



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 07:33 AM UTC

PDB ID : 9IL7 / pdb_00009il7
Title : Crystal Structure of SME E166A in Complex with Ceftaroline Fosamil
Authors : Dhankhar, K.; Hazra, S.
Deposited on : 2024-06-29
Resolution : 2.40 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

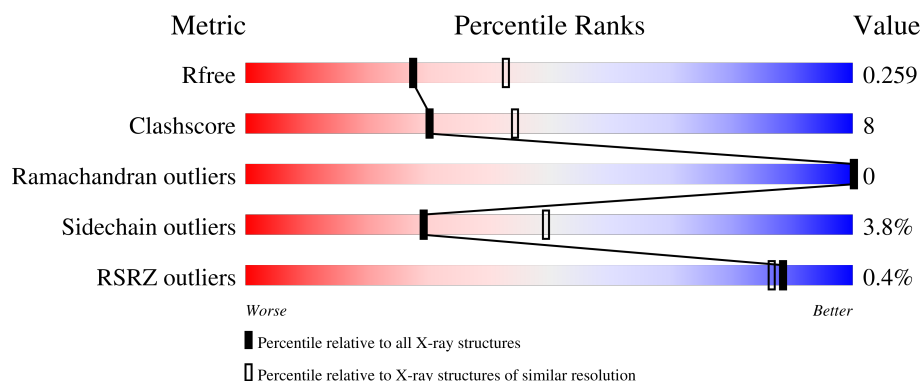
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	275	 81% 13% . .
1	B	275	 75% 17% . . .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	A1L2O	B	301	X	-	-	-
3	PEG	B	303	-	-	X	-
3	PEG	B	305	-	-	X	-
3	PEG	B	307	-	-	X	-
6	PG4	B	304	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 4425 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.

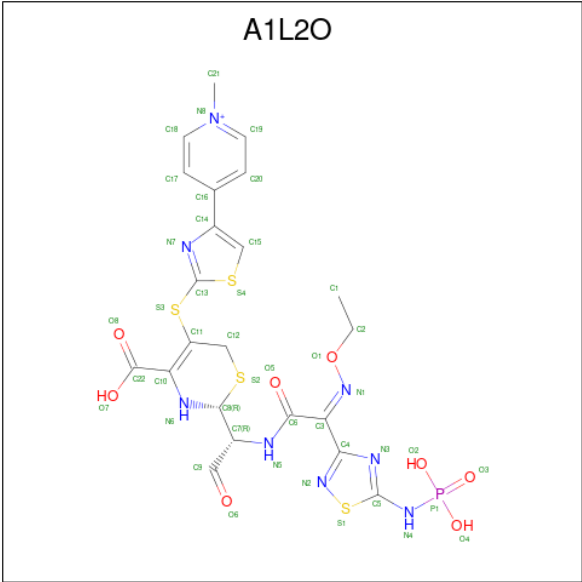
- Molecule 1 is a protein called beta-lactamase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	0	0	0
			2045	1280	363	395	7			
1	B	267	Total	C	N	O	S	0	2	0
			2054	1287	365	395	7			

There are 18 discrepancies between the modelled and reference sequences:

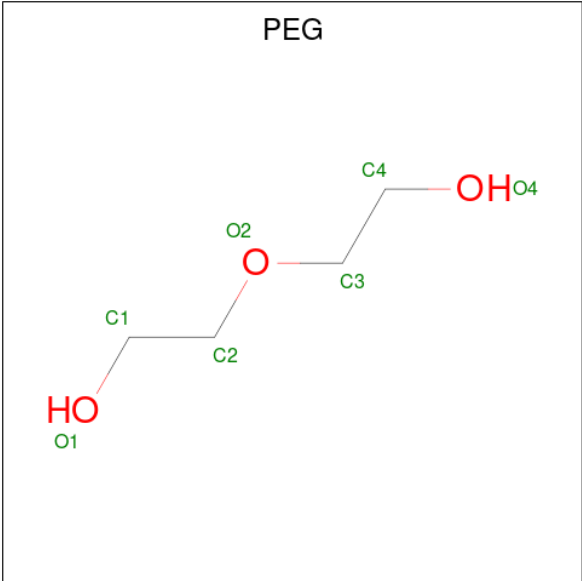
Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	MET	-	initiating methionine	UNP Q54488
A	0	GLY	-	expression tag	UNP Q54488
A	143	ALA	GLU	engineered mutation	UNP Q54488
A	268	HIS	-	expression tag	UNP Q54488
A	269	HIS	-	expression tag	UNP Q54488
A	270	HIS	-	expression tag	UNP Q54488
A	271	HIS	-	expression tag	UNP Q54488
A	272	HIS	-	expression tag	UNP Q54488
A	273	HIS	-	expression tag	UNP Q54488
B	-1	MET	-	initiating methionine	UNP Q54488
B	0	GLY	-	expression tag	UNP Q54488
B	143	ALA	GLU	engineered mutation	UNP Q54488
B	268	HIS	-	expression tag	UNP Q54488
B	269	HIS	-	expression tag	UNP Q54488
B	270	HIS	-	expression tag	UNP Q54488
B	271	HIS	-	expression tag	UNP Q54488
B	272	HIS	-	expression tag	UNP Q54488
B	273	HIS	-	expression tag	UNP Q54488

- Molecule 2 is (2 {R})-2-[(1 {R})-1-[(2 {Z})-2-ethoxyimino-2-[5-(phosphonoamino)-1,2,4-thiadiazol-3-yl]ethanoyl]amino]-2-oxidanylidene-ethyl]-5-[[4-(1-methylpyridin-4-yl)-1,3-thiazol-2-yl]sulfanyl]-3,6-dihydro-2 {H}-1,3-thiazine-4-carboxylic acid (CCD ID: A1L2O) (formula: C₂₂H₂₄N₈O₈PS₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	S	0	0
			43	22	8	8	1	4		
2	B	1	Total	C	N	O	P	S	0	0
			43	22	8	8	1	4		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		

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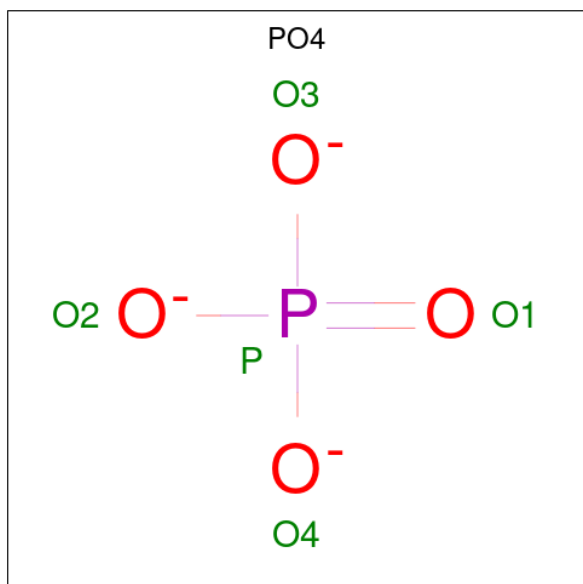
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		
3	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	7	Total	Cl	0	0
			7	7		
4	B	9	Total	Cl	0	0
			9	9		

- Molecule 5 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P) (labeled as "Ligand of Interest" by depositor).



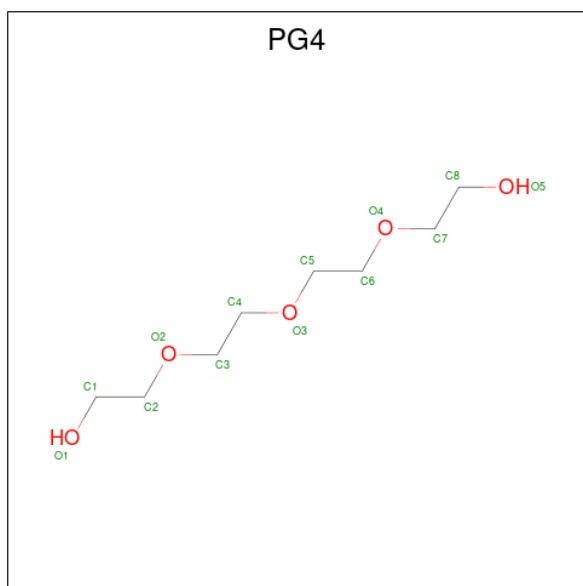
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	P	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	P	0	0
			5	4	1		
5	B	1	Total	O	P	0	0
			5	4	1		
5	B	1	Total	O	P	0	0
			5	4	1		
5	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 6 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			13	8	5		

- Molecule 7 is IODIDE ION (CCD ID: IOD) (formula: I) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	I	0	0
			1	1		

- Molecule 8 is SODIUM ION (CCD ID: NA) (formula: Na) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	B	1	Total 1	Na 1	0	0

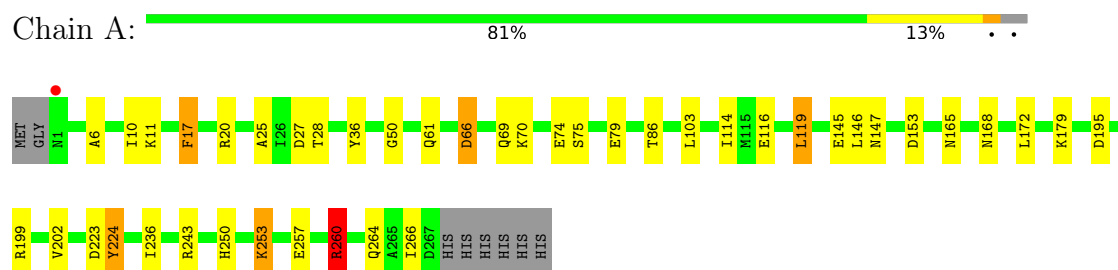
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	57	Total 57	O 57	0	0
9	B	85	Total 85	O 85	0	0

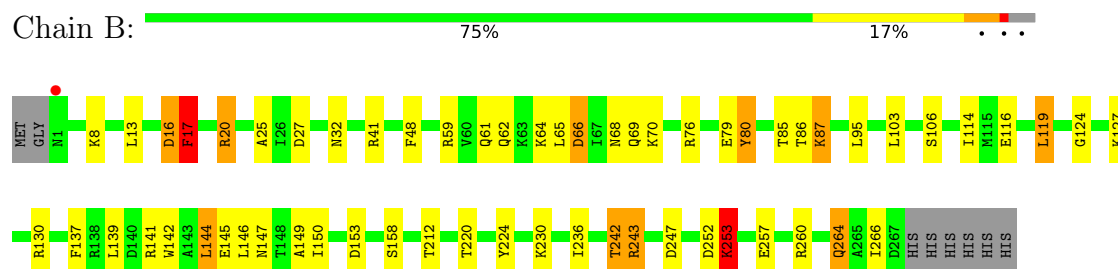
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry 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: beta-lactamase



- Molecule 1: beta-lactamase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	35.62Å 51.08Å 129.79Å 90.00° 90.12° 90.00°	Depositor
Resolution (Å)	21.74 – 2.40 21.74 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.4 (21.74-2.40) 98.4 (21.74-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.31 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.190 , 0.260 0.190 , 0.259	Depositor DCC
R_{free} test set	922 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	13.6	Xtriage
Anisotropy	0.204	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.046 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	4425	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 56.50 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.7111e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: IOD, NA, PEG, A1L2O, CL, PO4, PG4

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	0.72	0/2080	1.47	16/2807 (0.6%)
1	B	0.93	0/2095	1.60	24/2827 (0.8%)
All	All	0.83	0/4175	1.54	40/5634 (0.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3

There are no bond length outliers.

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	153	ASP	CA-CB-CG	7.50	120.10	112.60
1	B	76	ARG	CD-NE-CZ	-7.30	114.18	124.40
1	A	153	ASP	CA-CB-CG	7.26	119.86	112.60
1	B	130	ARG	NE-CZ-NH2	-7.08	112.83	119.20
1	A	224	TYR	CB-CA-C	6.96	123.05	109.33
1	A	20	ARG	N-CA-CB	-6.91	99.21	110.81
1	B	116	GLU	CB-CA-C	6.86	124.37	110.38
1	B	253	LYS	CB-CA-C	-6.71	99.44	110.85
1	B	17	PHE	CA-CB-CG	6.60	120.40	113.80
1	A	253	LYS	CB-CA-C	-6.60	99.63	110.85
1	B	264	GLN	CB-CA-C	-6.56	99.69	110.85
1	B	79	GLU	CB-CG-CD	6.43	123.54	112.60
1	A	145	GLU	CB-CG-CD	-6.40	101.72	112.60
1	A	224	TYR	N-CA-CB	-6.34	100.24	111.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	61	GLN	N-CA-CB	-6.32	100.83	110.12
1	B	20	ARG	CD-NE-CZ	6.28	133.19	124.40
1	B	16	ASP	CA-CB-CG	6.20	118.80	112.60
1	A	86	THR	CA-CB-OG1	-6.19	100.31	109.60
1	A	61	GLN	N-CA-CB	-6.16	101.07	110.12
1	B	212	THR	CA-CB-OG1	-5.96	100.66	109.60
1	B	66	ASP	CA-CB-CG	5.94	118.54	112.60
1	B	243	ARG	CD-NE-CZ	5.93	132.71	124.40
1	A	116	GLU	CB-CA-C	5.74	122.09	110.38
1	B	86	THR	CA-CB-OG1	-5.62	101.17	109.60
1	B	85	THR	N-CA-CB	5.61	118.37	110.12
1	B	264	GLN	N-CA-CB	5.59	118.43	110.16
1	B	137	PHE	N-CA-CB	-5.55	101.44	109.83
1	A	79	GLU	CB-CG-CD	5.52	121.98	112.60
1	B	242	THR	CA-CB-OG1	-5.51	101.33	109.60
1	B	224	TYR	CB-CA-C	5.46	120.80	109.38
1	B	80	TYR	N-CA-CB	-5.46	102.02	110.04
1	A	202	VAL	N-CA-CB	5.42	116.28	111.67
1	A	17	PHE	CA-CB-CG	5.30	119.10	113.80
1	B	127	LYS	CB-CG-CD	5.28	123.45	111.30
1	A	66	ASP	CA-CB-CG	5.17	117.77	112.60
1	A	223	ASP	CA-CB-CG	5.08	117.68	112.60
1	A	74	GLU	CB-CA-C	-5.07	102.33	110.74
1	A	36	TYR	CB-CA-C	5.05	117.92	110.14
1	B	20	ARG	CG-CD-NE	-5.03	100.94	112.00
1	B	68	ASN	CA-CB-CG	-5.00	107.60	112.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	199	ARG	Sidechain
1	A	243	ARG	Sidechain
1	A	260	ARG	Sidechain

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	2045	0	2028	22	0
1	B	2054	0	2048	46	0
2	A	43	0	0	1	0
2	B	43	0	0	2	0
3	A	7	0	10	0	0
3	B	35	0	50	13	0
4	A	7	0	0	0	0
4	B	9	0	0	1	0
5	A	10	0	0	0	0
5	B	15	0	0	1	0
6	B	13	0	18	7	0
7	B	1	0	0	0	0
8	B	1	0	0	0	0
9	A	57	0	0	5	0
9	B	85	0	0	1	0
All	All	4425	0	4154	70	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 8.

