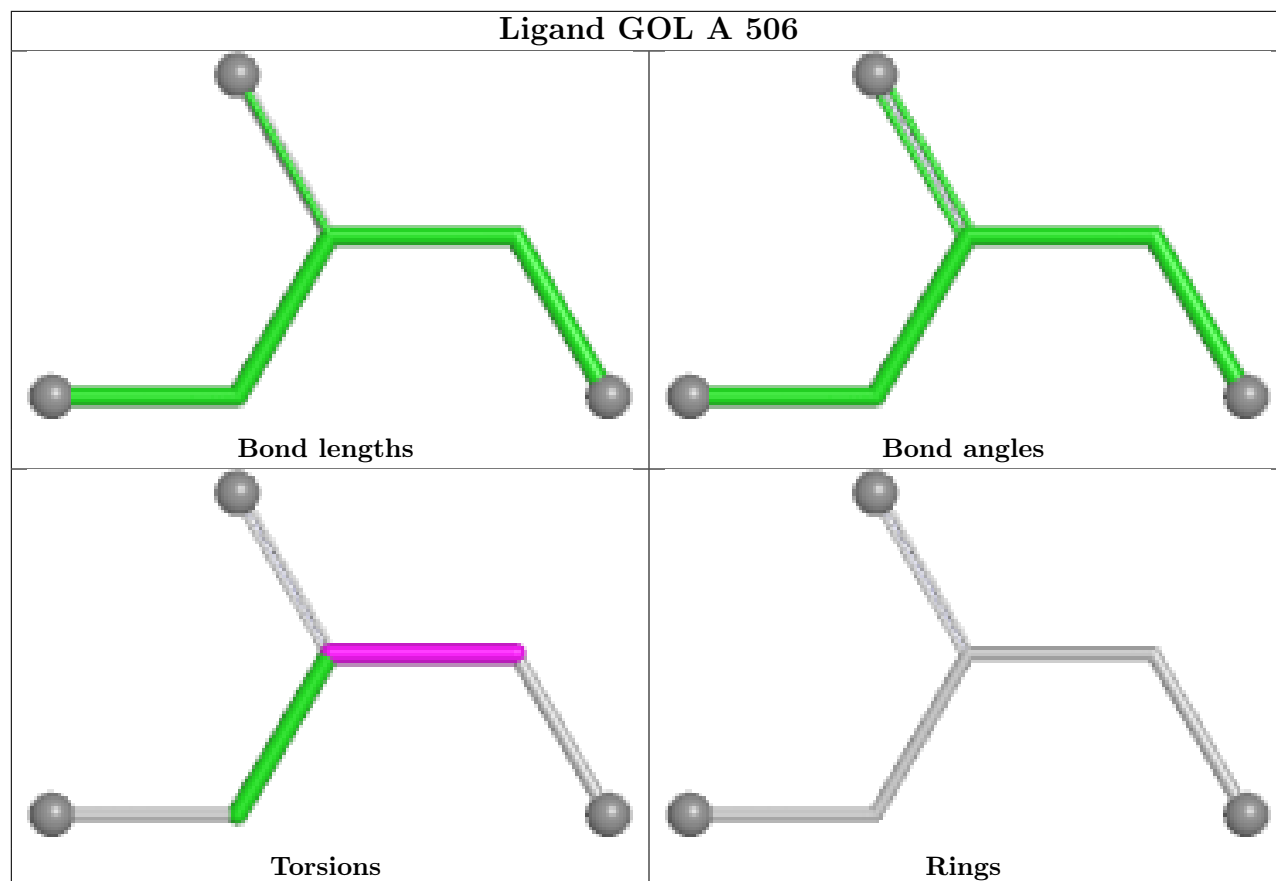


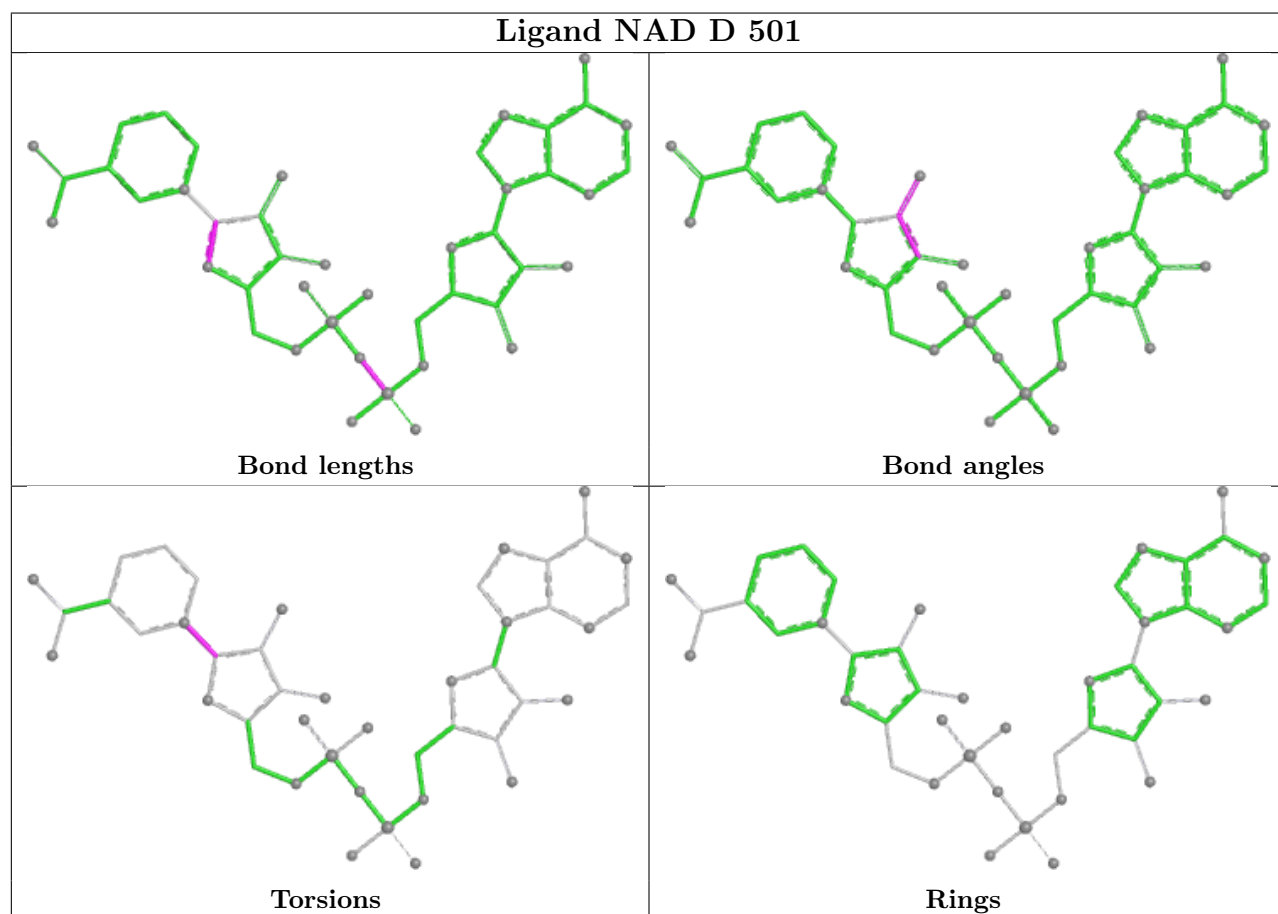
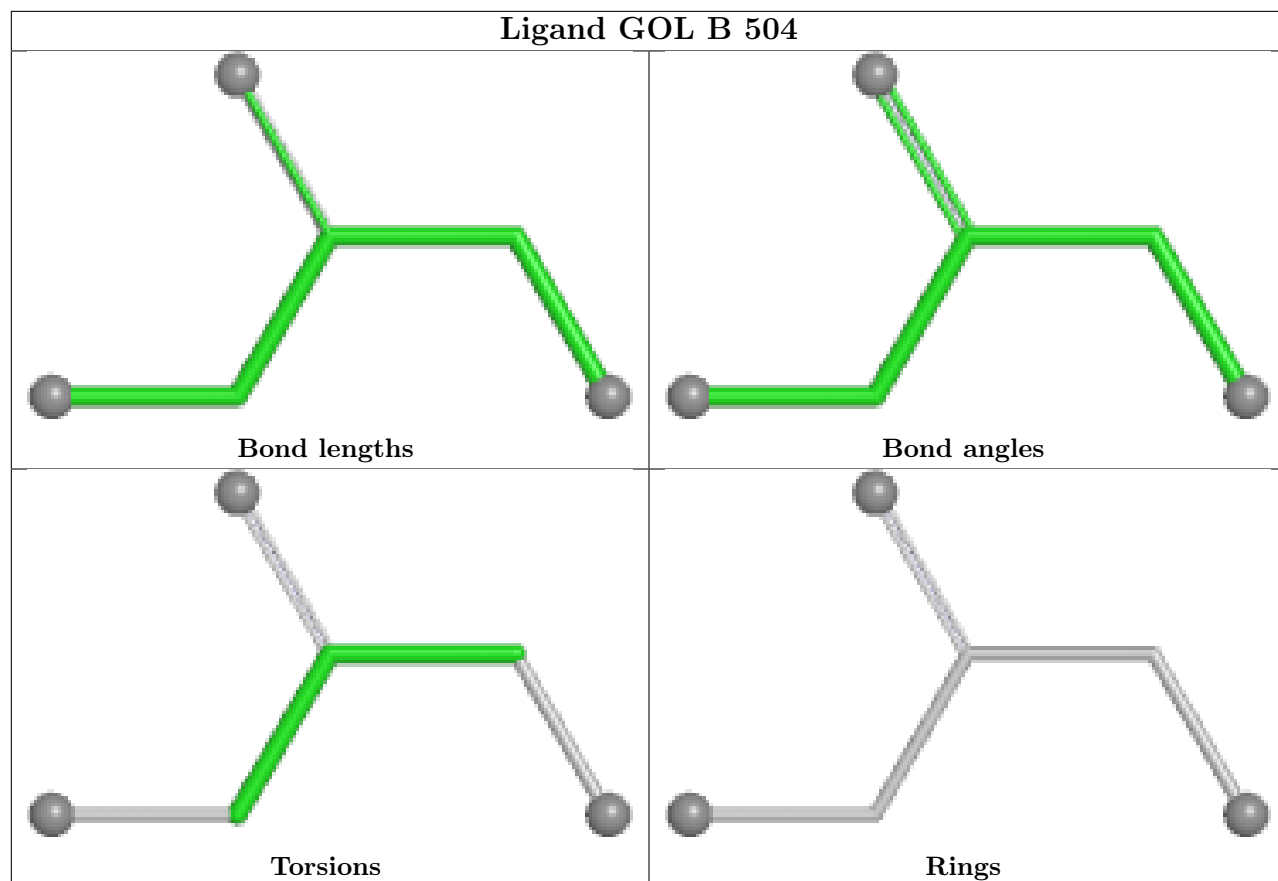


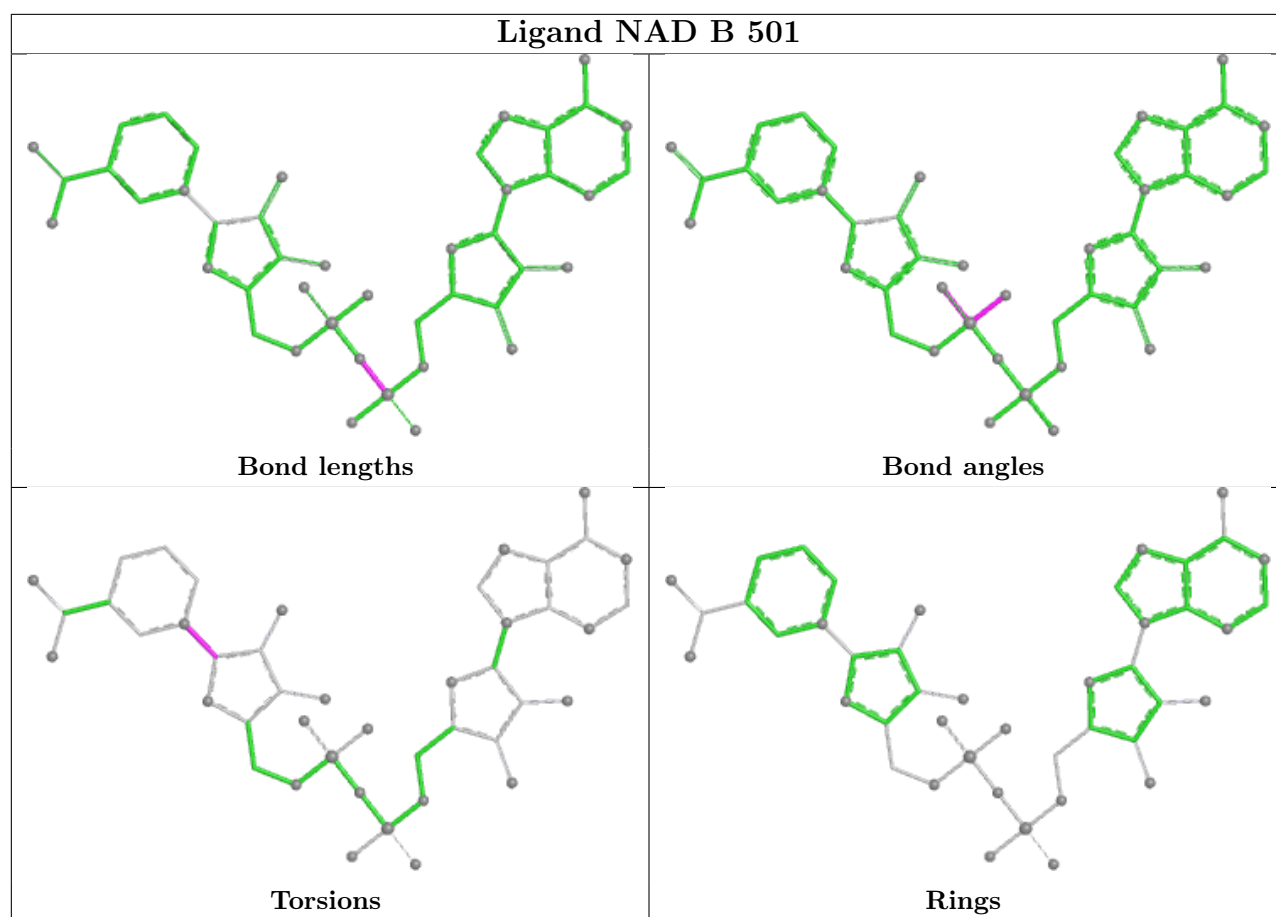












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	407/443 (91%)	-0.39	3 (0%) 84 87	7, 13, 22, 54	6 (1%)
1	B	413/443 (93%)	-0.34	2 (0%) 87 90	8, 13, 26, 49	5 (1%)
1	C	408/443 (92%)	0.24	7 (1%) 69 73	13, 19, 33, 57	1 (0%)
1	D	407/443 (91%)	0.09	4 (0%) 79 83	9, 18, 30, 52	3 (0%)
All	All	1635/1772 (92%)	-0.10	16 (0%) 79 83	7, 16, 29, 57	15 (0%)

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	273	GLU	3.6
1	D	332[A]	PHE	3.1
1	A	332[A]	PHE	3.0
1	B	9	ALA	3.0
1	C	383	ALA	2.8
1	D	15	GLY	2.8
1	A	152	SER	2.5
1	A	153	ASP	2.5
1	D	273	GLU	2.5
1	C	150	ILE	2.5
1	C	369	VAL	2.3
1	D	152	SER	2.3
1	C	365	ALA	2.1
1	C	152	SER	2.1
1	C	14	ASN	2.0
1	C	153	ASP	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 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
1	SNC	C	265	8/9	0.90	0.12	18,20,24,32	0
1	SNC	B	265	8/9	0.95	0.07	13,15,26,33	0
1	SNC	D	265	8/9	0.96	0.07	14,17,19,24	0
1	SNC	A	265	8/9	0.97	0.07	12,13,18,29	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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
4	GOL	B	507	6/6	0.75	0.17	34,40,43,46	0
4	GOL	A	506	6/6	0.78	0.15	41,43,48,50	0
7	BTB	A	512	14/14	0.78	0.14	24,31,34,34	0
6	EDO	B	509	4/4	0.79	0.15	28,31,36,38	0
7	BTB	A	513	14/14	0.79	0.17	37,51,66,68	0
4	GOL	A	507	6/6	0.80	0.14	34,37,40,43	0
6	EDO	C	506	4/4	0.82	0.16	32,35,39,40	0
6	EDO	D	507	4/4	0.82	0.14	36,37,40,44	0
4	GOL	D	505	6/6	0.83	0.14	34,43,45,46	0
4	GOL	A	505	6/6	0.83	0.16	22,32,37,42	0
6	EDO	A	510	4/4	0.86	0.13	33,34,36,40	0
6	EDO	B	512	4/4	0.86	0.13	36,37,37,45	0
4	GOL	B	506	6/6	0.87	0.12	25,36,38,39	0
6	EDO	A	511	4/4	0.87	0.12	35,35,36,42	0
6	EDO	B	511	4/4	0.89	0.14	22,30,34,35	0
6	EDO	B	510	4/4	0.89	0.12	22,29,32,35	0
7	BTB	B	513	14/14	0.89	0.11	22,29,34,34	0
6	EDO	A	509	4/4	0.91	0.12	19,25,26,32	0
4	GOL	B	505	6/6	0.91	0.10	27,28,29,29	0
4	GOL	C	504	6/6	0.94	0.07	22,25,26,27	0
4	GOL	D	504	6/6	0.94	0.08	23,26,28,30	0
4	GOL	B	504	6/6	0.94	0.08	16,19,20,22	0

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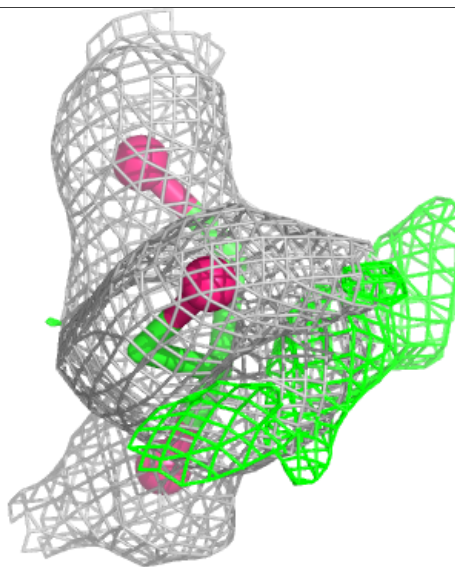
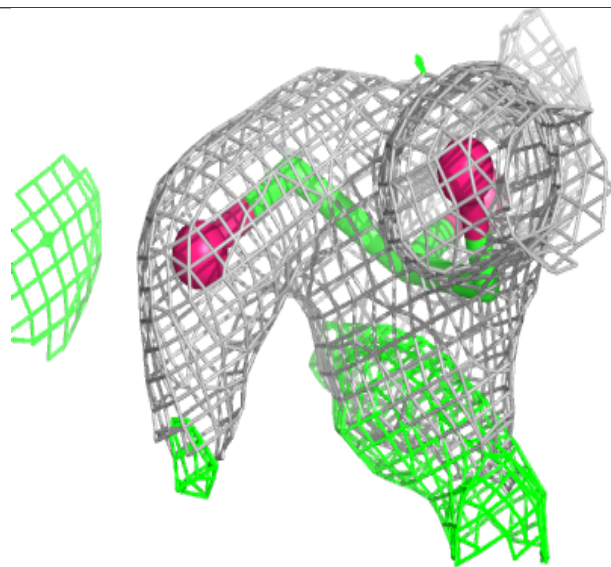
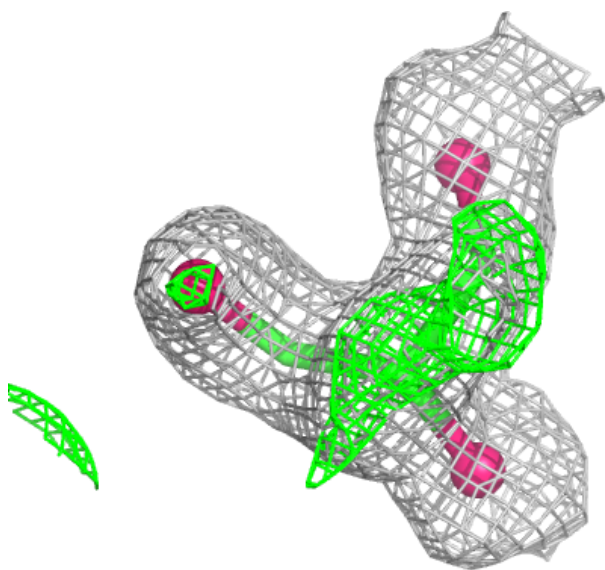
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MG	D	506	1/1	0.96	0.04	16,16,16,16	0
4	GOL	A	504	6/6	0.97	0.06	19,21,21,22	0
2	NAD	C	501	44/44	0.98	0.05	12,15,18,19	0
2	NAD	D	501	44/44	0.98	0.05	12,15,17,18	0
3	ZN	D	502	1/1	0.98	0.04	14,14,14,14	1
5	MG	C	505	1/1	0.99	0.04	17,17,17,17	0
3	ZN	D	503	1/1	0.99	0.02	14,14,14,14	1
2	NAD	B	501	44/44	0.99	0.04	8,11,13,15	0
3	ZN	C	503	1/1	0.99	0.03	13,13,13,13	1
2	NAD	A	501	44/44	0.99	0.04	8,10,12,15	0
5	MG	A	508	1/1	0.99	0.07	9,9,9,9	0
5	MG	B	508	1/1	0.99	0.05	10,10,10,10	0
3	ZN	A	502	1/1	1.00	0.01	12,12,12,12	0
3	ZN	A	503	1/1	1.00	0.01	10,10,10,10	1
3	ZN	B	502	1/1	1.00	0.01	12,12,12,12	0
3	ZN	B	503	1/1	1.00	0.03	11,11,11,11	0
3	ZN	C	502	1/1	1.00	0.01	16,16,16,16	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.