All (70) 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:B:150:ILE:HD11	3:B:303:PEG:O4	1.68	0.93
1:B:142:TRP:HB2	6:B:304:PG4:H82	1.53	0.88
3:B:303:PEG:H12	6:B:304:PG4:H42	1.61	0.81
1:A:66:ASP:H	1:A:69:GLN:HE21	1.34	0.76
1:B:17:PHE:O	1:B:243:ARG:NH1	2.20	0.74
1:B:13:LEU:HD12	3:B:307:PEG:H42	1.69	0.74
1:B:252:ASP:H	3:B:305:PEG:H32	1.52	0.73
1:B:150:ILE:CD1	3:B:303:PEG:O4	2.36	0.73
1:A:28:THR:O	1:B:62:GLN:HG2	1.89	0.72
1:A:260:ARG:HH22	1:A:264:GLN:NE2	1.87	0.72
1:B:17:PHE:CD2	1:B:243:ARG:HD2	2.26	0.69
1:B:66:ASP:H	1:B:69:GLN:HE21	1.38	0.69
1:B:16:ASP:C	3:B:307:PEG:H11	2.19	0.67
1:B:17:PHE:CE2	1:B:243:ARG:HD2	2.29	0.67
1:A:250:HIS:NE2	9:A:401:HOH:O	2.28	0.66
1:A:66:ASP:H	1:A:69:GLN:NE2	1.93	0.65
1:B:220[B]:THR:HG22	1:B:242:THR:HG22	1.79	0.65
1:B:141:ARG:HD3	6:B:304:PG4:H41	1.79	0.64
1:B:141:ARG:HD2	6:B:304:PG4:H51	1.79	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:260:ARG:HH22	1:A:264:GLN:HE22	1.47	0.61
1:B:252:ASP:H	3:B:305:PEG:C3	2.13	0.61
1:B:149:ALA:HB1	1:B:220[B]:THR:HG21	1.83	0.60
1:B:247:ASP:OD1	3:B:302:PEG:H41	2.02	0.59
1:B:16:ASP:O	3:B:307:PEG:H11	2.03	0.59
1:B:27:ASP:HB2	1:B:266:ILE:HD13	1.86	0.58
1:A:260:ARG:NH2	1:A:264:GLN:NE2	2.52	0.57
1:B:32:ASN:ND2	9:B:401:HOH:O	2.37	0.57
1:A:11:LYS:HE3	9:A:453:HOH:O	2.04	0.57
1:B:66:ASP:H	1:B:69:GLN:NE2	2.03	0.56
1:A:27:ASP:HB2	1:A:266:ILE:HD13	1.87	0.55
1:B:253:LYS:HB2	3:B:305:PEG:H11	1.88	0.55
1:B:145:GLU:CB	6:B:304:PG4:O5	2.56	0.54
1:B:141:ARG:CD	6:B:304:PG4:H51	2.38	0.54
1:B:87:LYS:NZ	5:B:321:PO4:O2	2.42	0.53
1:A:103:LEU:C	1:A:103:LEU:HD23	2.34	0.52
2:B:301:A1L2O:O3	4:B:315:CL:CL	2.66	0.51
1:A:257:GLU:OE2	1:A:257:GLU:HA	2.09	0.50
1:B:114:ILE:HG22	1:B:119:LEU:HD22	1.93	0.50
1:B:13:LEU:HD12	3:B:307:PEG:C4	2.41	0.50
1:A:66:ASP:O	1:A:69:GLN:HG2	2.12	0.49
1:B:257:GLU:HA	1:B:257:GLU:OE2	2.13	0.49
1:A:168:ASN:HD22	1:A:172:LEU:HD12	1.78	0.48
1:B:103:LEU:C	1:B:103:LEU:HD23	2.38	0.48
1:A:114:ILE:HG22	1:A:119:LEU:HD22	1.95	0.47
1:B:147:ASN:ND2	2:B:301:A1L2O:N2	2.60	0.47
1:B:260:ARG:O	1:B:264:GLN:HG2	2.16	0.46
1:B:59:ARG:HH11	1:B:64:LYS:HD3	1.82	0.45
1:B:220[B]:THR:HG22	1:B:242:THR:CG2	2.44	0.45
1:B:66:ASP:O	1:B:69:GLN:HG2	2.17	0.45
1:B:87:LYS:HE3	1:B:87:LYS:HB3	1.54	0.44
1:A:146:LEU:C	1:A:146:LEU:HD12	2.42	0.44
1:B:65:LEU:HD13	1:B:95:LEU:HD21	1.99	0.43
1:B:80:TYR:HB3	1:B:144:LEU:HD21	2.00	0.43
1:A:165:ASN:HA	9:A:406:HOH:O	2.17	0.43
1:B:252:ASP:N	3:B:305:PEG:H21	2.34	0.43
1:A:6:ALA:O	1:A:10:ILE:HG12	2.19	0.43
1:B:119:LEU:O	1:B:124:GLY:HA3	2.18	0.42
1:B:144:LEU:HD12	1:B:144:LEU:HA	1.90	0.42
1:A:147:ASN:ND2	2:A:301:A1L2O:N2	2.66	0.42
1:A:253:LYS:HD2	9:A:433:HOH:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:220[B]:THR:HG22	1:B:242:THR:CB	2.50	0.42
1:B:41:ARG:HA	1:B:158:SER:O	2.20	0.41
1:B:139:LEU:HD11	1:B:146:LEU:HD22	2.02	0.41
1:A:50:GLY:HA3	1:A:224:TYR:OH	2.21	0.41
1:A:195:ASP:HB2	9:A:424:HOH:O	2.21	0.41
1:B:25:ALA:HA	1:B:236:ILE:O	2.21	0.41
1:B:146:LEU:HD12	1:B:146:LEU:C	2.46	0.41
3:B:303:PEG:C1	6:B:304:PG4:H42	2.40	0.41
1:A:25:ALA:HA	1:A:236:ILE:O	2.21	0.40
1:B:48:PHE:CE1	1:B:146:LEU:HD21	2.56	0.40

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	265/275 (96%)	260 (98%)	5 (2%)	0	100	100
1	B	267/275 (97%)	260 (97%)	7 (3%)	0	100	100
All	All	532/550 (97%)	520 (98%)	12 (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	212/223 (95%)	206 (97%)	6 (3%)	38	60
1	B	214/223 (96%)	204 (95%)	10 (5%)	23	41
All	All	426/446 (96%)	410 (96%)	16 (4%)	29	49

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	PHE
1	A	70	LYS
1	A	75	SER
1	A	119	LEU
1	A	179	LYS
1	A	260	ARG
1	B	8	LYS
1	B	17	PHE
1	B	20	ARG
1	B	70	LYS
1	B	87	LYS
1	B	106	SER
1	B	119	LEU
1	B	144	LEU
1	B	230	LYS
1	B	253	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	69	GLN
1	A	147	ASN
1	A	168	ASN
1	A	264	GLN
1	B	32	ASN
1	B	69	GLN
1	B	147	ASN
1	B	168	ASN

5.3.3 RNA ⓘ

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 32 ligands modelled in this entry, 18 are monoatomic - leaving 14 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	PEG	B	307	-	6,6,6	0.43	0	5,5,5	0.41	0
5	PO4	B	321	-	4,4,4	1.21	1 (25%)	6,6,6	0.36	0
2	A1L2O	B	301	1	43,46,46	1.95	7 (16%)	47,65,65	3.91	21 (44%)
3	PEG	A	302	-	6,6,6	0.17	0	5,5,5	0.12	0
6	PG4	B	304	-	12,12,12	0.65	0	11,11,11	0.62	0
3	PEG	B	303	-	6,6,6	0.28	0	5,5,5	0.20	0
3	PEG	B	302	-	6,6,6	0.38	0	5,5,5	0.19	0
3	PEG	B	305	-	6,6,6	0.89	0	5,5,5	0.59	0
5	PO4	A	310	-	4,4,4	0.67	0	6,6,6	0.59	0
5	PO4	B	319	-	4,4,4	1.34	1 (25%)	6,6,6	0.77	0
2	A1L2O	A	301	1	43,46,46	1.70	7 (16%)	47,65,65	2.77	15 (31%)
3	PEG	B	306	-	6,6,6	0.54	0	5,5,5	0.44	0
5	PO4	A	311	-	4,4,4	0.54	0	6,6,6	0.74	0
5	PO4	B	320	-	4,4,4	1.63	1 (25%)	6,6,6	0.77	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
3	PEG	B	307	-	-	2/4/4/4	-
2	A1L2O	B	301	1	1/1/8/12	12/24/52/52	0/3/4/4
3	PEG	A	302	-	-	3/4/4/4	-
6	PG4	B	304	-	-	6/10/10/10	-
3	PEG	B	303	-	-	2/4/4/4	-
3	PEG	B	302	-	-	2/4/4/4	-
3	PEG	B	305	-	-	2/4/4/4	-
2	A1L2O	A	301	1	-	6/24/52/52	0/3/4/4
3	PEG	B	306	-	-	1/4/4/4	-

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	A1L2O	P1-O3	6.77	1.56	1.46
2	B	301	A1L2O	P1-O3	6.60	1.56	1.46
2	B	301	A1L2O	P1-N4	6.38	1.73	1.63
2	B	301	A1L2O	C3-C4	-3.54	1.41	1.46
2	B	301	A1L2O	C16-C14	3.35	1.52	1.47
2	A	301	A1L2O	C10-C22	3.17	1.53	1.48
2	B	301	A1L2O	C5-N3	3.16	1.36	1.31
2	A	301	A1L2O	P1-N4	2.77	1.67	1.63
2	A	301	A1L2O	C16-C14	2.67	1.51	1.47
2	A	301	A1L2O	C14-N7	2.60	1.43	1.39
2	A	301	A1L2O	C5-N3	2.52	1.35	1.31
5	B	320	PO4	P-O1	2.26	1.55	1.50
5	B	321	PO4	P-O1	2.26	1.55	1.50
2	B	301	A1L2O	C15-S4	2.25	1.77	1.71
5	B	319	PO4	P-O1	2.14	1.55	1.50
2	B	301	A1L2O	P1-O4	2.13	1.62	1.56
2	A	301	A1L2O	C3-C6	2.08	1.52	1.48

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	A1L2O	C3-C6-N5	13.02	134.81	114.27
2	B	301	A1L2O	C7-N5-C6	12.03	145.74	121.97
2	B	301	A1L2O	O5-C6-N5	-9.85	105.74	123.09
2	A	301	A1L2O	C3-C6-N5	8.84	128.21	114.27
2	B	301	A1L2O	O3-P1-N4	-7.16	99.56	111.03
2	A	301	A1L2O	S1-C5-N4	6.28	128.30	124.42
2	A	301	A1L2O	O5-C6-C3	-6.07	111.84	121.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	A1L2O	C17-C16-C14	6.02	127.61	120.87
2	B	301	A1L2O	C12-S2-C8	5.55	104.53	94.36
2	A	301	A1L2O	C20-C16-C14	5.39	126.91	120.87
2	A	301	A1L2O	C12-S2-C8	5.19	103.87	94.36
2	A	301	A1L2O	O1-C2-C1	5.09	121.83	109.10
2	A	301	A1L2O	C17-C16-C14	-4.47	115.86	120.87
2	A	301	A1L2O	O1-N1-C3	4.35	117.65	111.59
2	B	301	A1L2O	C9-C7-N5	4.29	116.57	109.80
2	B	301	A1L2O	S4-C13-N7	-3.99	112.61	115.15
2	B	301	A1L2O	C20-C16-C14	-3.97	116.42	120.87
2	B	301	A1L2O	O6-C9-C7	-3.91	114.51	124.86
2	B	301	A1L2O	C4-N2-S1	3.67	112.58	107.62
2	A	301	A1L2O	O8-C22-C10	-3.59	113.94	120.20
2	B	301	A1L2O	S1-C5-N4	-3.48	122.27	124.42
2	B	301	A1L2O	C5-S1-N2	-3.44	89.01	91.80
2	B	301	A1L2O	O8-C22-C10	-3.36	114.35	120.20
2	B	301	A1L2O	O7-C22-O8	3.30	131.75	123.90
2	A	301	A1L2O	O6-C9-C7	-2.92	117.15	124.86
2	B	301	A1L2O	O1-C2-C1	2.79	116.07	109.10
2	A	301	A1L2O	C6-C3-N1	2.68	133.81	120.64
2	A	301	A1L2O	O4-P1-O3	-2.58	106.99	113.45
2	B	301	A1L2O	C8-C7-N5	-2.55	104.53	109.93
2	B	301	A1L2O	C15-C14-N7	-2.40	111.77	114.53
2	A	301	A1L2O	C9-C7-N5	2.36	113.52	109.80
2	A	301	A1L2O	C11-S3-C13	2.33	105.93	101.81
2	A	301	A1L2O	C4-N2-S1	2.30	110.73	107.62
2	B	301	A1L2O	C20-C16-C17	-2.07	115.94	118.57
2	B	301	A1L2O	O1-N1-C3	2.06	114.46	111.59
2	B	301	A1L2O	O2-P1-O4	2.04	113.08	107.59

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	301	A1L2O	C7

All (36) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	A1L2O	N6-C10-C22-O7
2	A	301	A1L2O	C1-C2-O1-N1
2	B	301	A1L2O	C4-C3-N1-O1
2	B	301	A1L2O	C6-C3-N1-O1

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Mol	Chain	Res	Type	Atoms
2	B	301	A1L2O	C3-C6-N5-C7
2	B	301	A1L2O	O5-C6-N5-C7
2	B	301	A1L2O	N6-C10-C22-O7
2	B	301	A1L2O	C1-C2-O1-N1
6	B	304	PG4	O2-C3-C4-O3
3	B	307	PEG	O2-C3-C4-O4
3	B	302	PEG	O1-C1-C2-O2
6	B	304	PG4	O3-C5-C6-O4
3	A	302	PEG	O2-C3-C4-O4
2	B	301	A1L2O	C4-C3-C6-N5
2	A	301	A1L2O	N1-C3-C6-N5
2	B	301	A1L2O	N1-C3-C6-N5
3	B	303	PEG	O1-C1-C2-O2
3	B	302	PEG	O2-C3-C4-O4
2	A	301	A1L2O	C4-C3-C6-O5
2	A	301	A1L2O	C4-C3-C6-N5
2	B	301	A1L2O	C4-C3-C6-O5
6	B	304	PG4	C3-C4-O3-C5
3	B	305	PEG	O2-C3-C4-O4
6	B	304	PG4	C6-C5-O3-C4
6	B	304	PG4	O1-C1-C2-O2
3	B	306	PEG	C4-C3-O2-C2
3	B	303	PEG	C4-C3-O2-C2
2	B	301	A1L2O	N1-C3-C6-O5
6	B	304	PG4	C4-C3-O2-C2
3	A	302	PEG	C4-C3-O2-C2
3	A	302	PEG	C1-C2-O2-C3
3	B	305	PEG	C4-C3-O2-C2
2	B	301	A1L2O	C11-C10-C22-O8
2	A	301	A1L2O	N1-C3-C6-O5
3	B	307	PEG	C1-C2-O2-C3
2	B	301	A1L2O	N6-C10-C22-O8

There are no ring outliers.

8 monomers are involved in 22 short contacts:

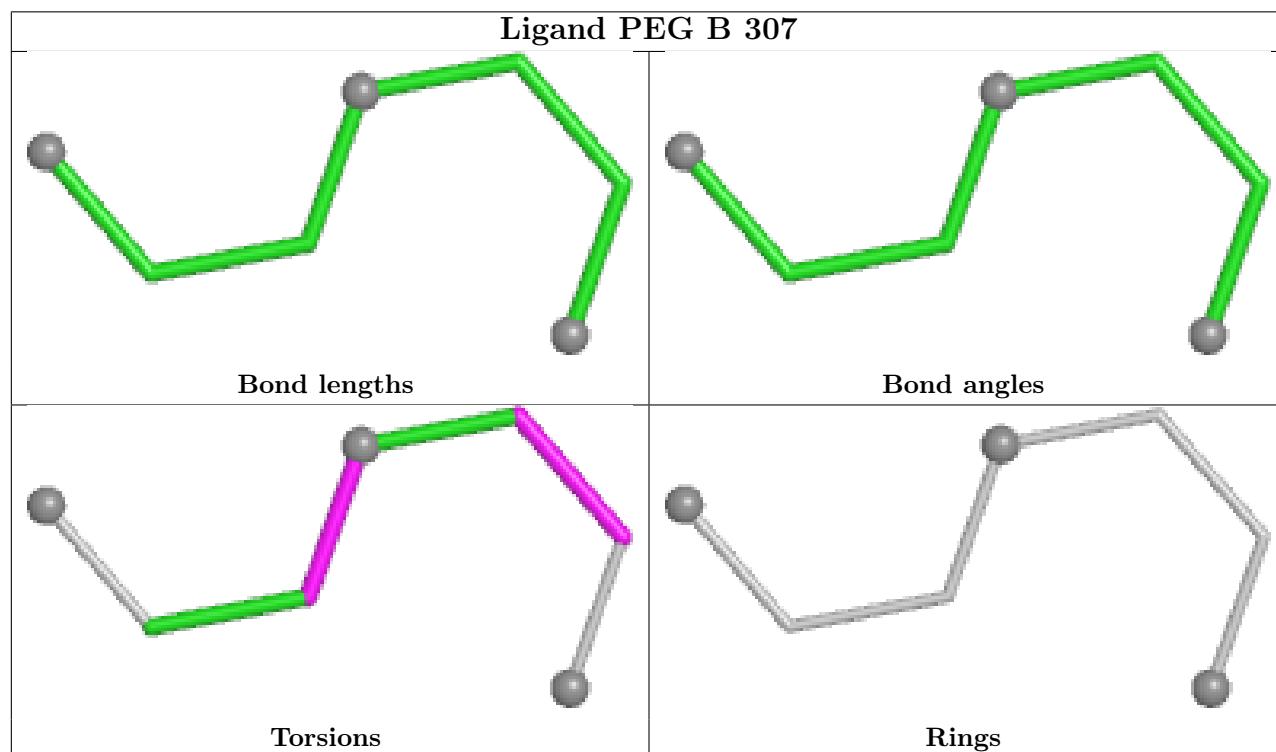
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	307	PEG	4	0
5	B	321	PO4	1	0
2	B	301	A1L2O	2	0
6	B	304	PG4	7	0
3	B	303	PEG	4	0