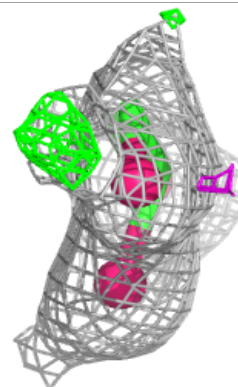
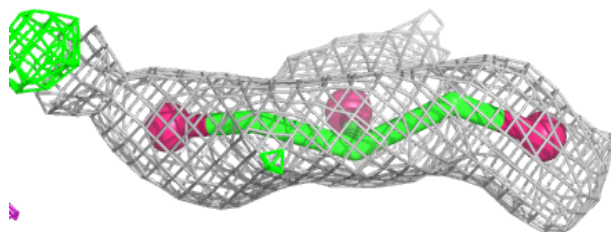
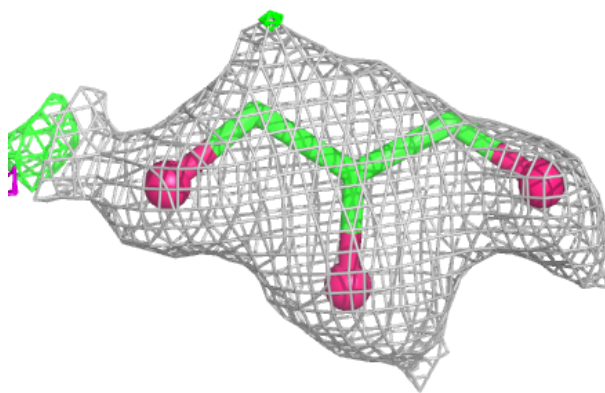
Electron density around GOL B 507:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



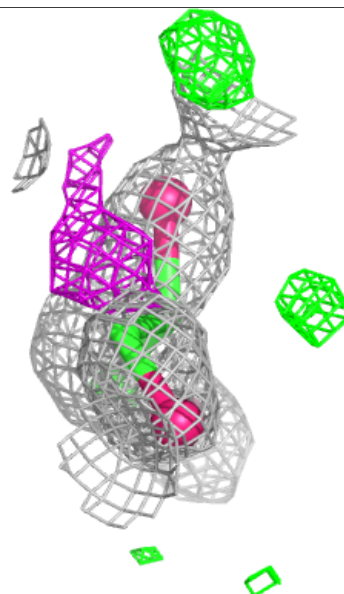
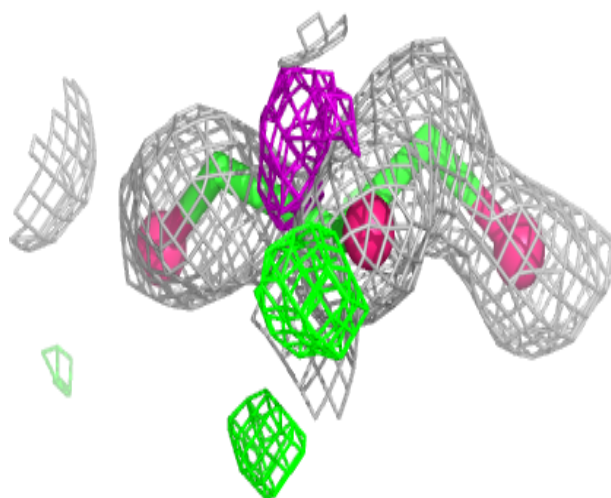
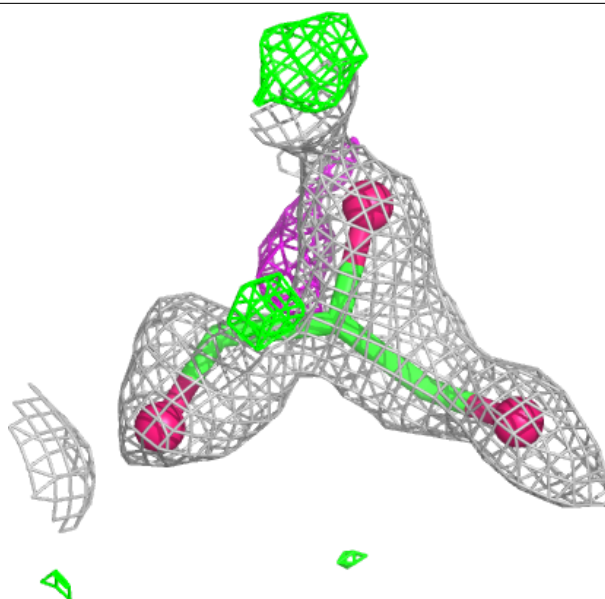
Electron density around GOL A 506:

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and green (positive)



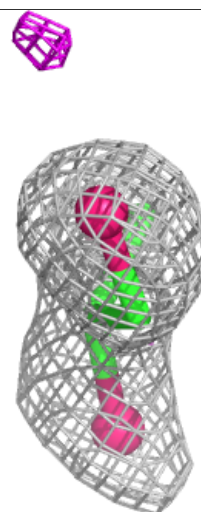
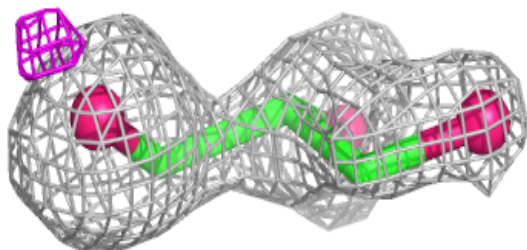
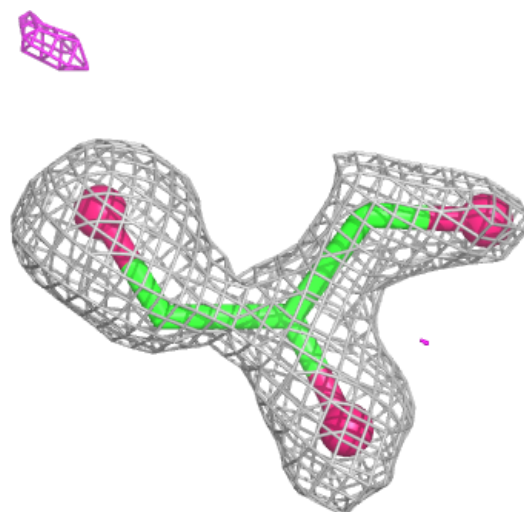
Electron density around GOL A 507:

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and green (positive)



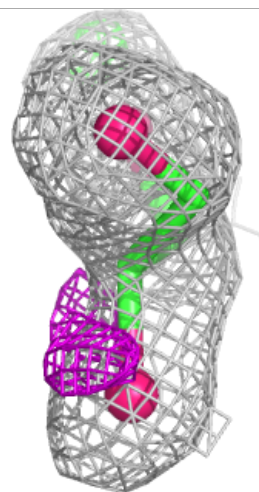
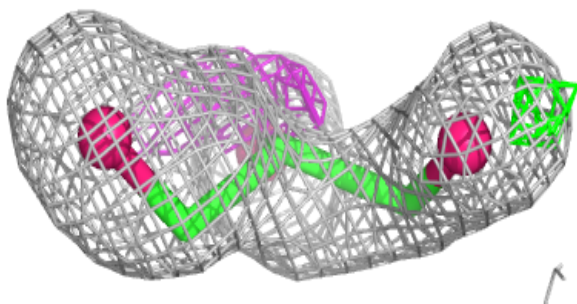
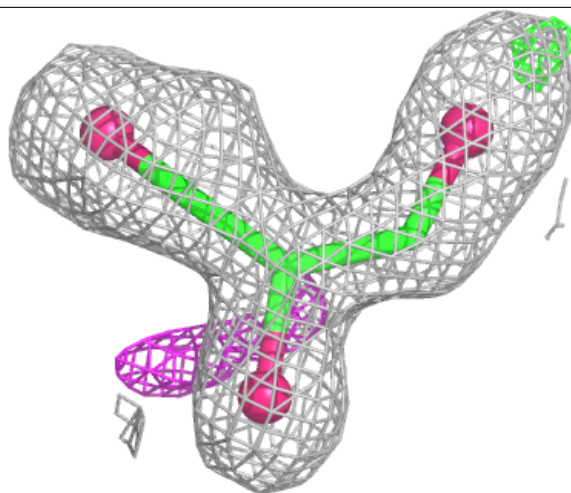
Electron density around GOL D 505:

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and green (positive)



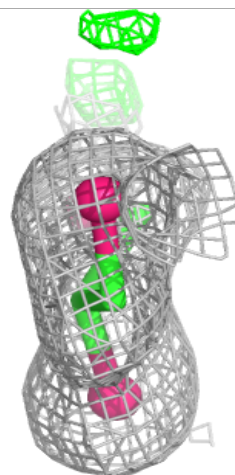
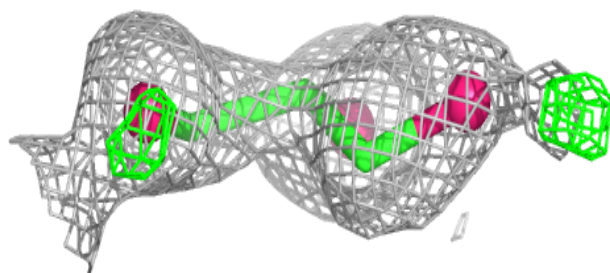
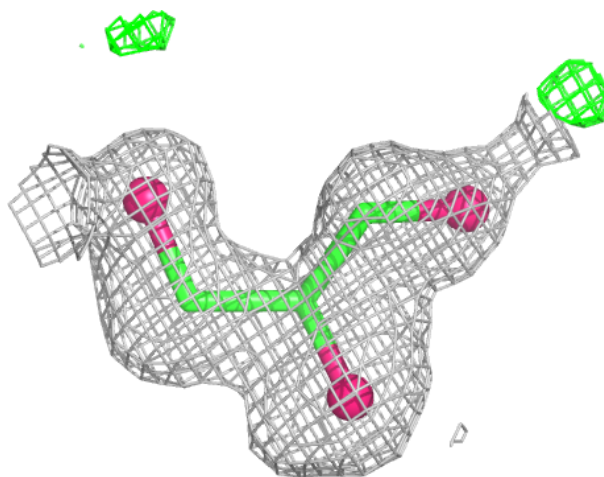
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and green (positive)



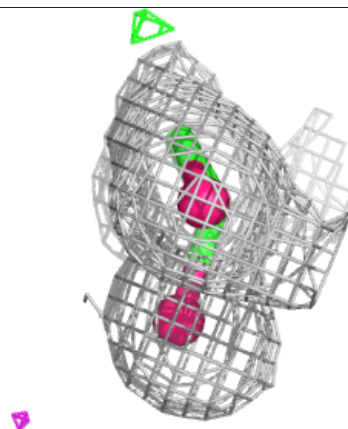
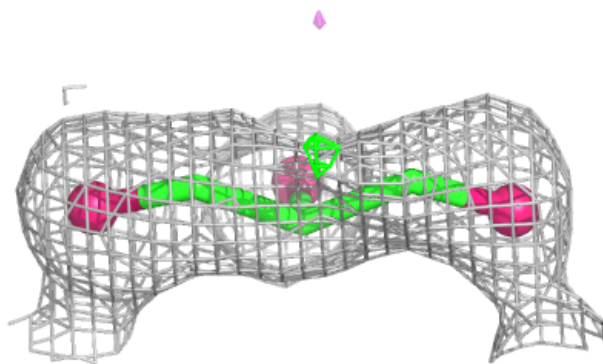
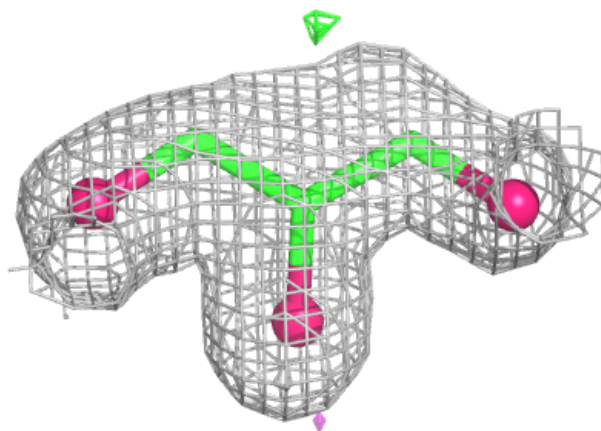
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and green (positive)



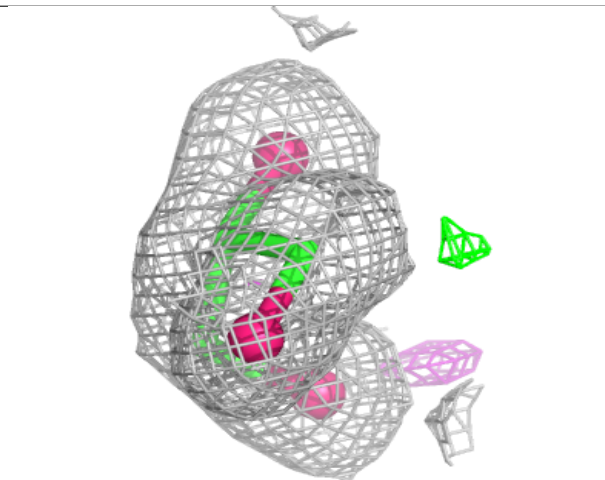
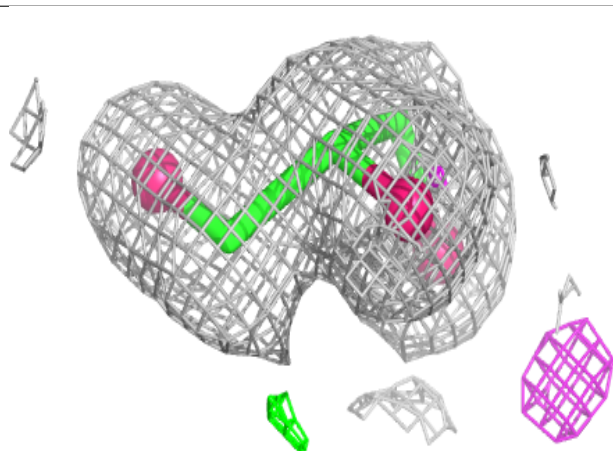
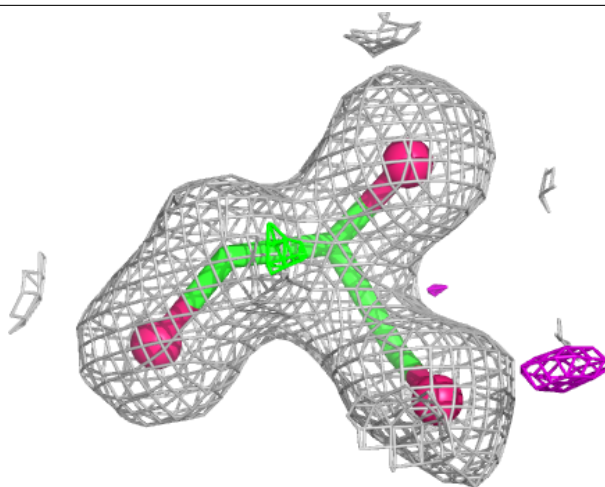
Electron density around GOL B 505:

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and green (positive)



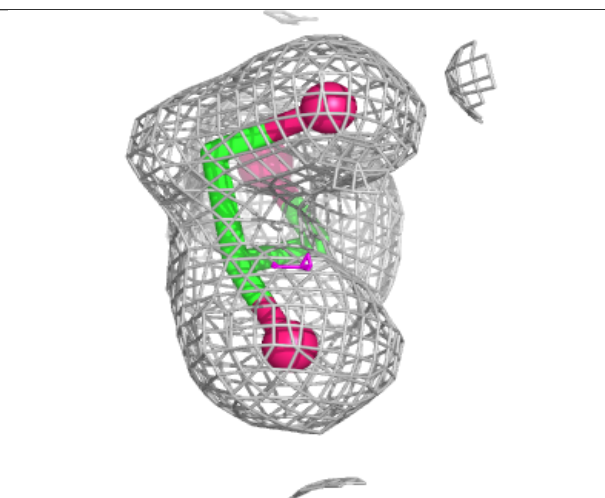
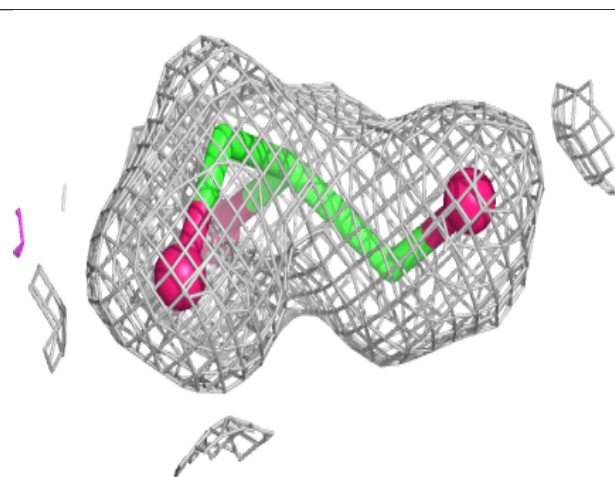
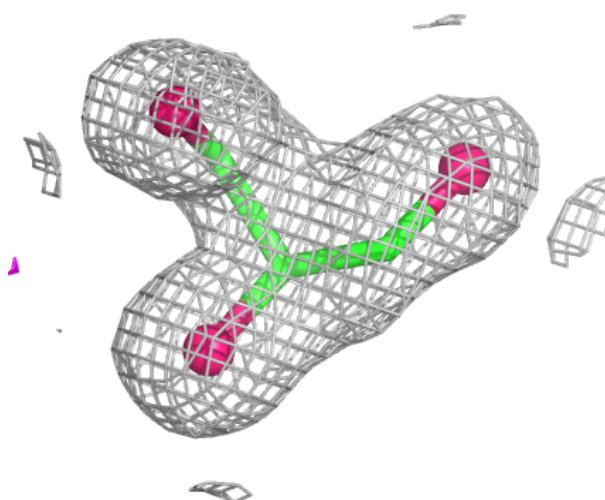
Electron density around GOL C 504:

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and green (positive)



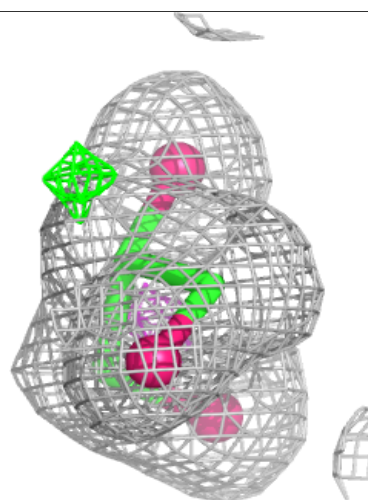
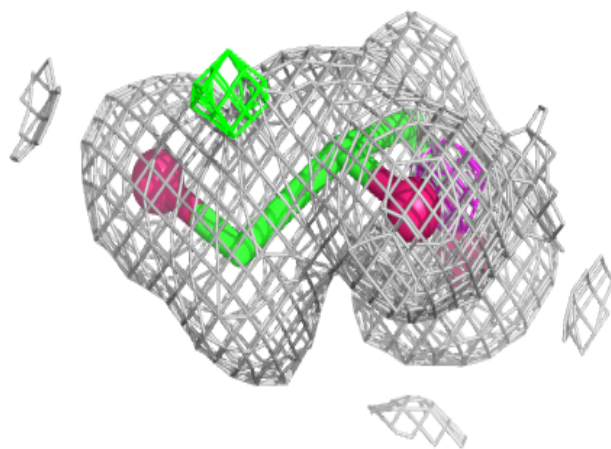
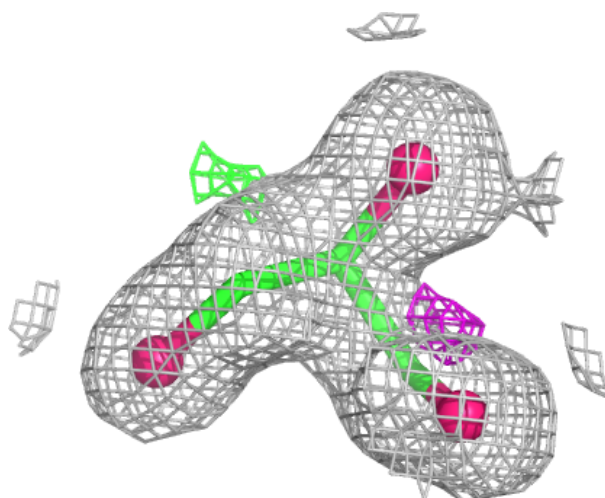
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and green (positive)



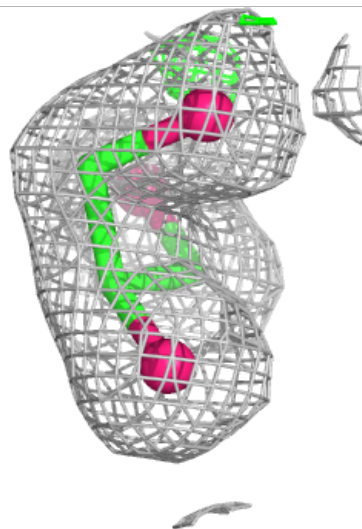
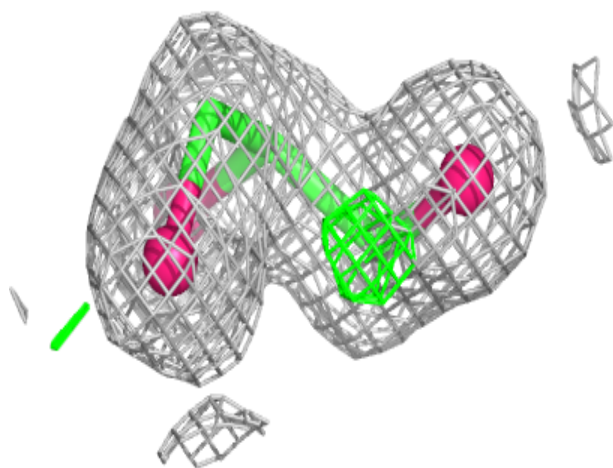
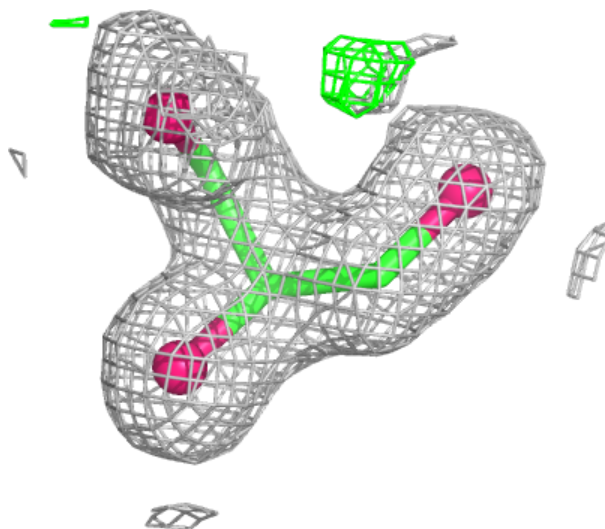
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and green (positive)



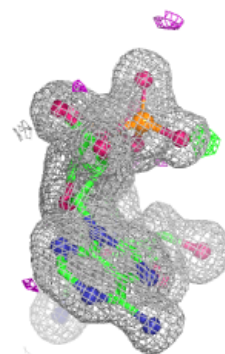
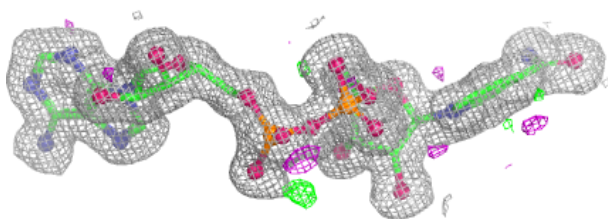
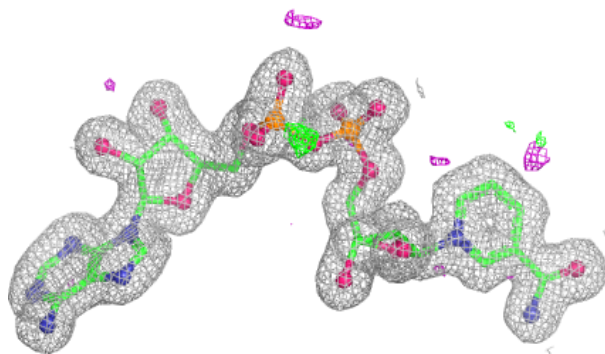
Electron density around GOL A 504:

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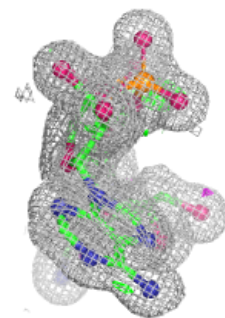
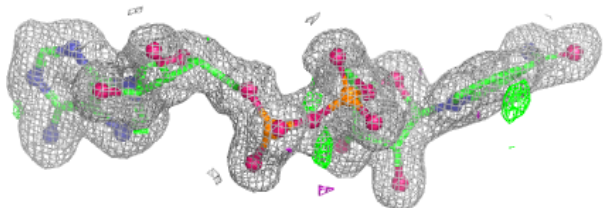
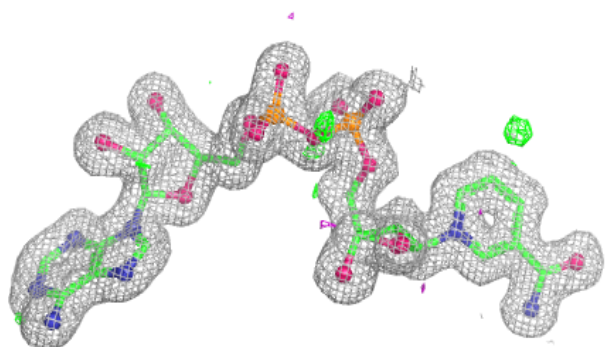


Electron density around NAD C 501:

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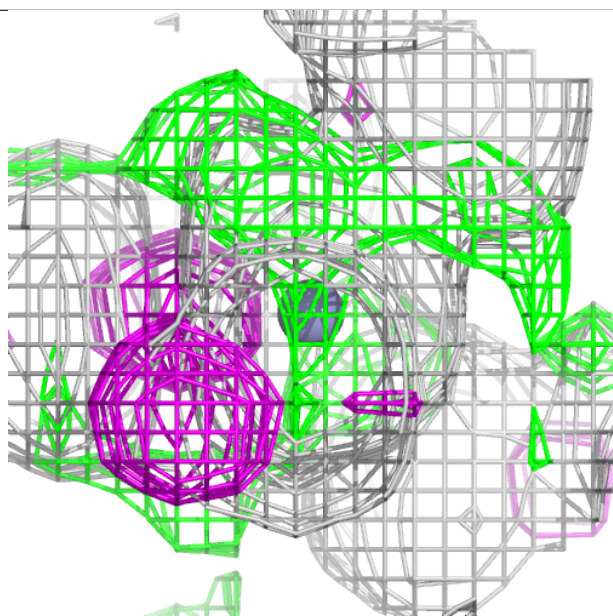
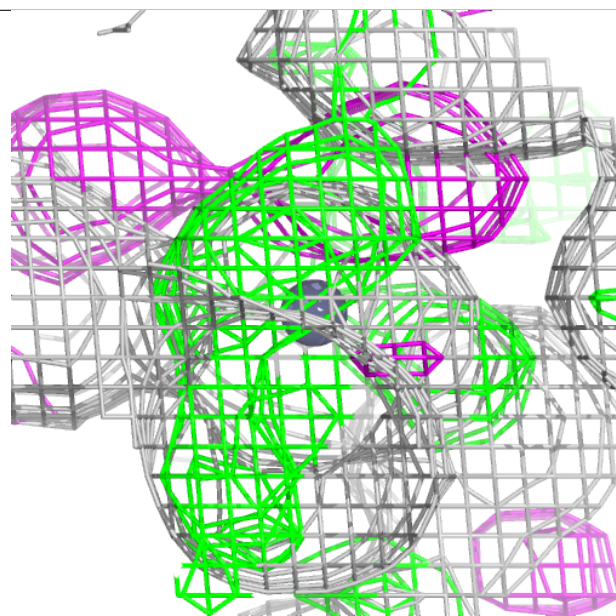
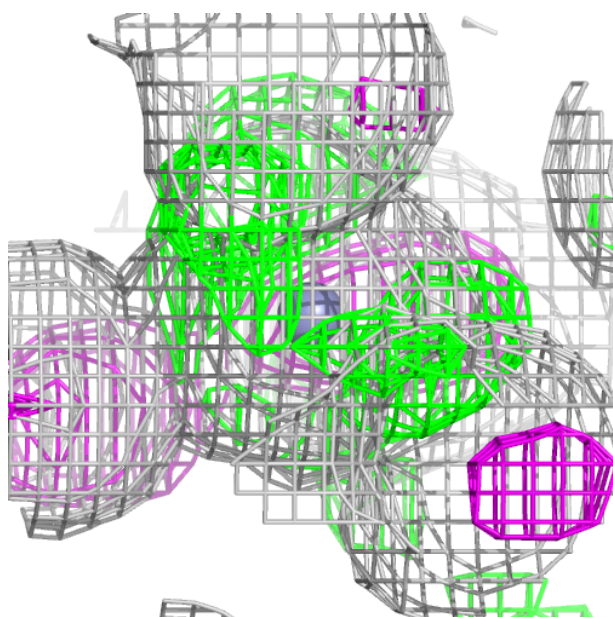
**Electron density around NAD D 501:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