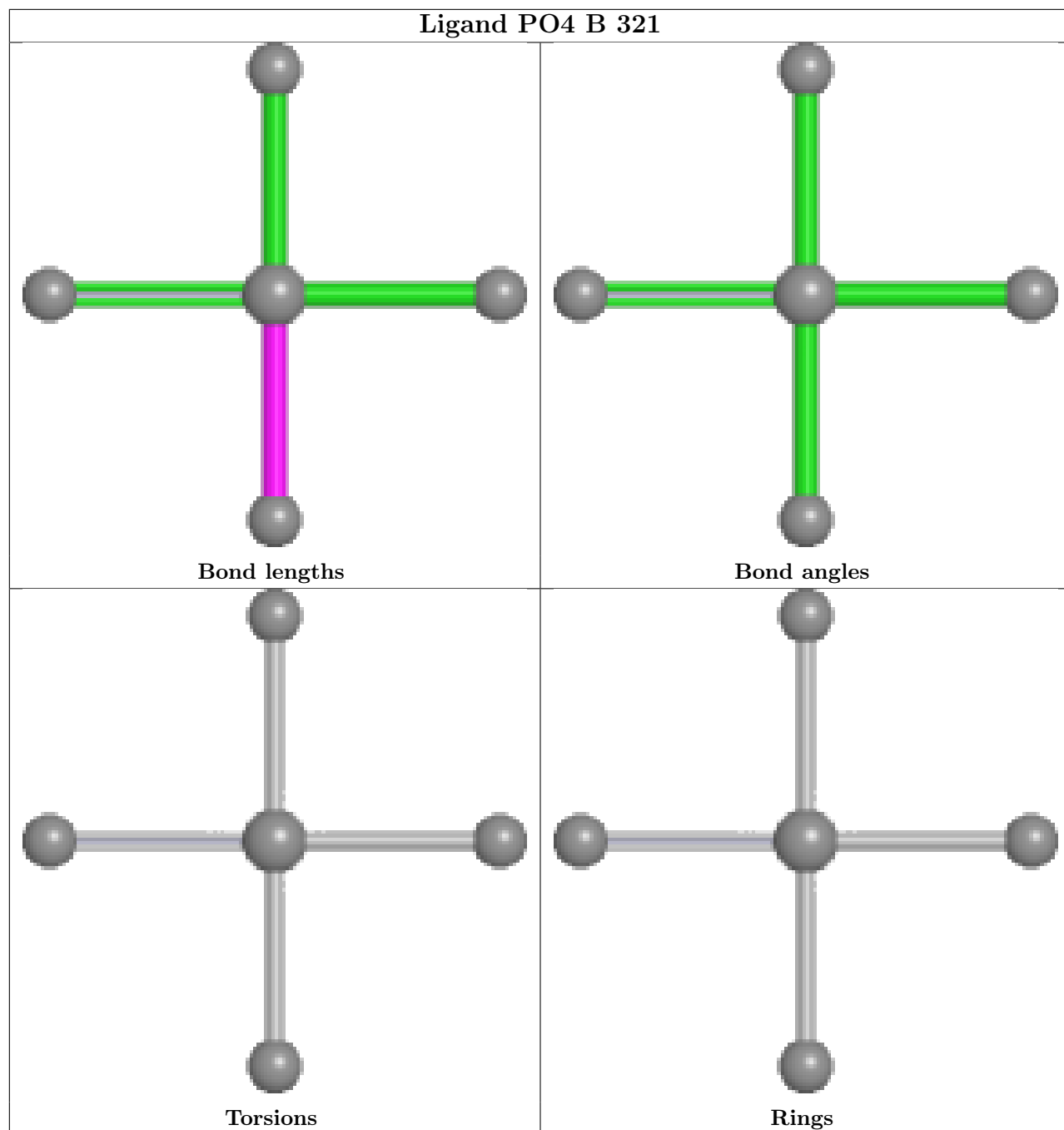
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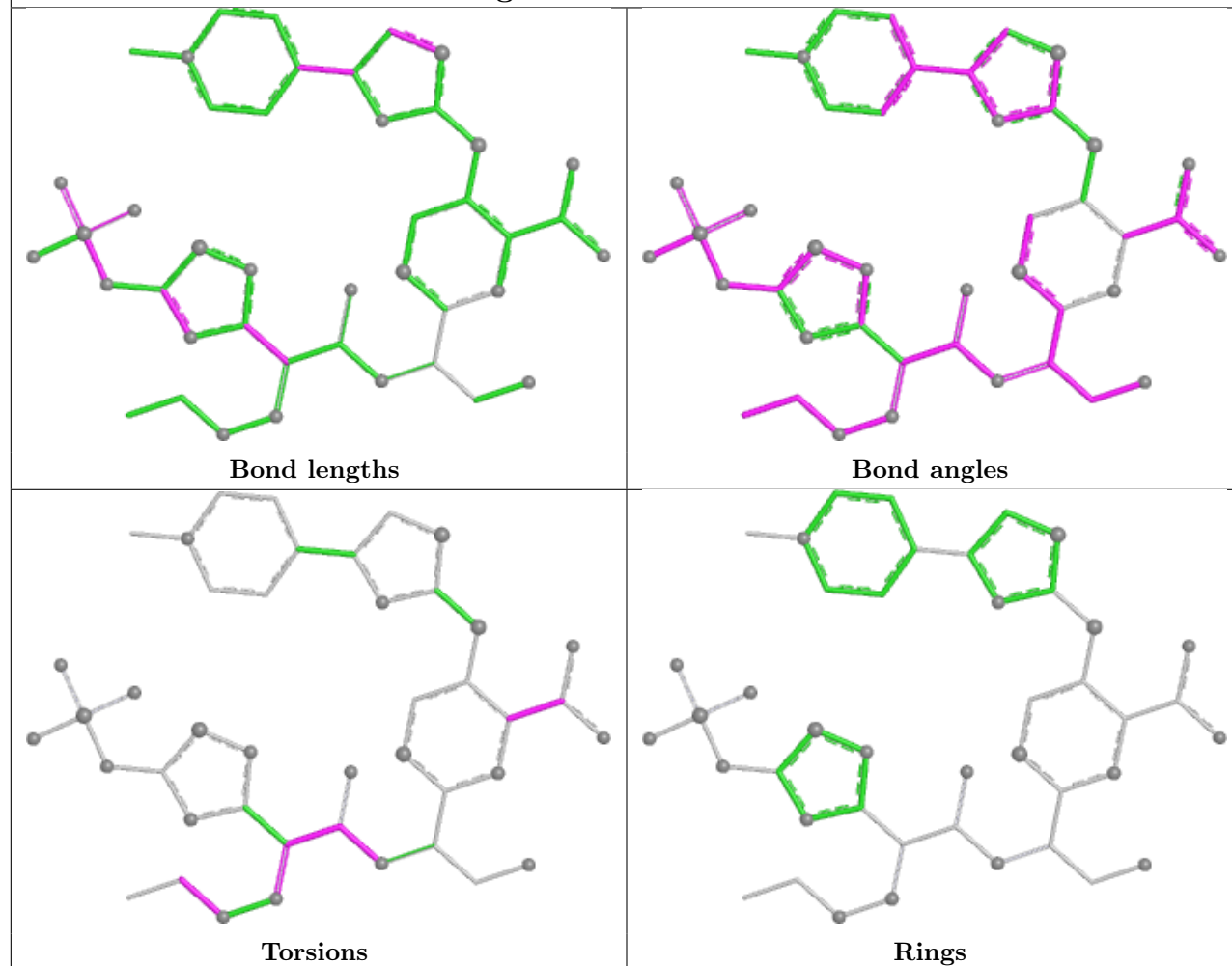
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	302	PEG	1	0
3	B	305	PEG	4	0
2	A	301	A1L2O	1	0

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

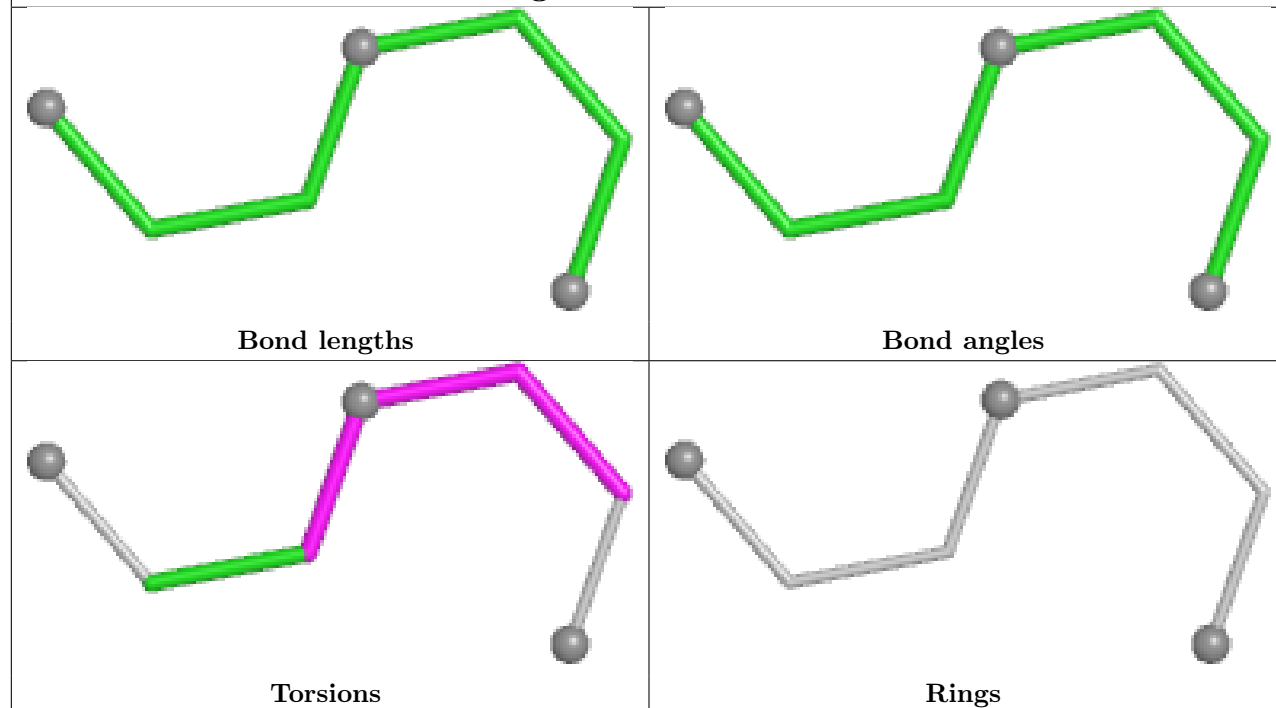


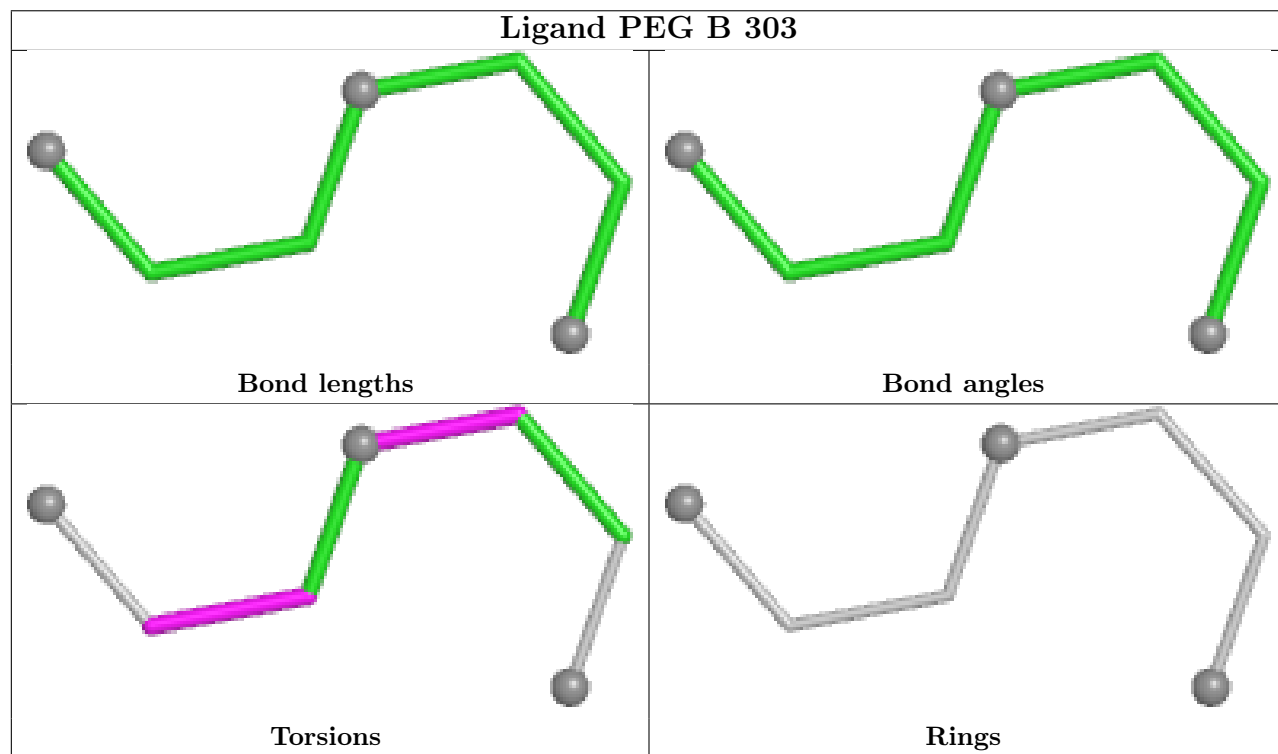
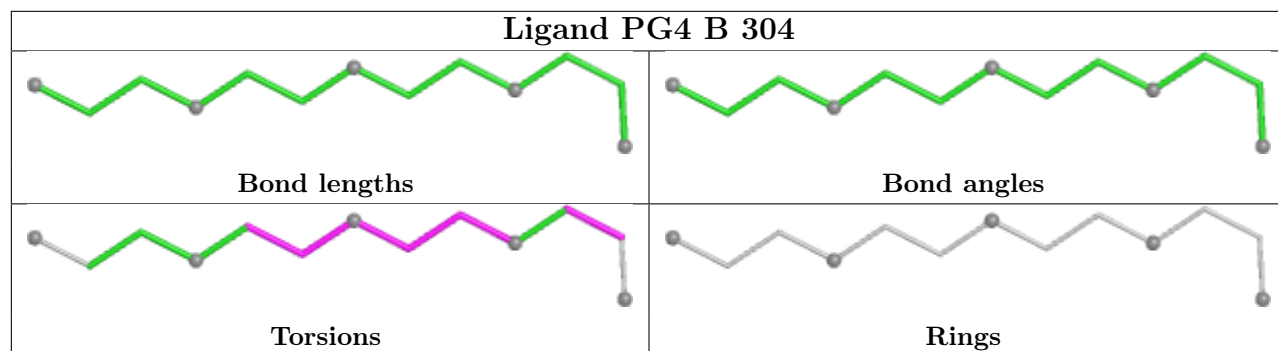