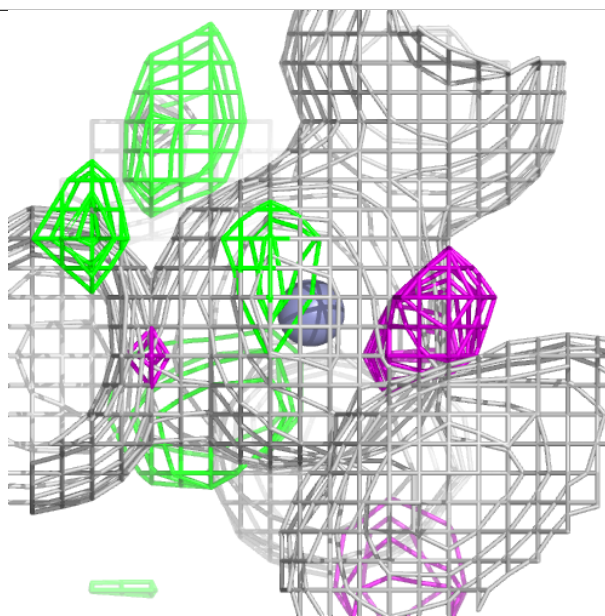
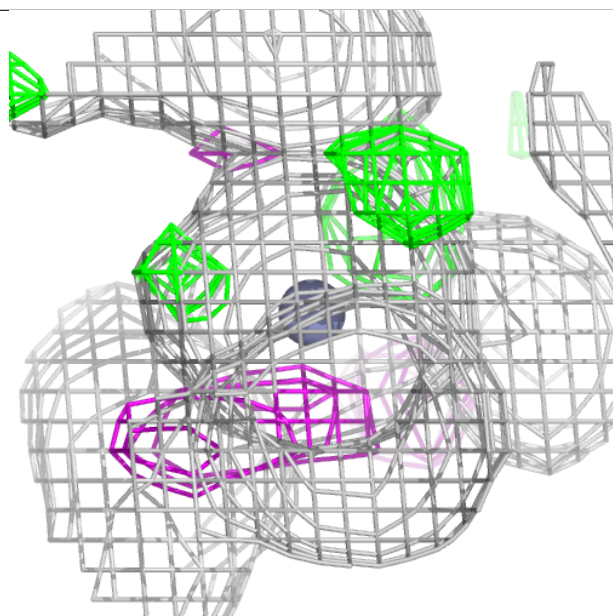
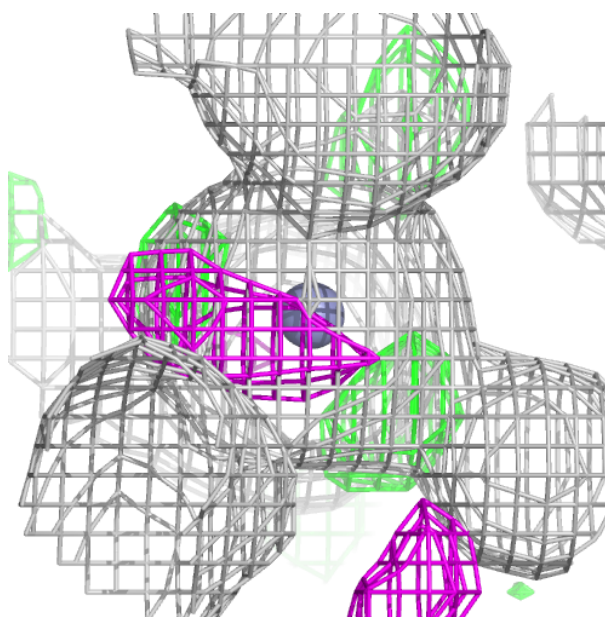
Electron density around ZN D 502:

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and green (positive)



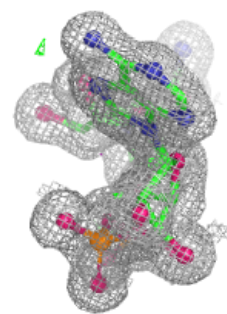
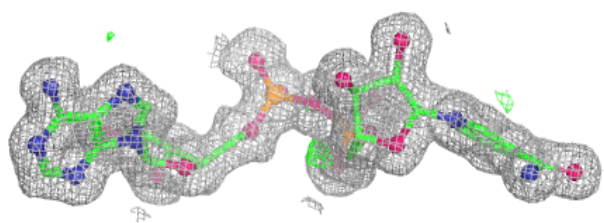
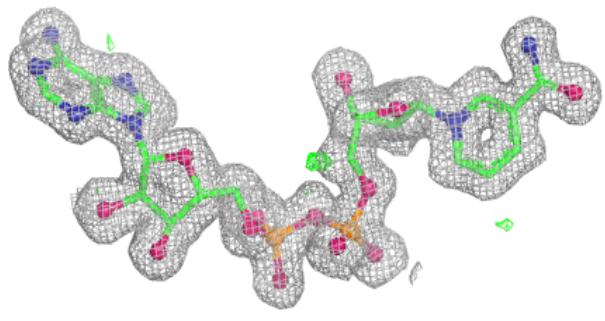
Electron density around ZN D 503:

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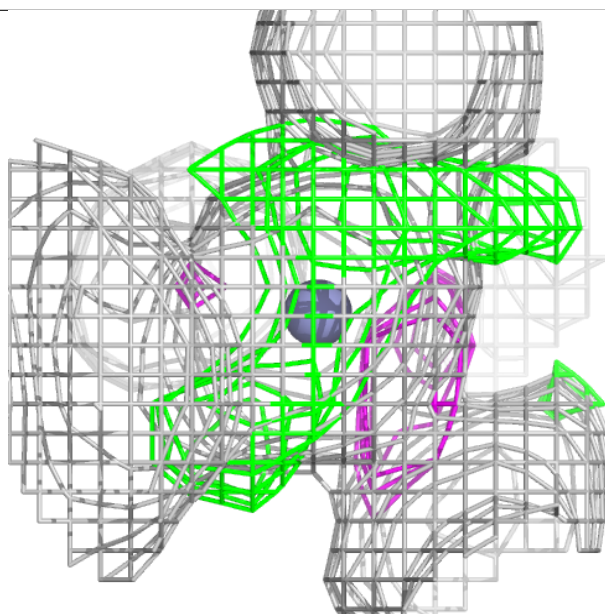
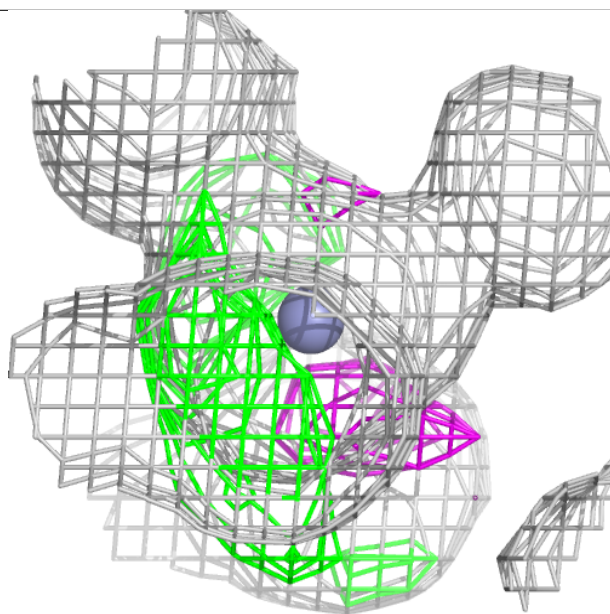
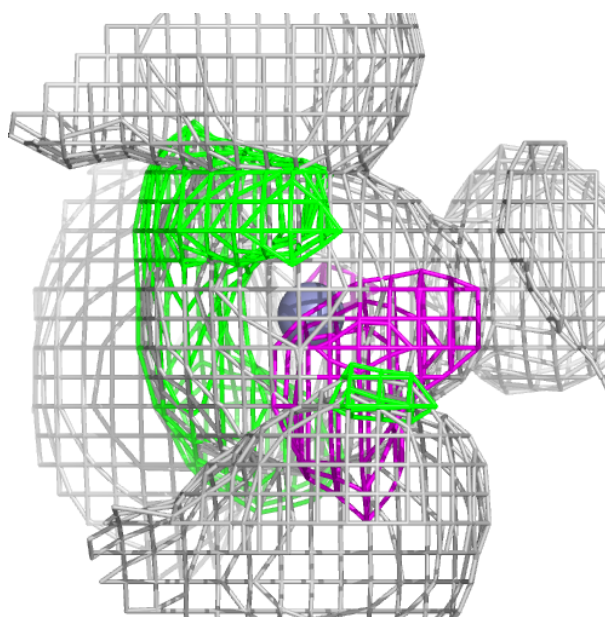
Electron density around NAD B 501:

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and green (positive)



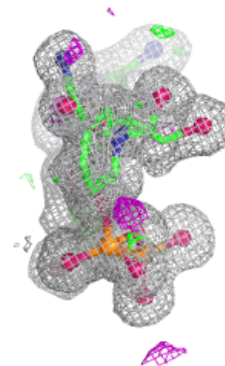
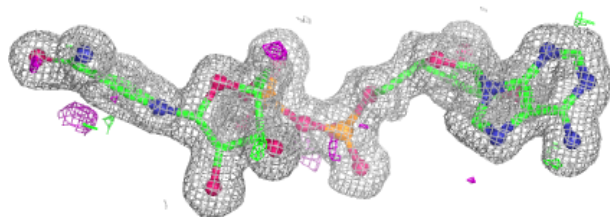
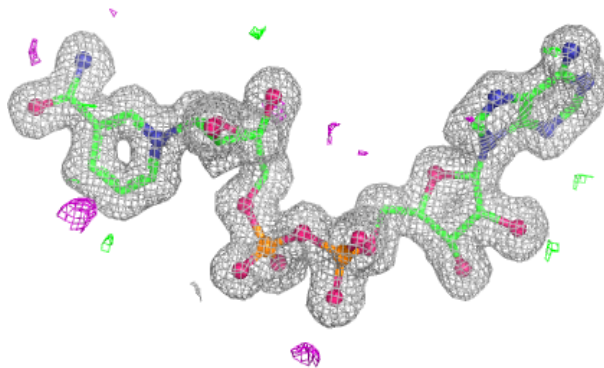
Electron density around ZN C 503:

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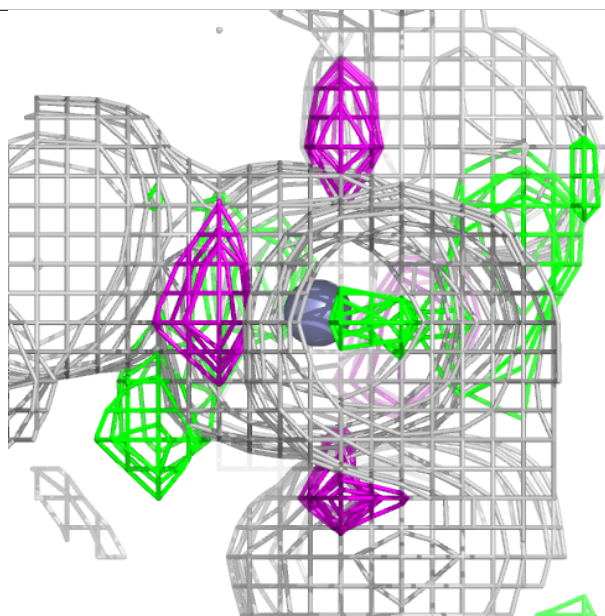
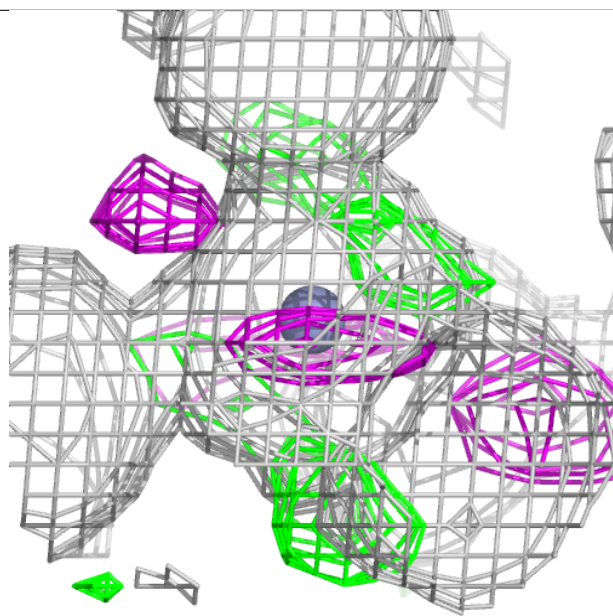
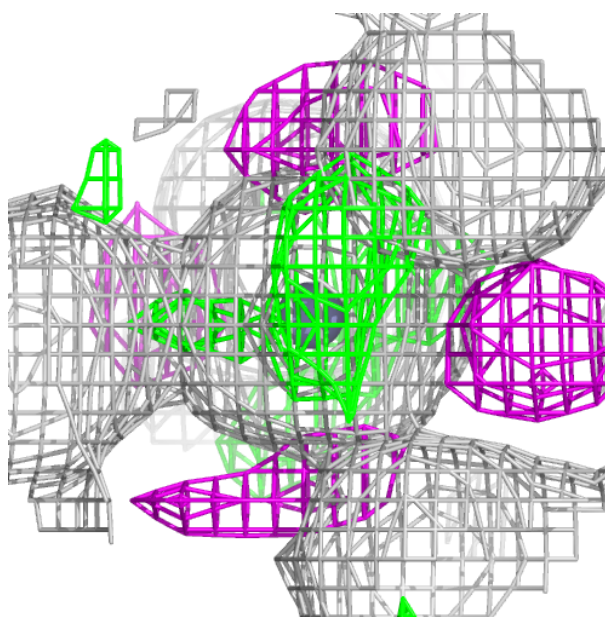
Electron density around NAD A 501:

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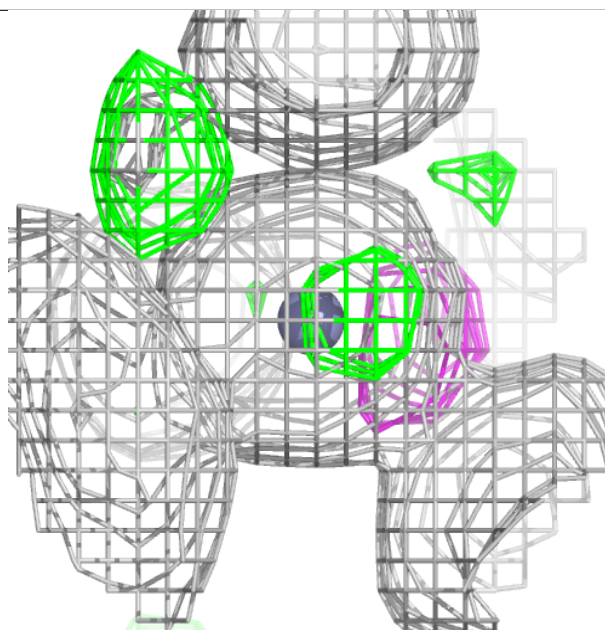
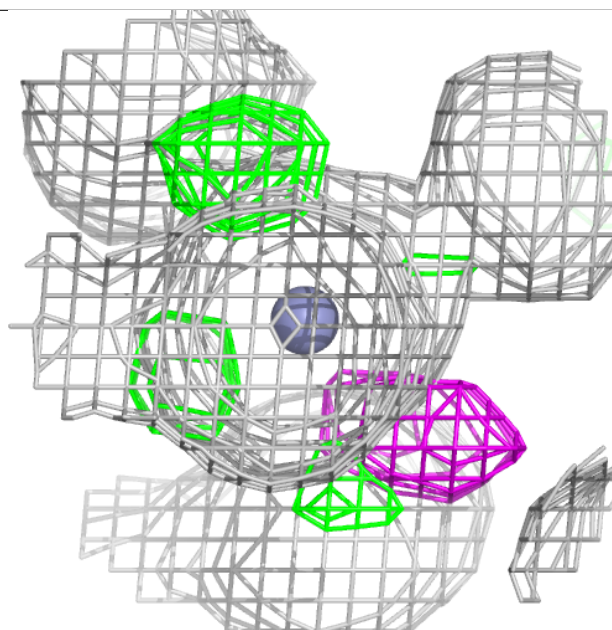
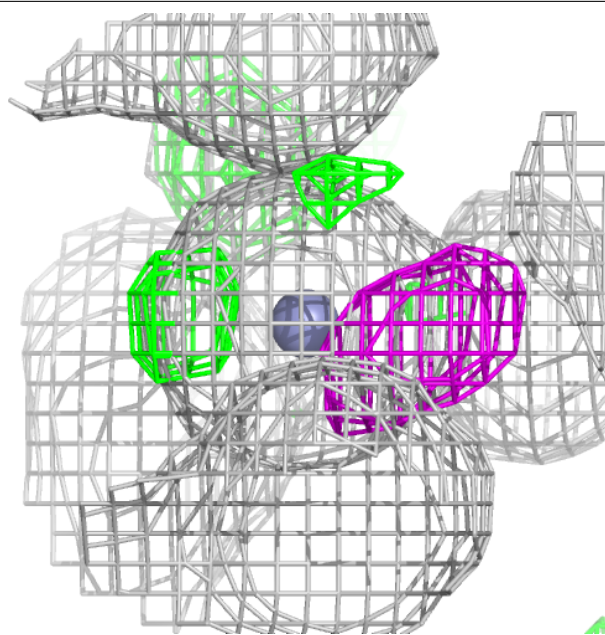
Electron density around ZN A 502:

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and green (positive)