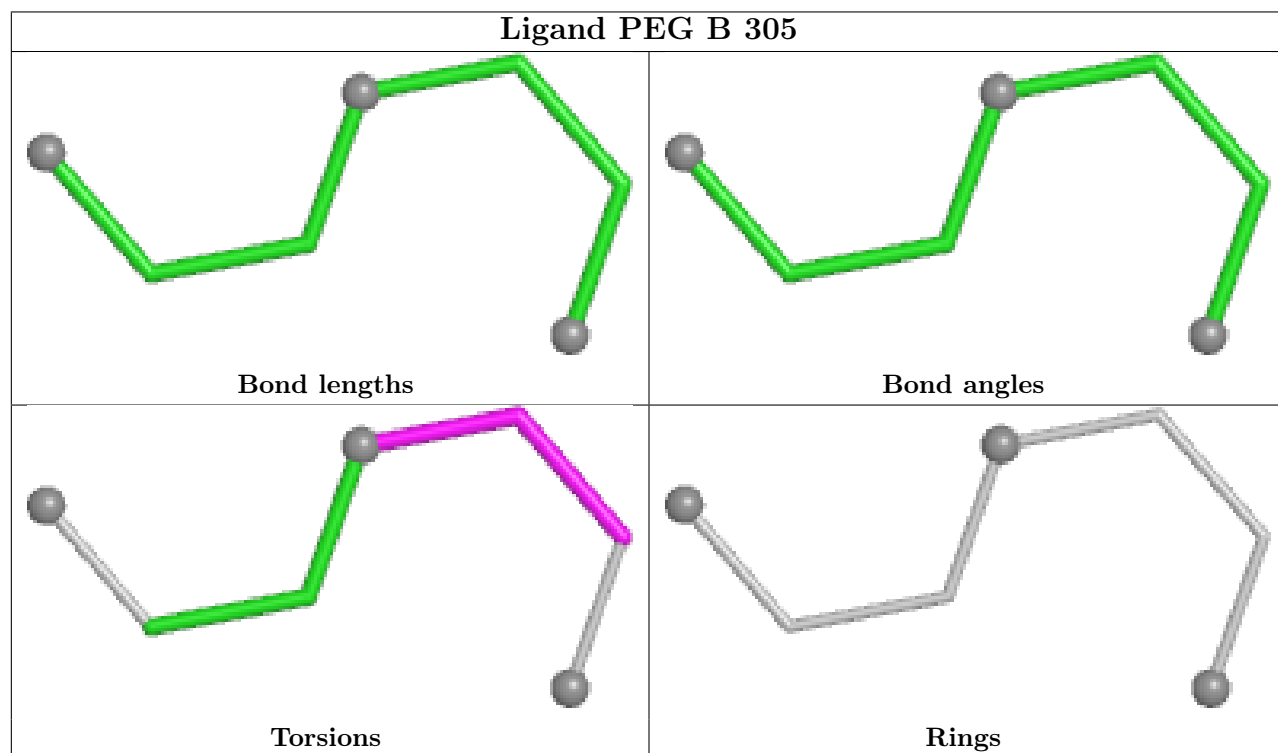
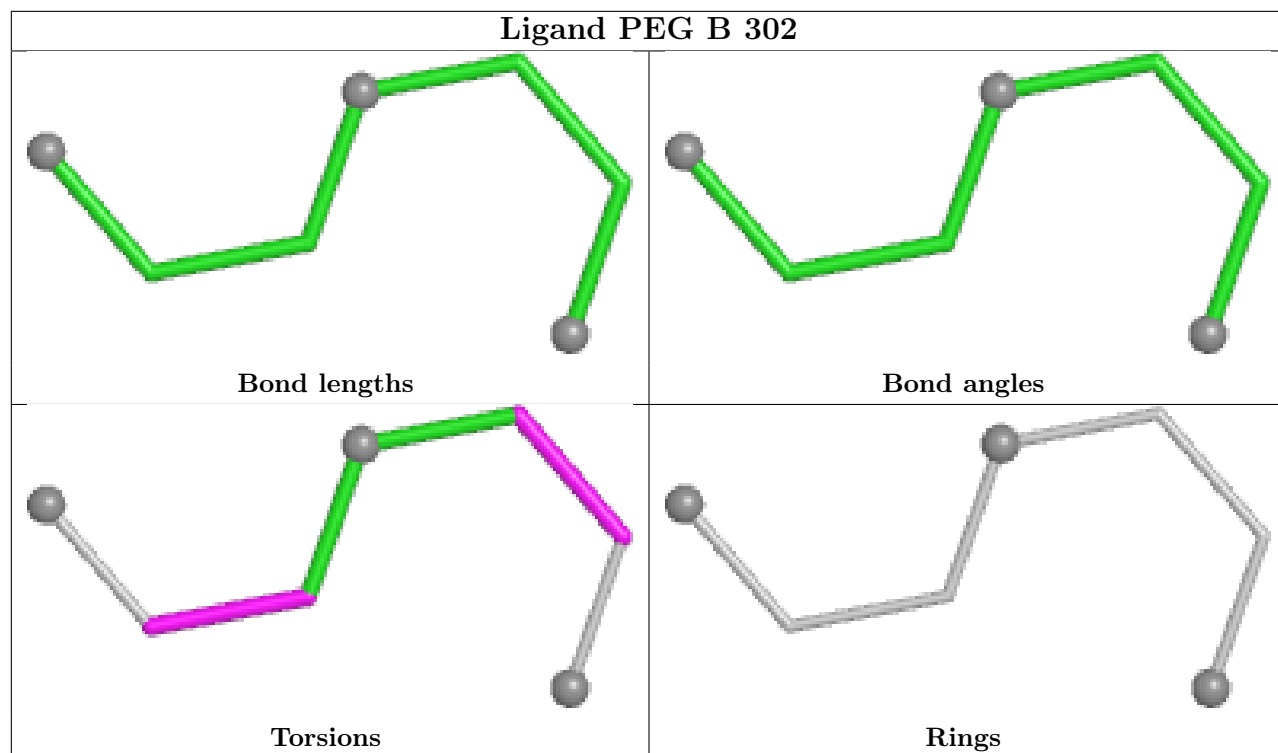
Ligand A1L2O B 301