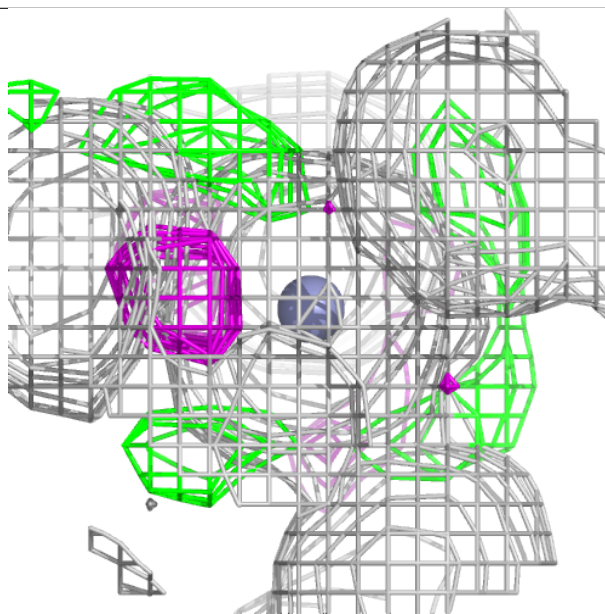
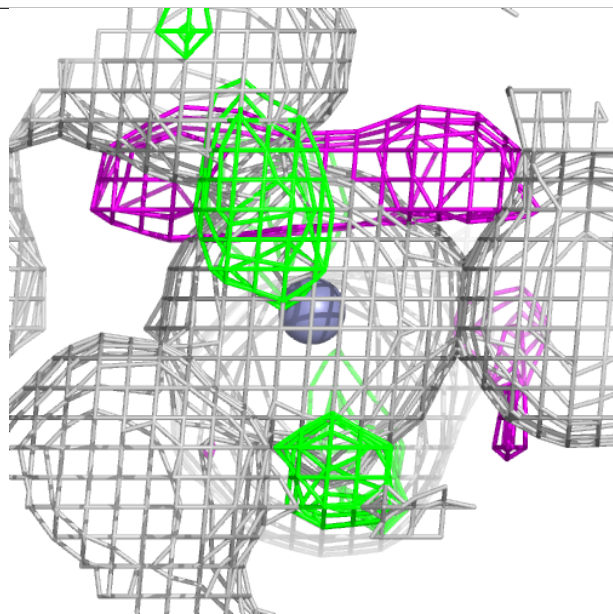
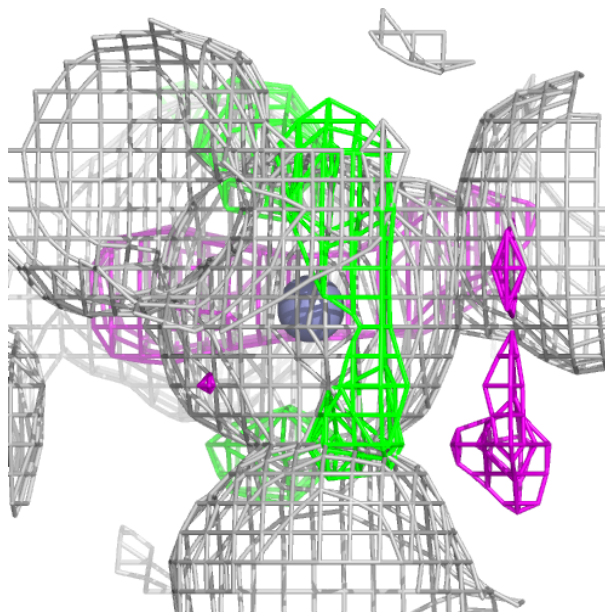
Electron density around ZN A 503:

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



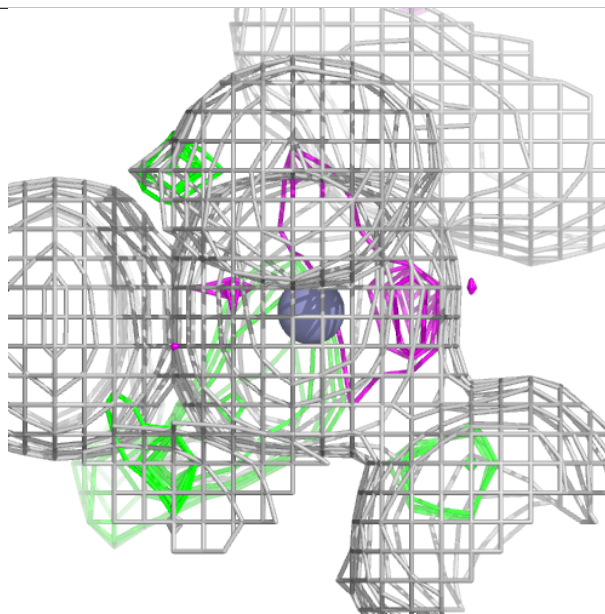
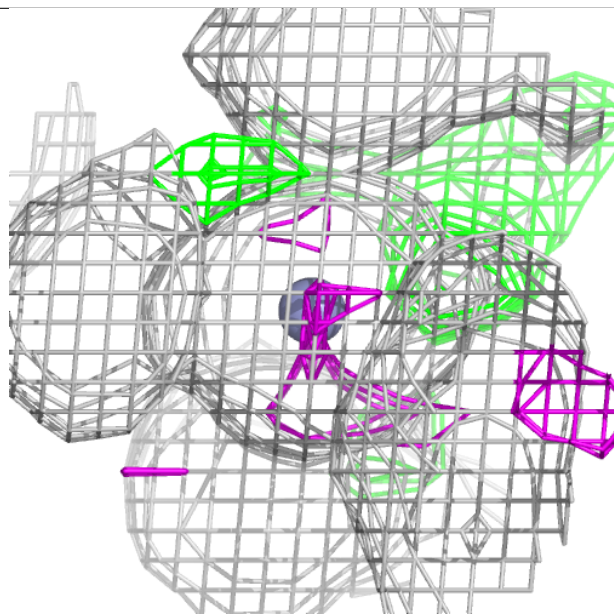
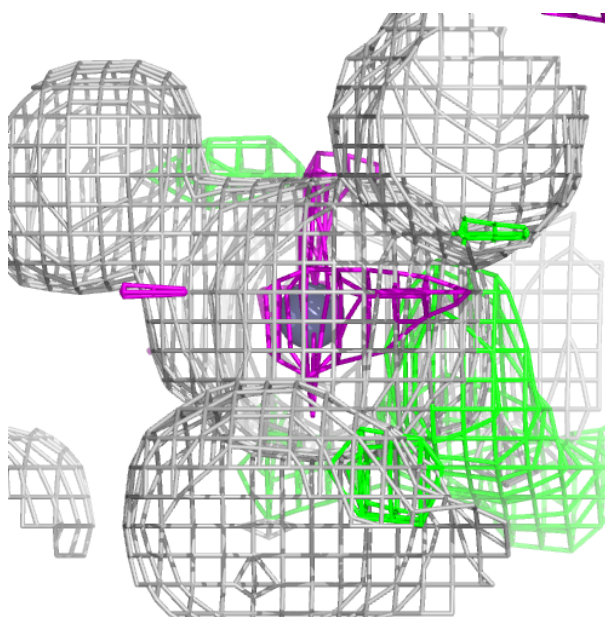
Electron density around ZN B 502:

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



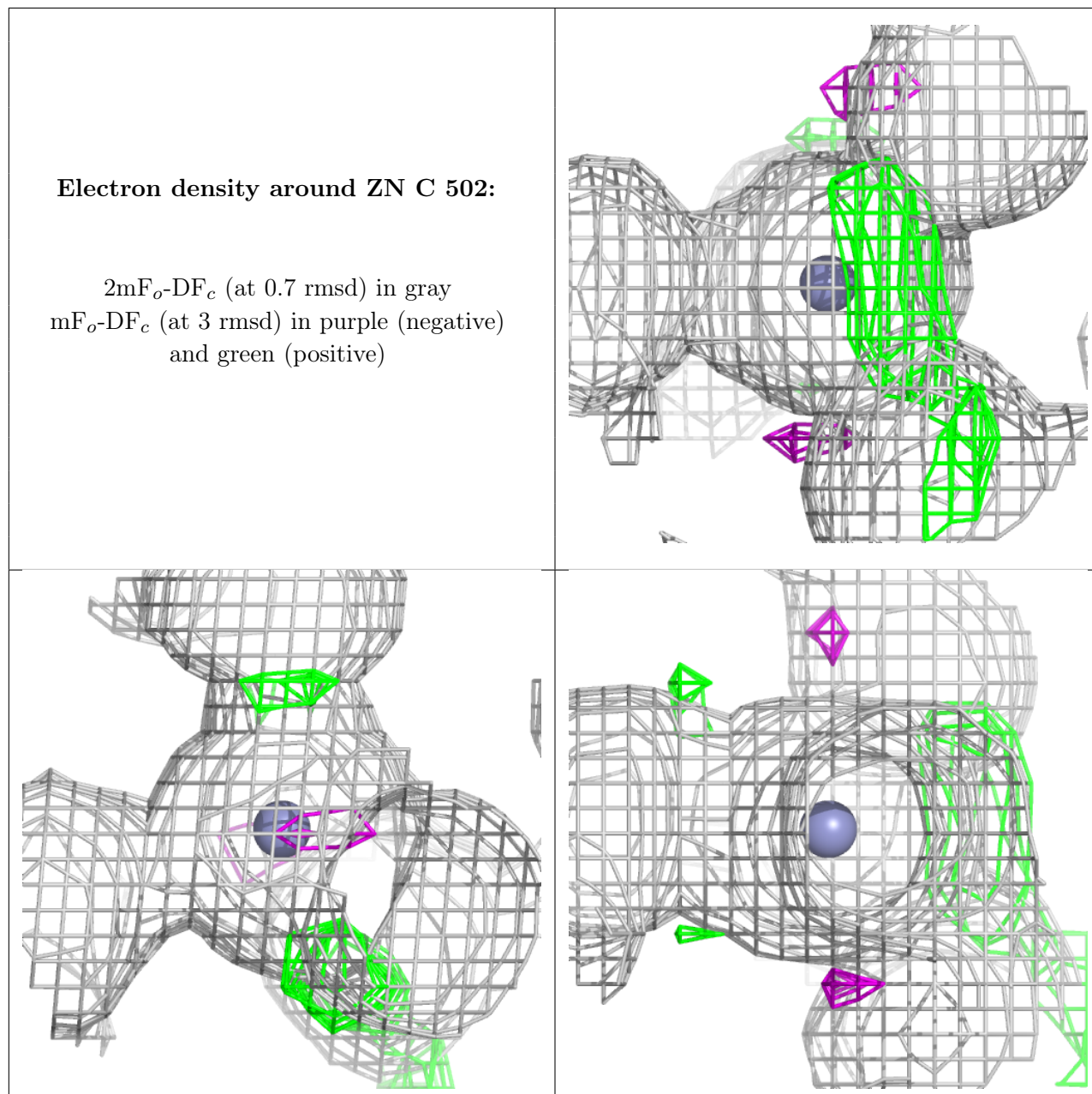
Electron density around ZN B 503:

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



Electron density around ZN C 502:

$2mF_o-DF_c$ (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.