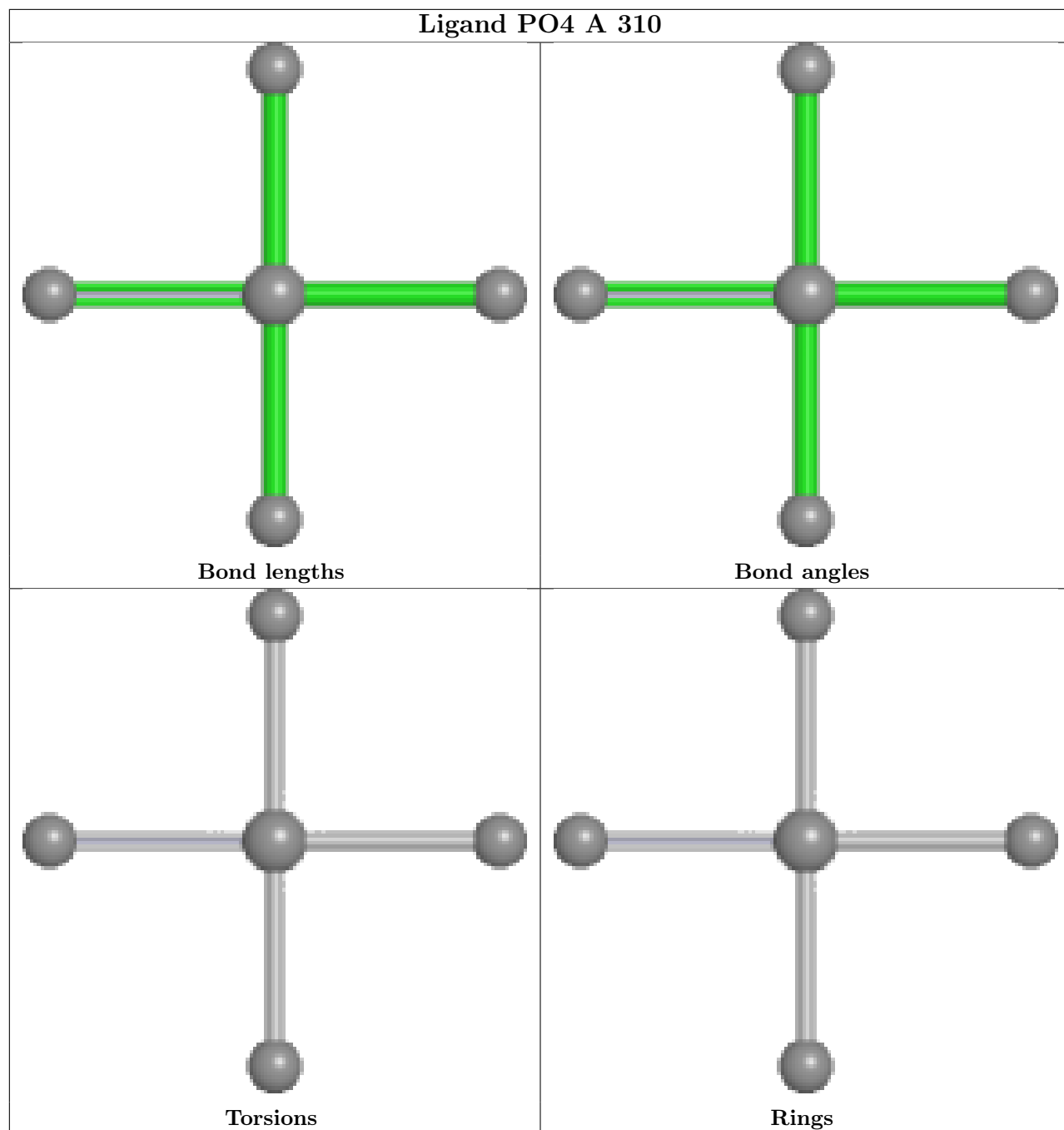


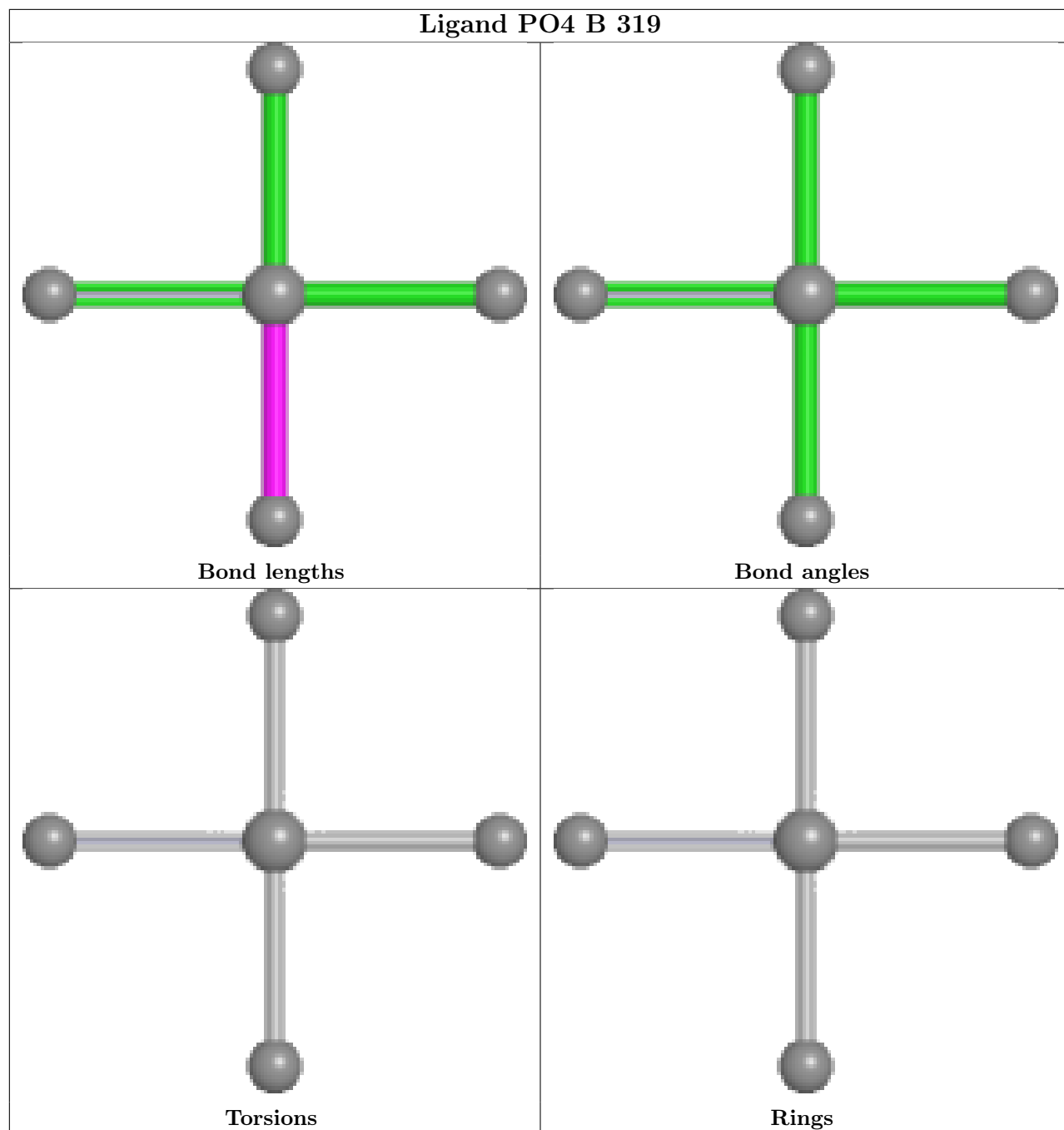
Ligand PEG A 302



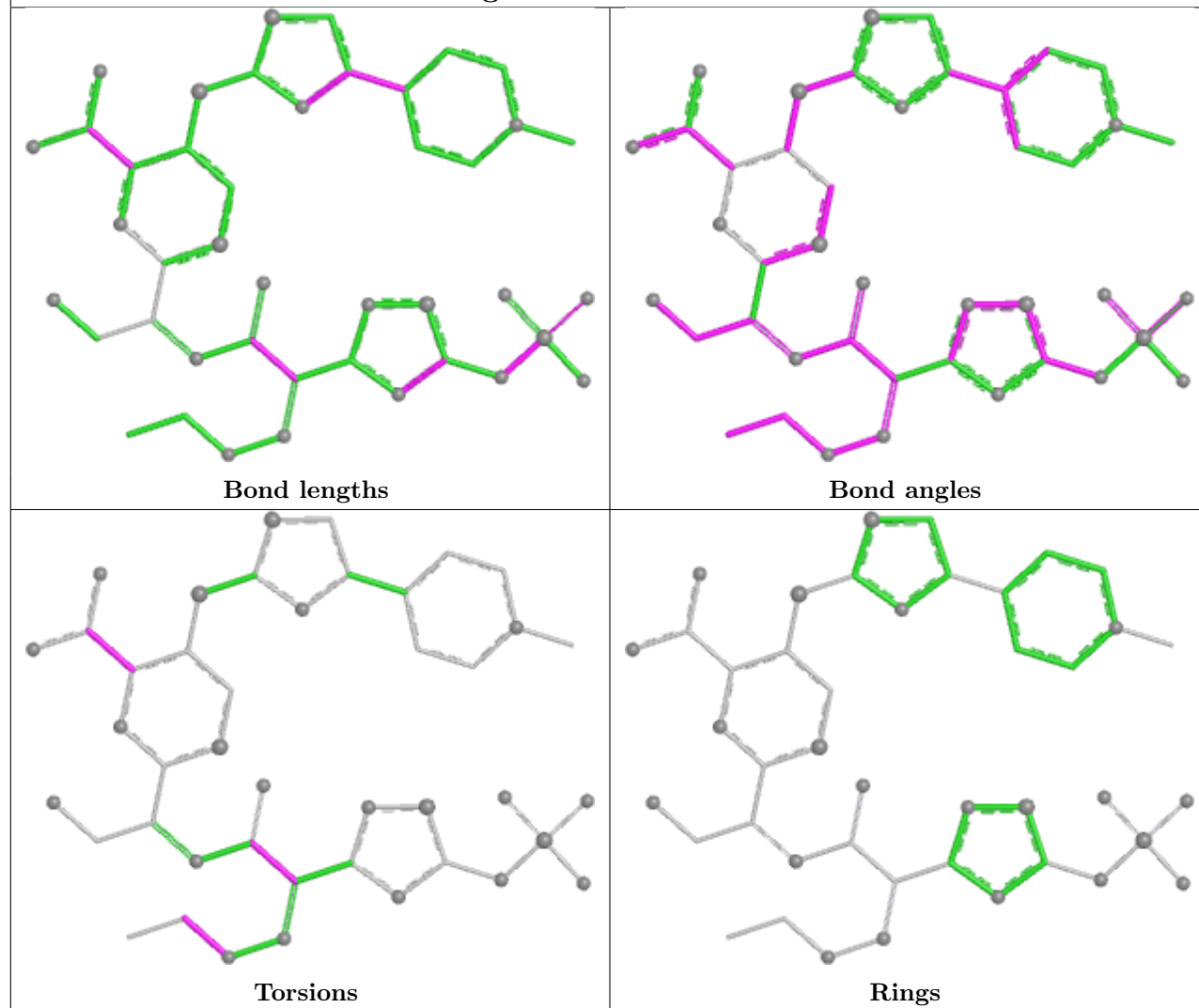


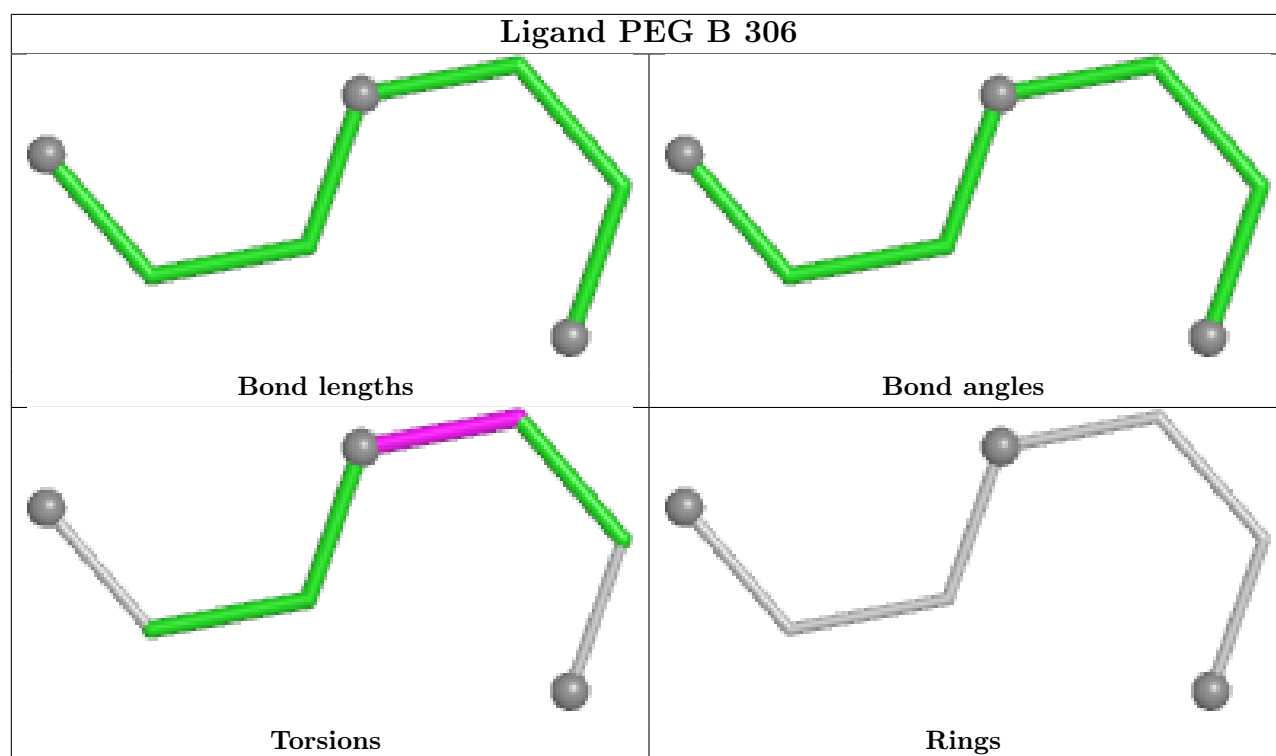


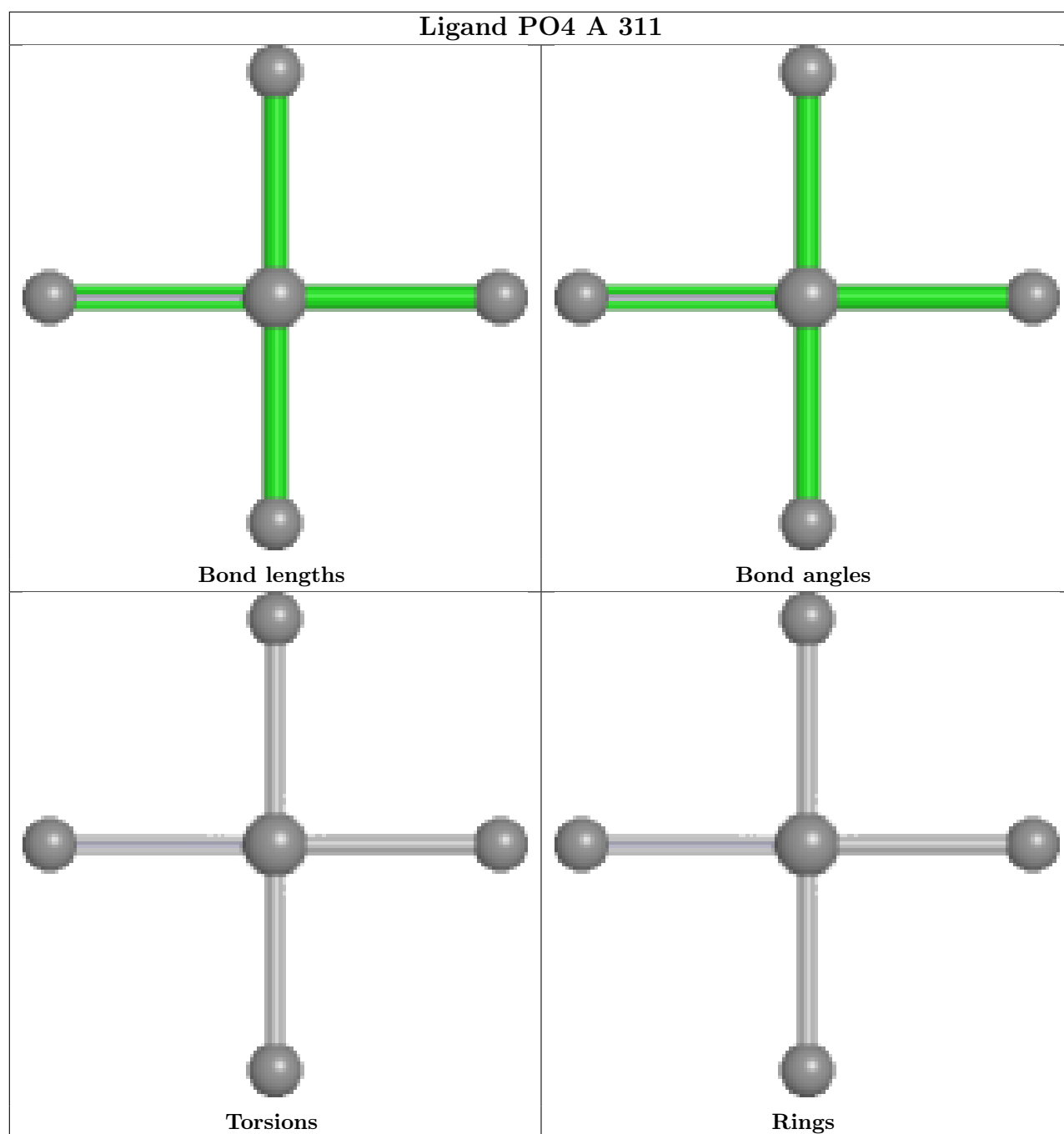


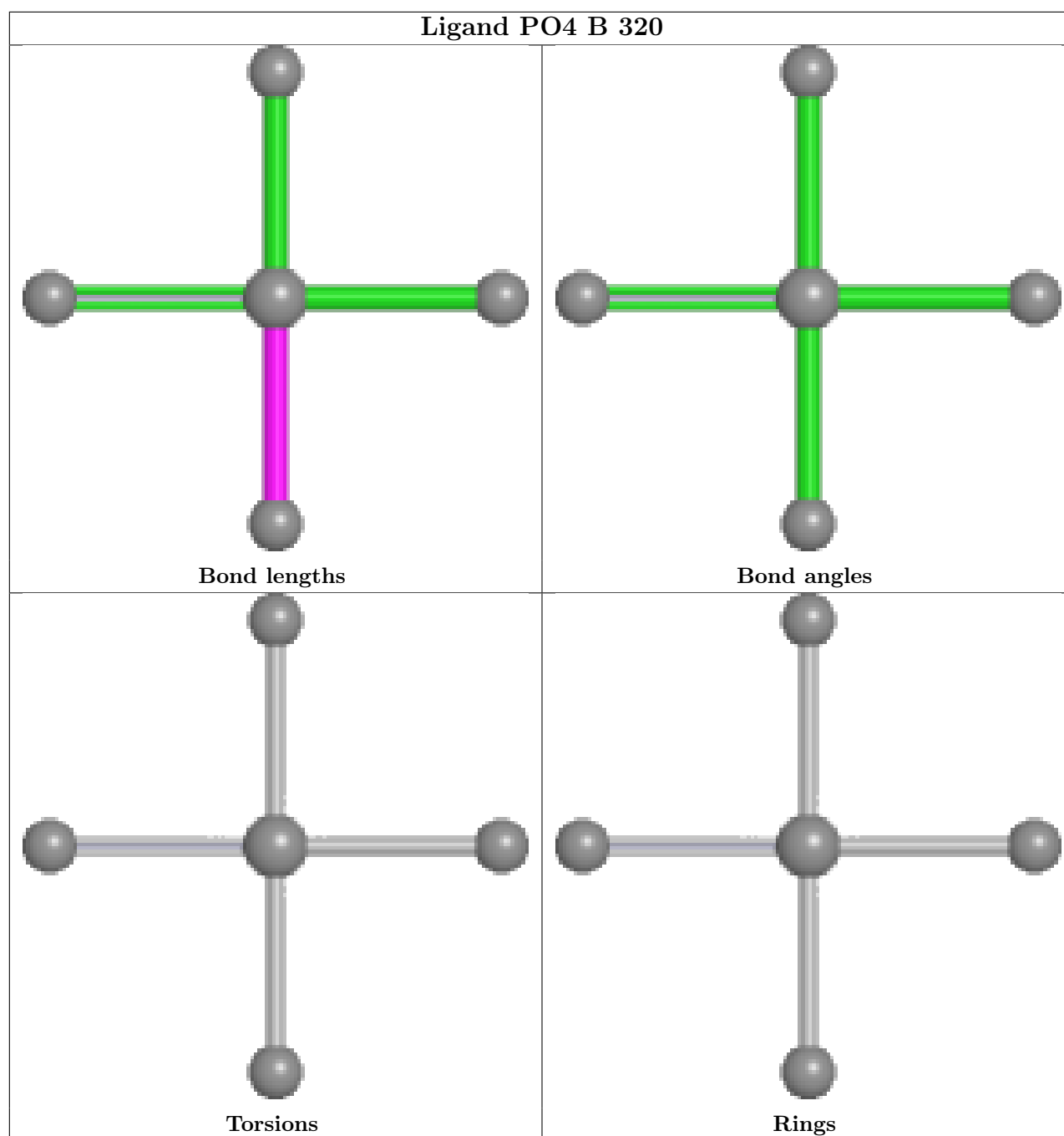


Ligand A1L2O A 301









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	267/275 (97%)	-0.23	1 (0%) 88 86	13, 24, 45, 100	0
1	B	267/275 (97%)	-0.71	1 (0%) 88 86	1, 7, 24, 72	2 (0%)
All	All	534/550 (97%)	-0.47	2 (0%) 88 86	1, 17, 40, 100	2 (0%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	ASN	2.9
1	B	1	ASN	2.1

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [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
5	PO4	A	310	5/5	0.75	0.22	60,73,84,85	0
3	PEG	B	303	7/7	0.78	0.14	47,53,58,62	0
3	PEG	B	305	7/7	0.80	0.15	17,22,24,26	0

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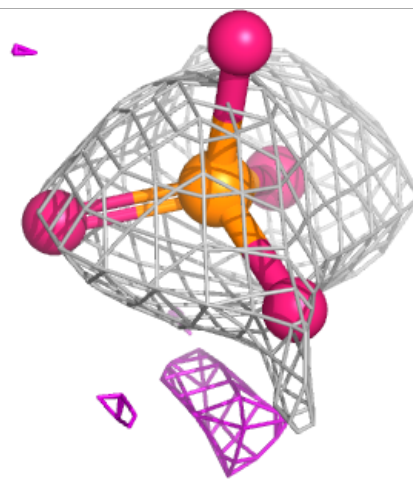
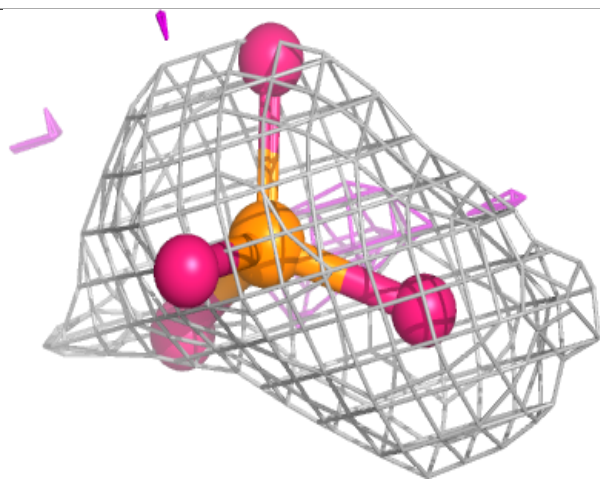
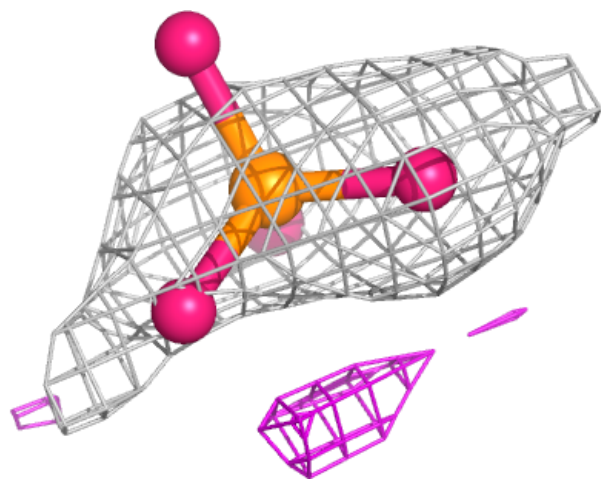
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	PG4	B	304	13/13	0.80	0.16	25,44,50,55	0
3	PEG	B	306	7/7	0.84	0.12	24,29,36,38	0
3	PEG	A	302	7/7	0.86	0.11	36,41,46,47	0
3	PEG	B	302	7/7	0.87	0.13	30,43,51,52	0
3	PEG	B	307	7/7	0.89	0.14	20,24,30,31	0
5	PO4	B	321	5/5	0.89	0.17	53,54,69,74	0
4	CL	A	308	1/1	0.89	0.09	45,45,45,45	0
4	CL	A	303	1/1	0.90	0.15	62,62,62,62	0
4	CL	A	309	1/1	0.91	0.10	48,48,48,48	0
5	PO4	A	311	5/5	0.91	0.15	38,54,59,59	0
7	IOD	B	308	1/1	0.91	0.11	103,103,103,103	0
4	CL	A	307	1/1	0.92	0.06	40,40,40,40	0
2	A1L2O	A	301	43/43	0.92	0.09	22,37,55,61	0
8	NA	B	318	1/1	0.92	0.12	30,30,30,30	0
4	CL	A	305	1/1	0.93	0.08	54,54,54,54	0
4	CL	B	317	1/1	0.93	0.07	54,54,54,54	0
4	CL	B	310	1/1	0.94	0.06	34,34,34,34	0
5	PO4	B	320	5/5	0.94	0.16	29,31,40,40	0
4	CL	B	313	1/1	0.94	0.07	37,37,37,37	0
2	A1L2O	B	301	43/43	0.95	0.08	6,16,32,37	0
4	CL	B	314	1/1	0.95	0.05	22,22,22,22	0
5	PO4	B	319	5/5	0.95	0.11	24,35,39,40	0
4	CL	B	312	1/1	0.95	0.08	46,46,46,46	0
4	CL	A	304	1/1	0.96	0.07	44,44,44,44	0
4	CL	A	306	1/1	0.97	0.08	39,39,39,39	0
4	CL	B	315	1/1	0.97	0.04	29,29,29,29	0
4	CL	B	316	1/1	0.97	0.13	39,39,39,39	0
4	CL	B	311	1/1	0.97	0.05	15,15,15,15	0
4	CL	B	309	1/1	0.98	0.05	33,33,33,33	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.

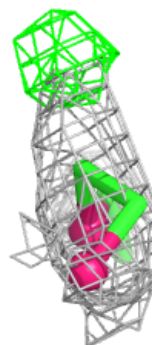
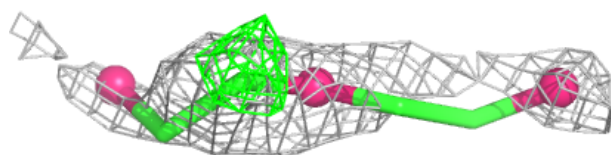
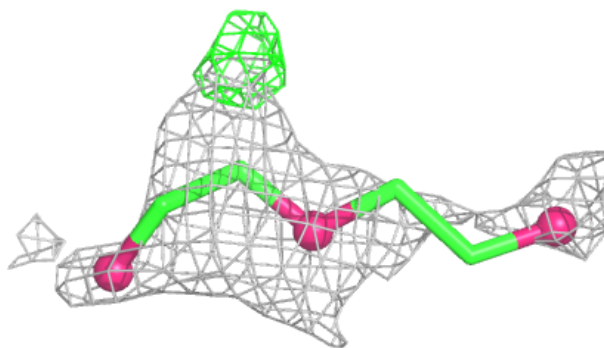
Electron density around PO4 A 310:

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

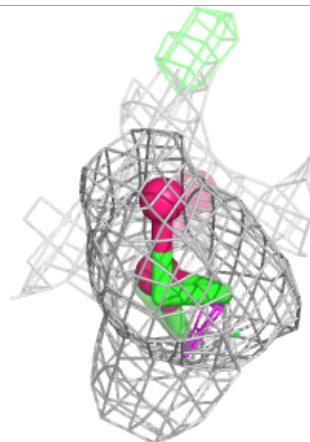
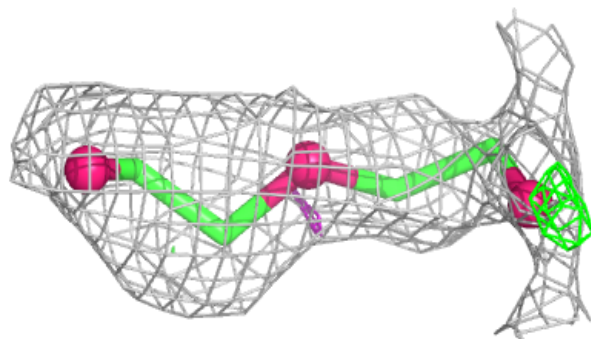
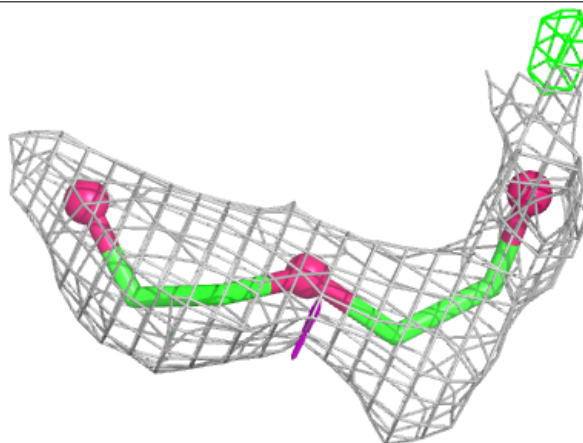


Electron density around PEG B 303:

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

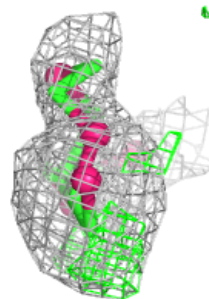
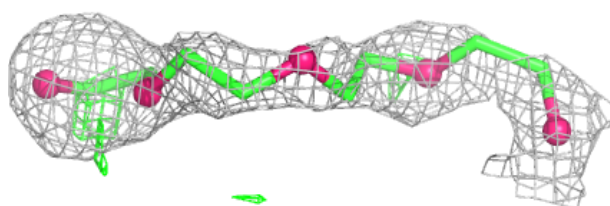
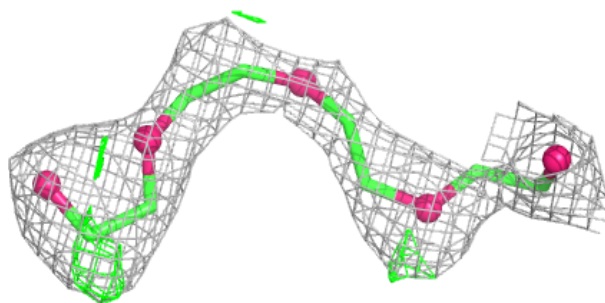
**Electron density around PEG B 305:**

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

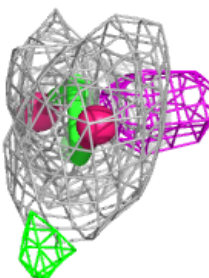
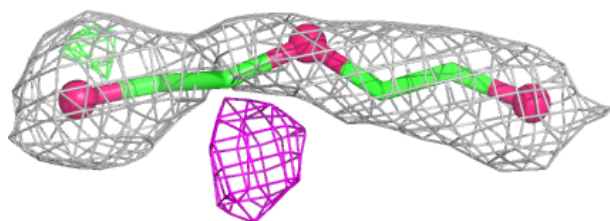
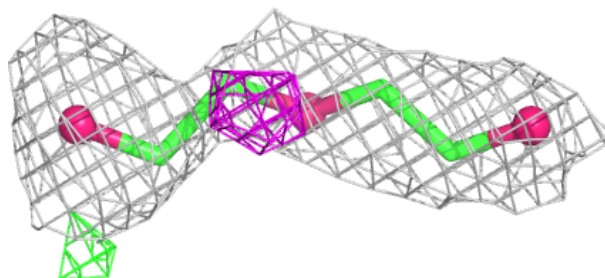


Electron density around PG4 B 304:

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

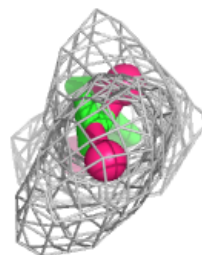
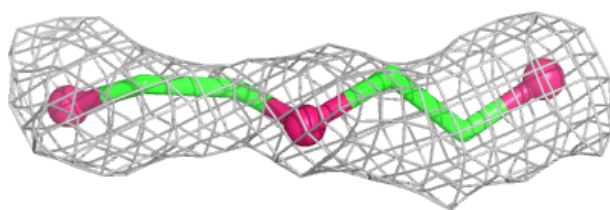
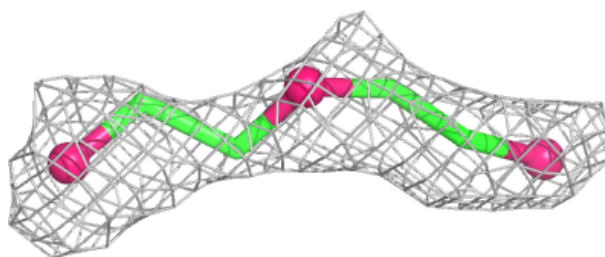
**Electron density around PEG B 306:**

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

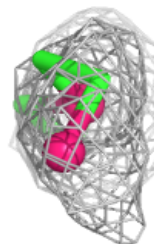
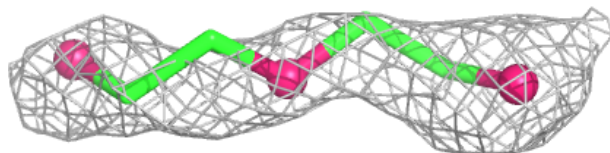
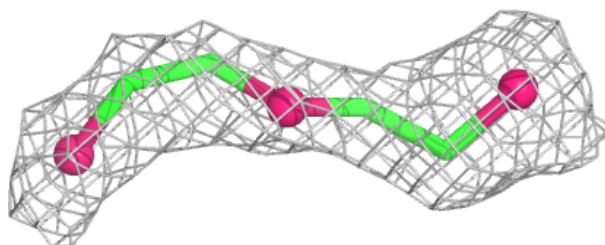


Electron density around PEG A 302:

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

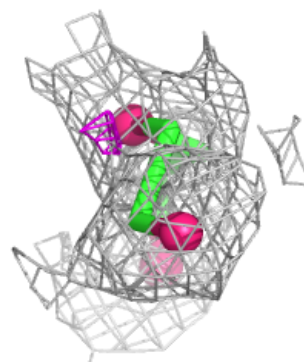
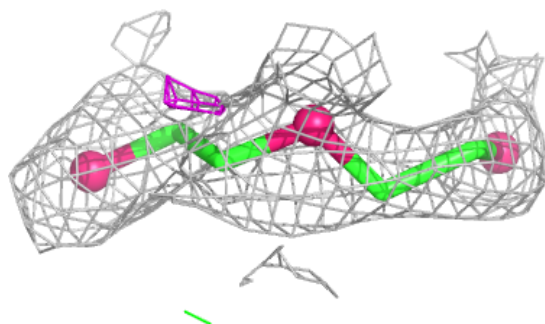
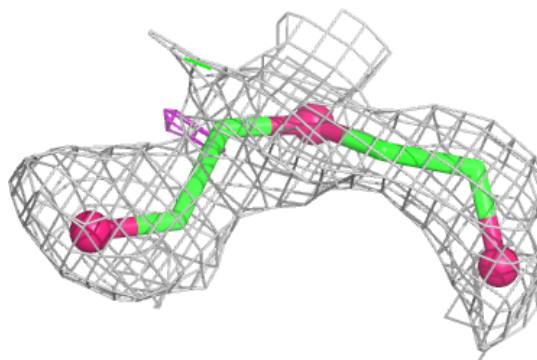
**Electron density around PEG B 302:**

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



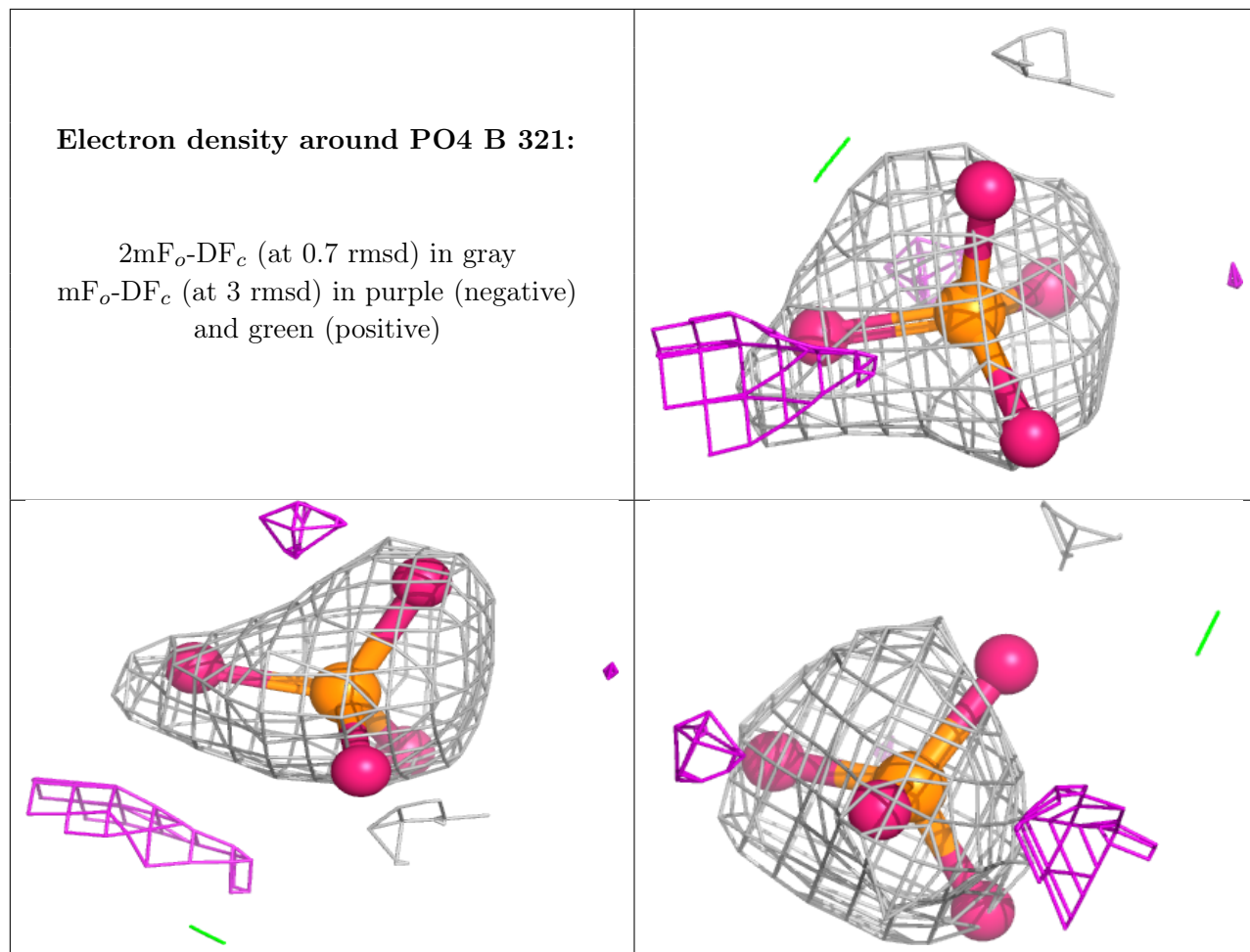
Electron density around PEG B 307:

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



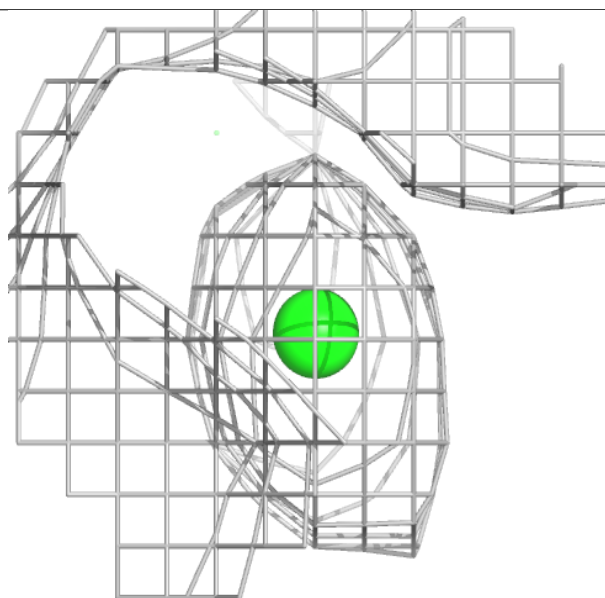
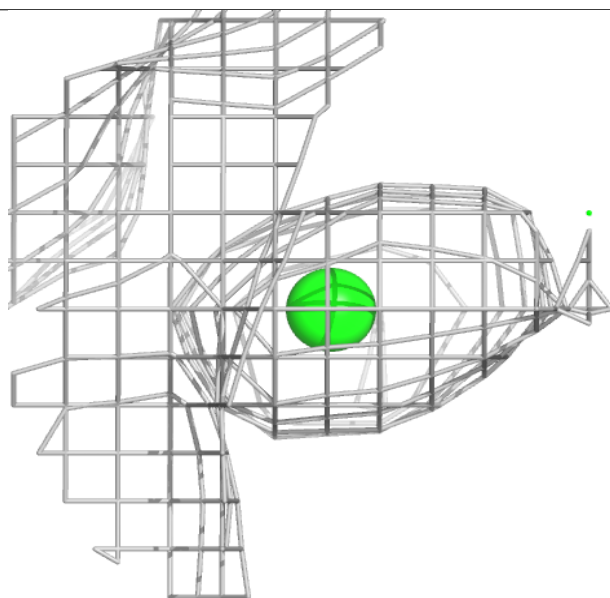
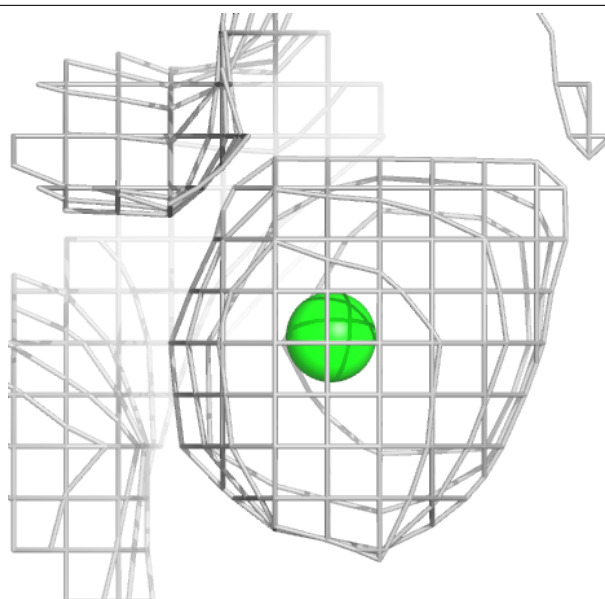
Electron density around PO4 B 321:

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



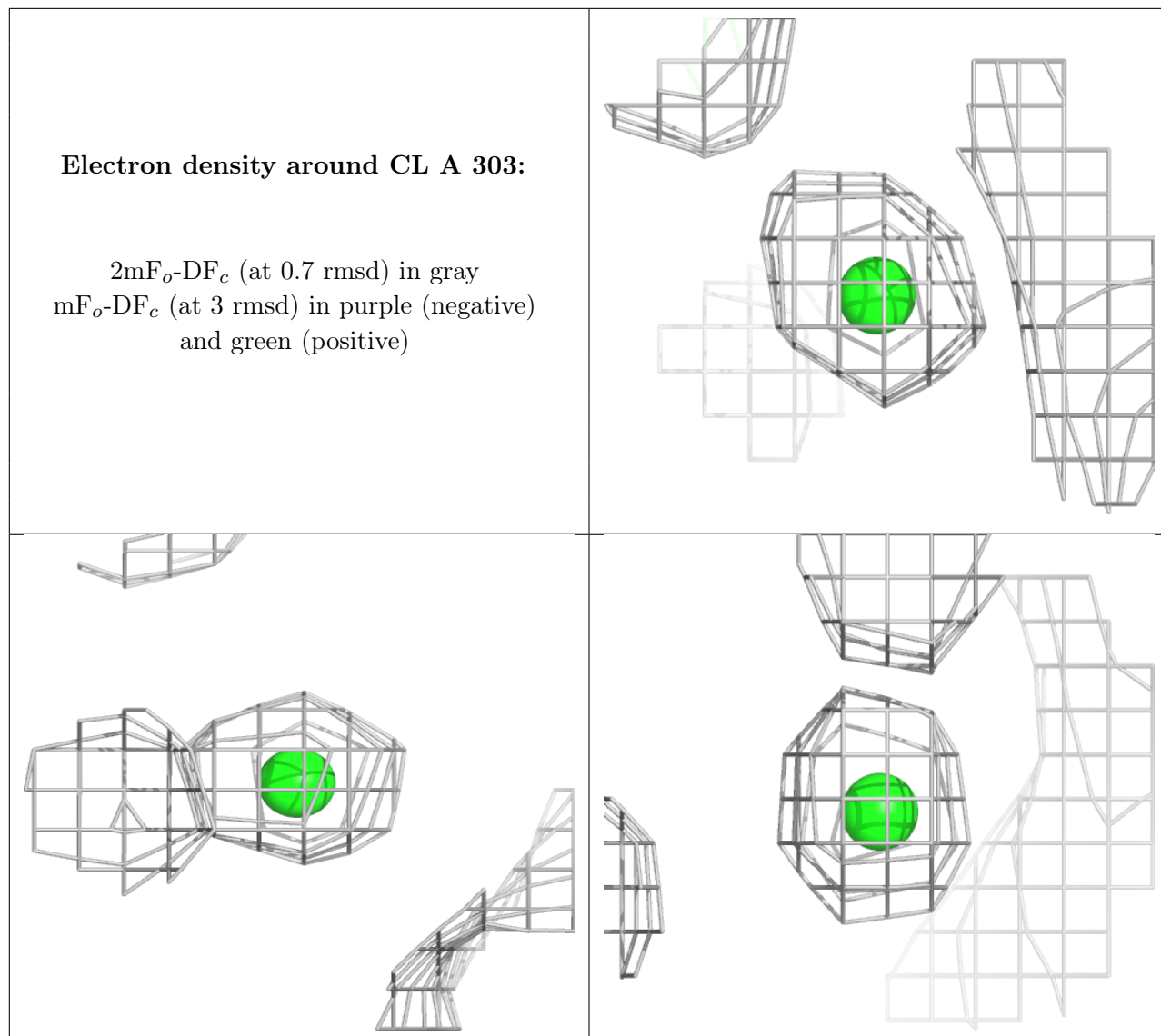
Electron density around CL A 308:

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



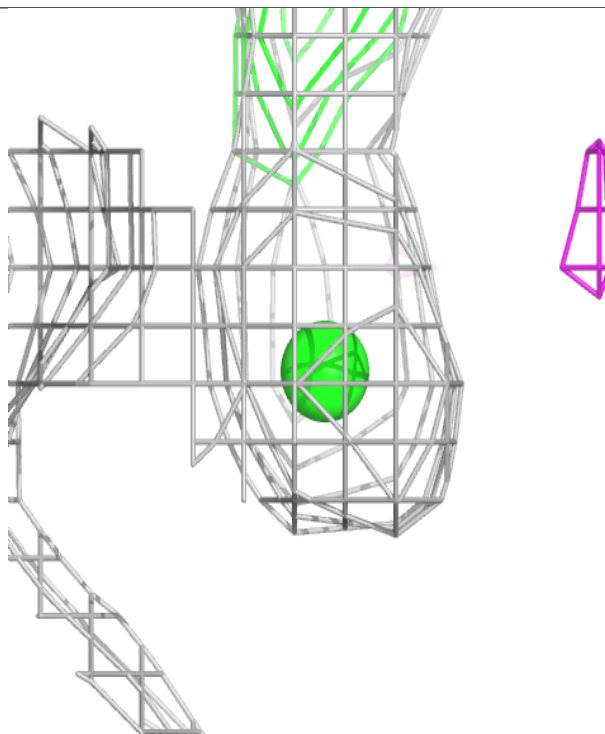
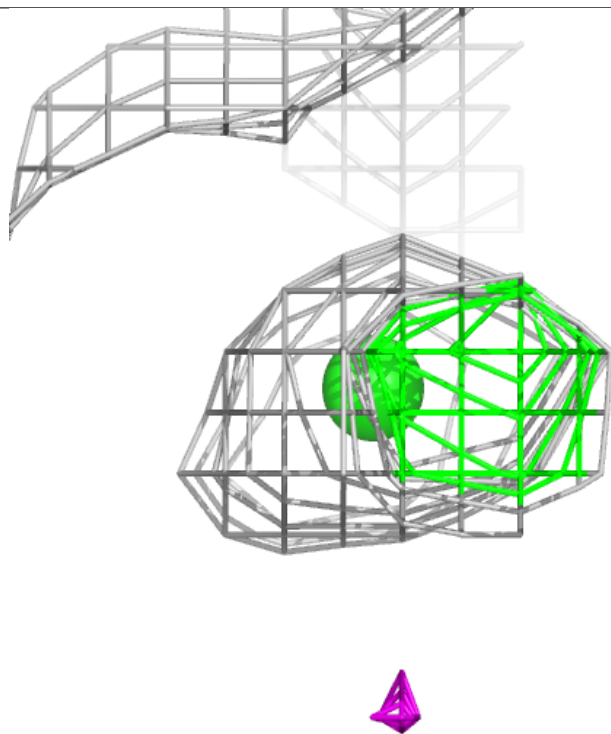
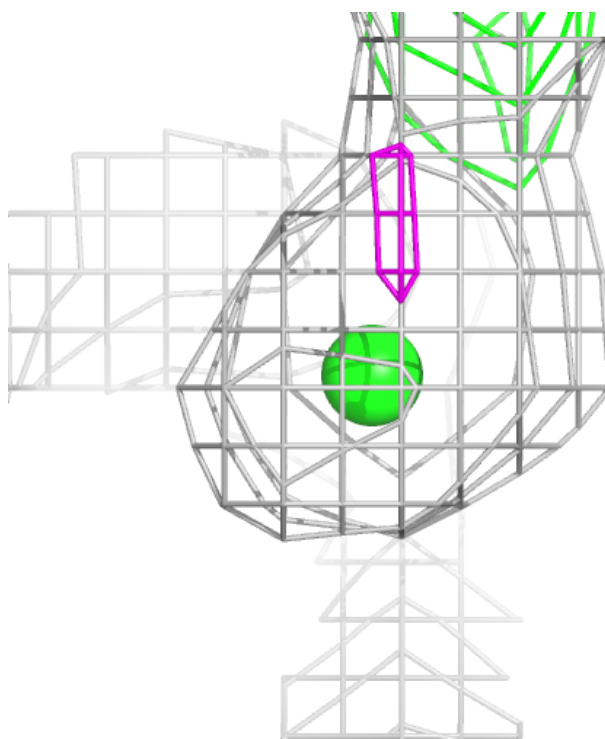
Electron density around CL A 303:

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



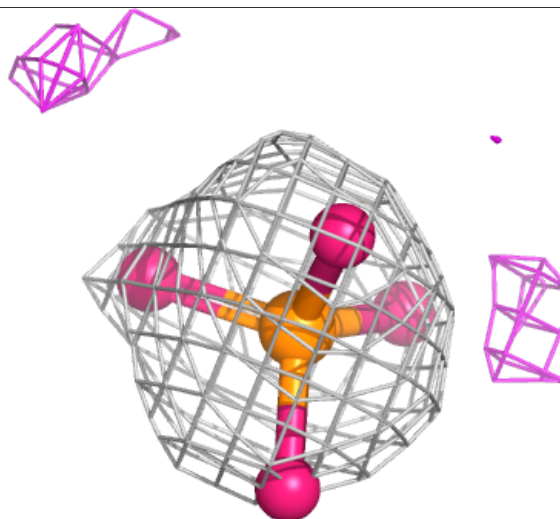
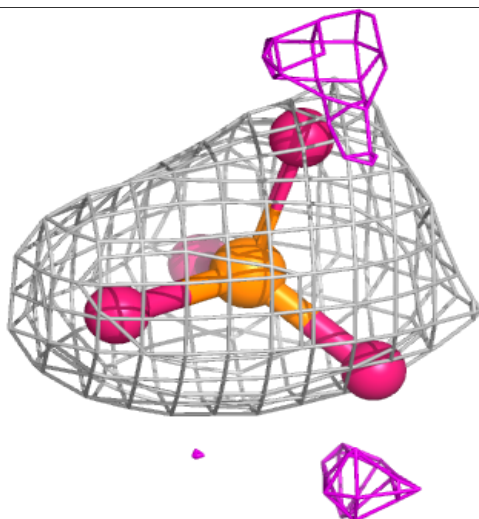
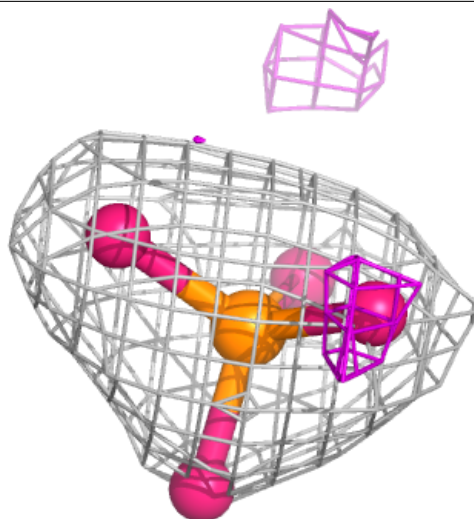
Electron density around CL A 309:

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



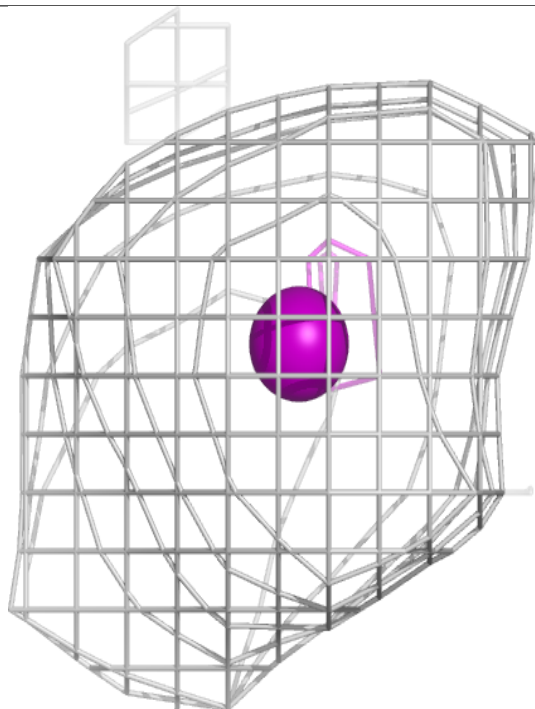
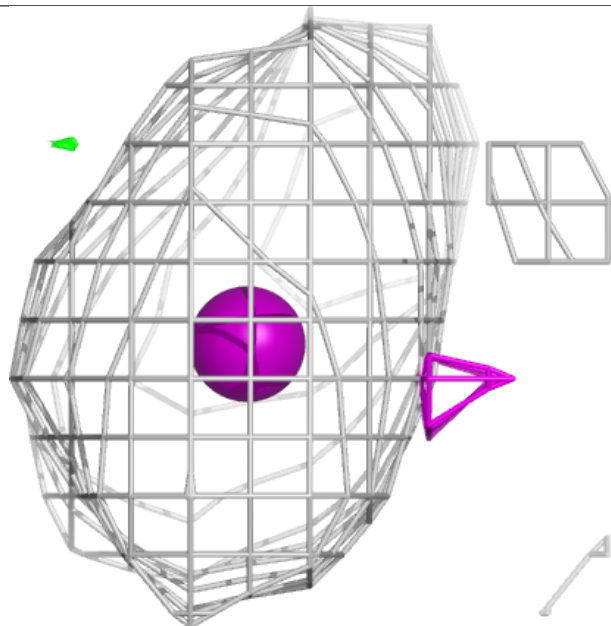
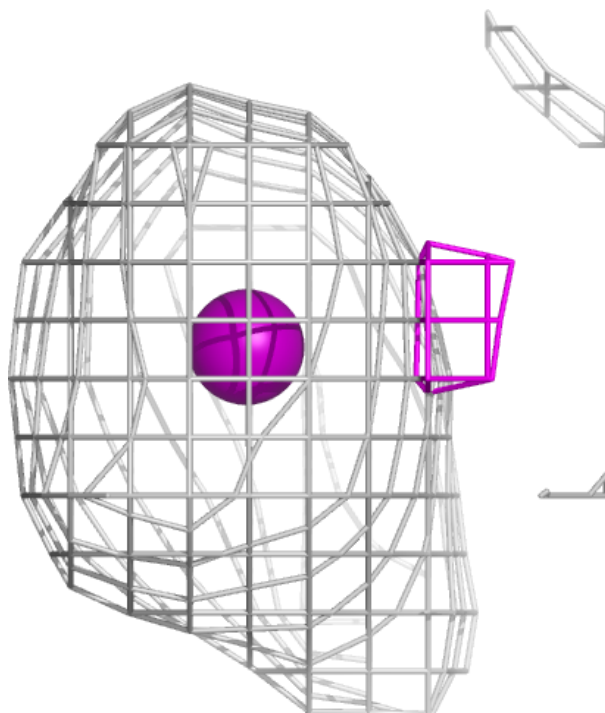
Electron density around PO4 A 311:

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



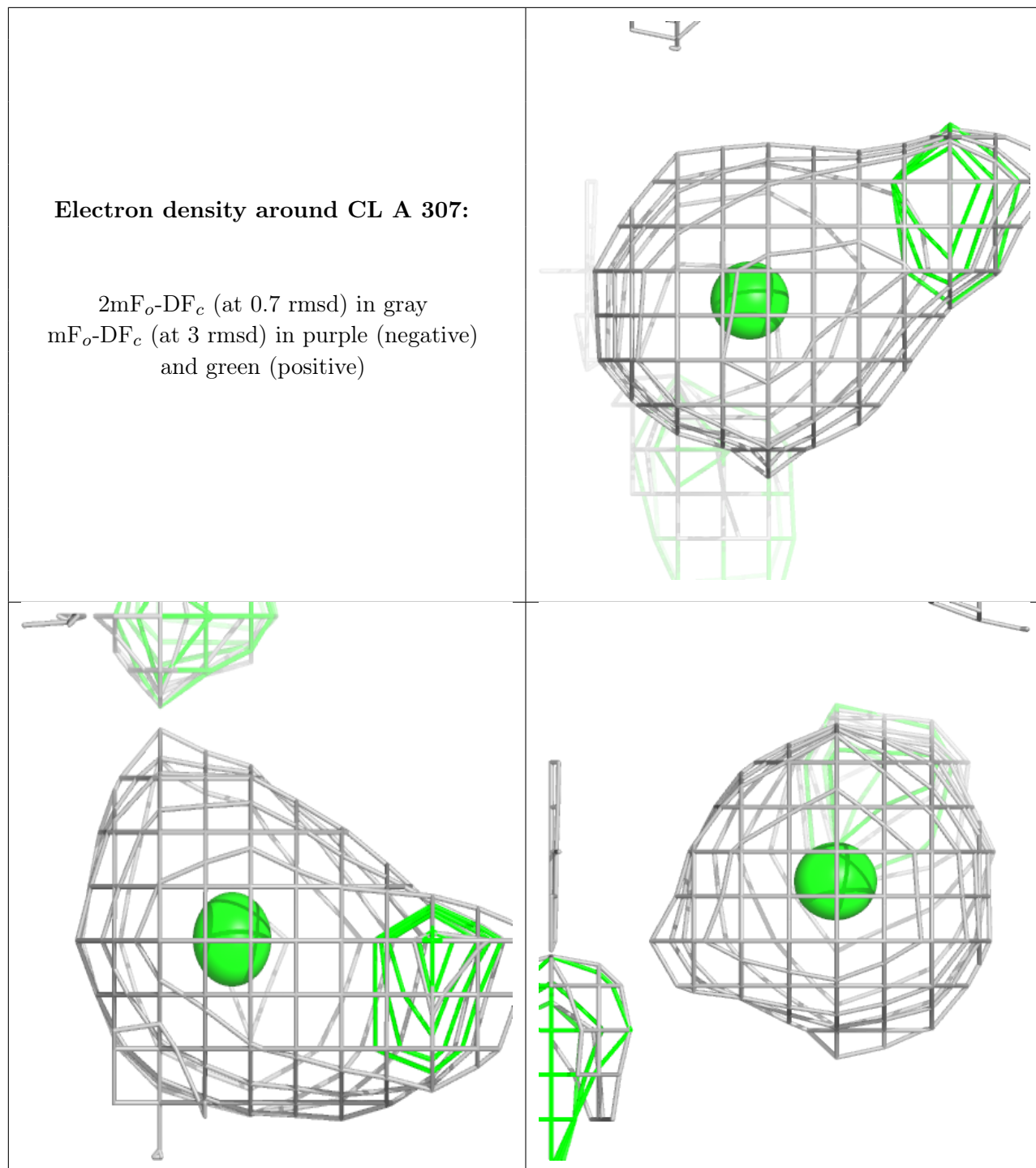
Electron density around IOD B 308:

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



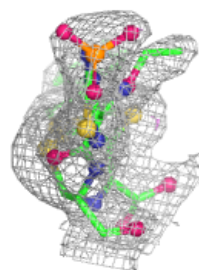
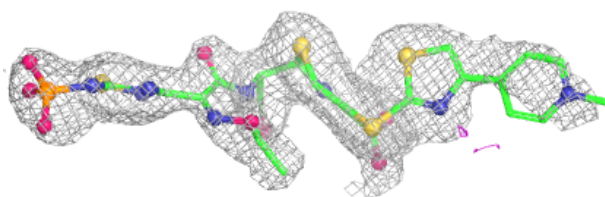
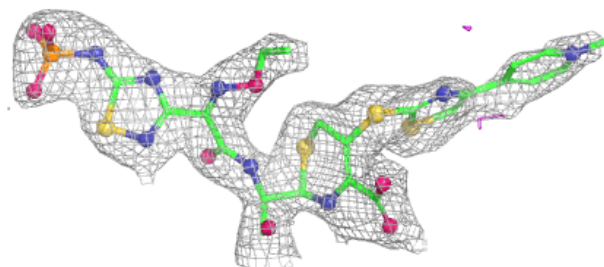
Electron density around CL A 307:

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



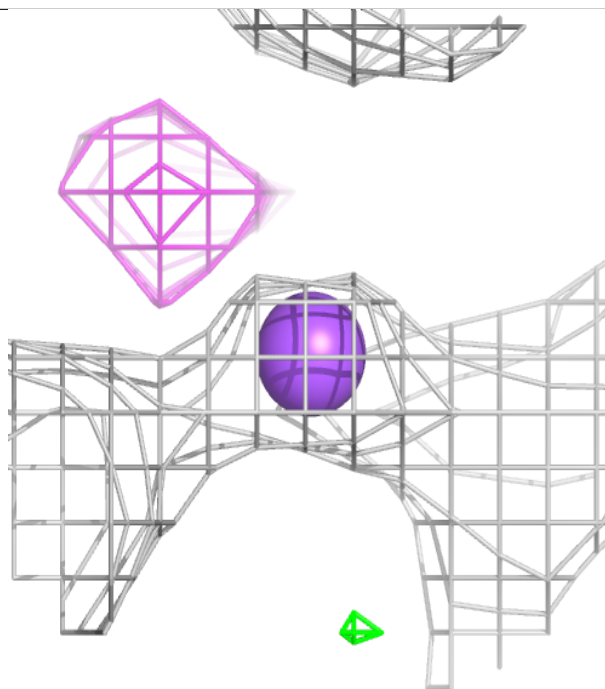
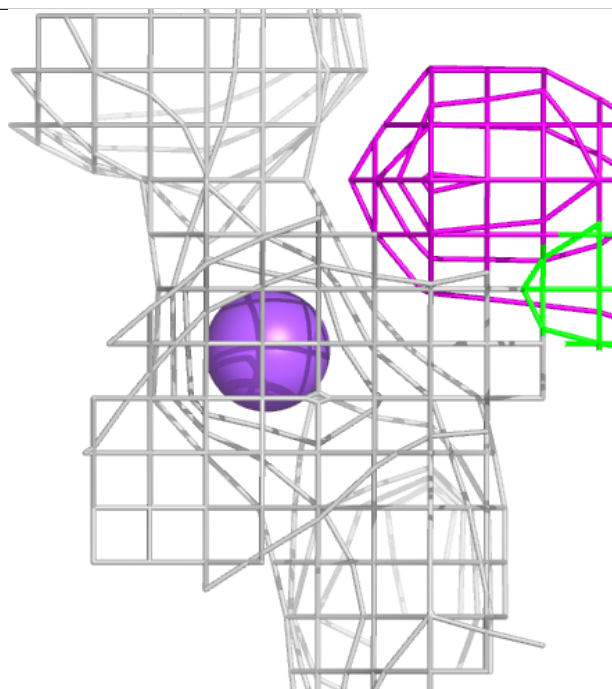
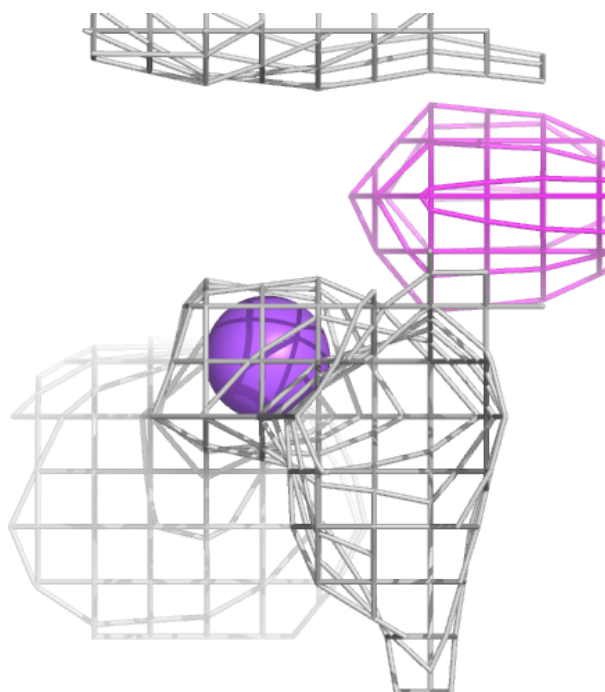
Electron density around A1L2O A 301:

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



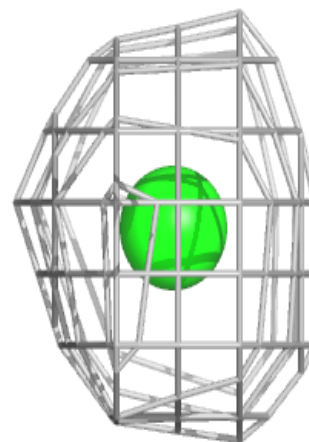
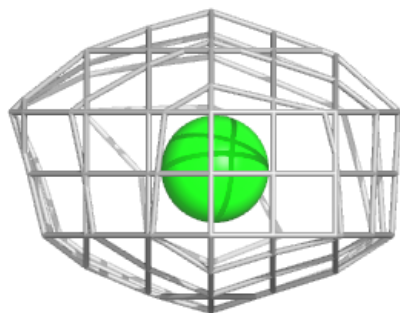
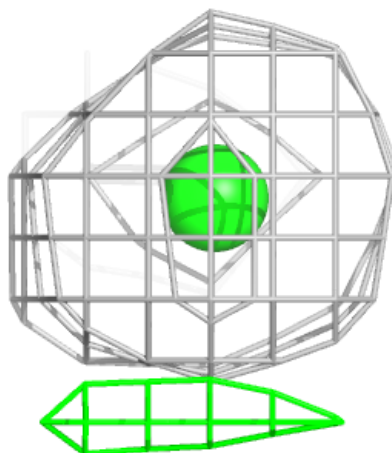
Electron density around NA B 318:

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



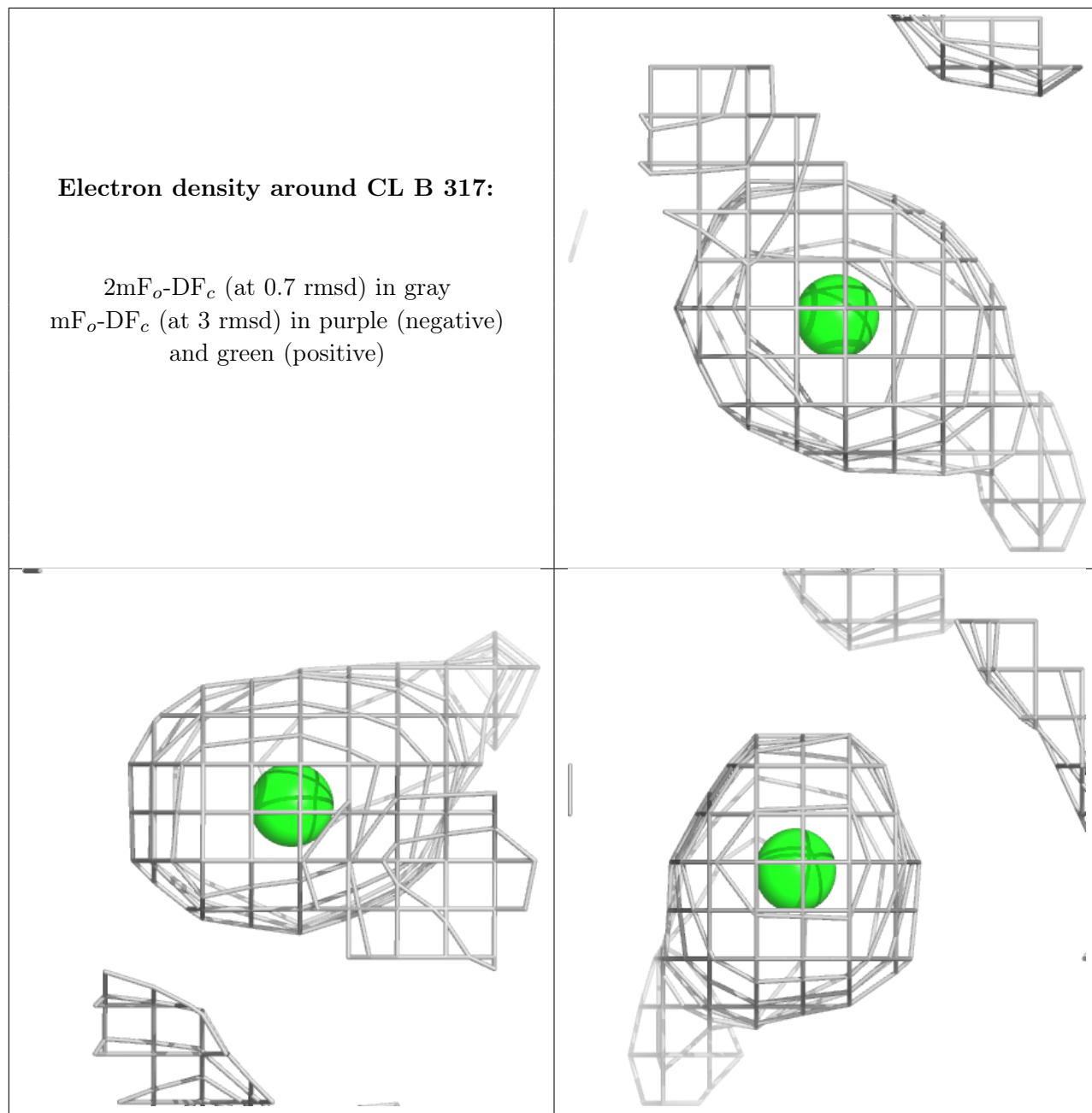
Electron density around CL A 305:

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



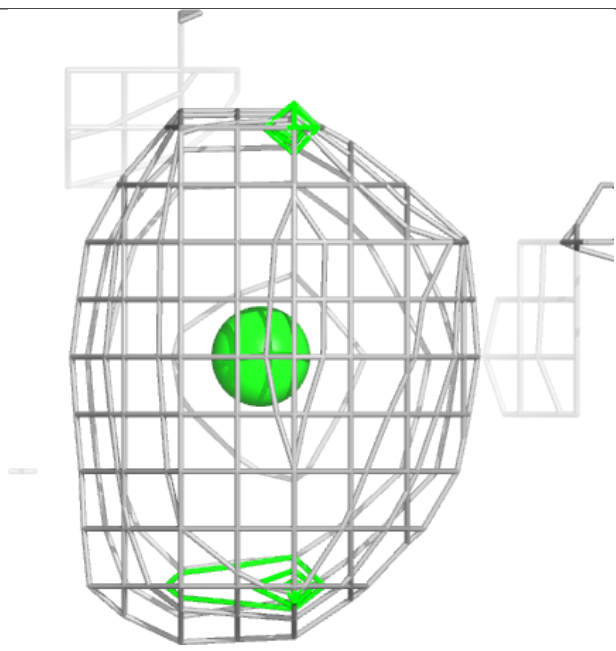
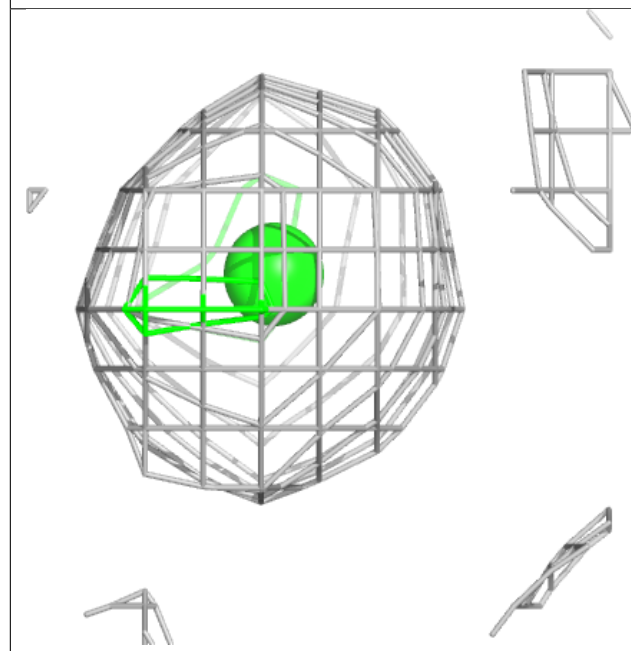
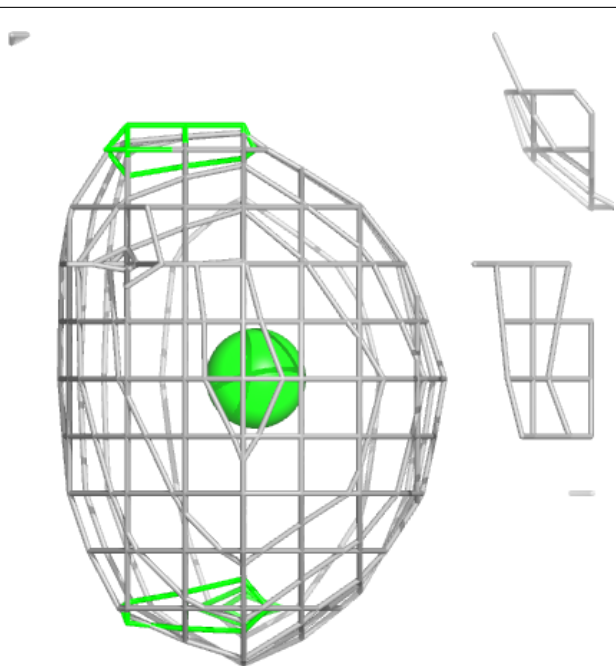
Electron density around CL B 317:

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



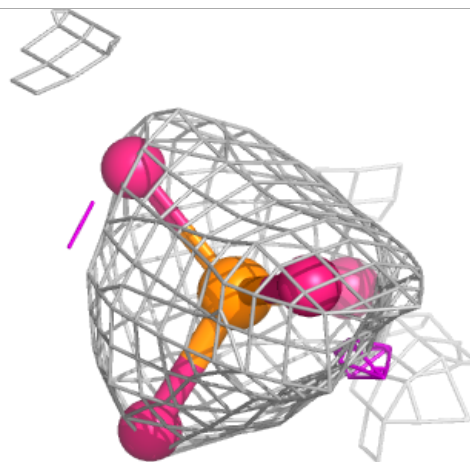
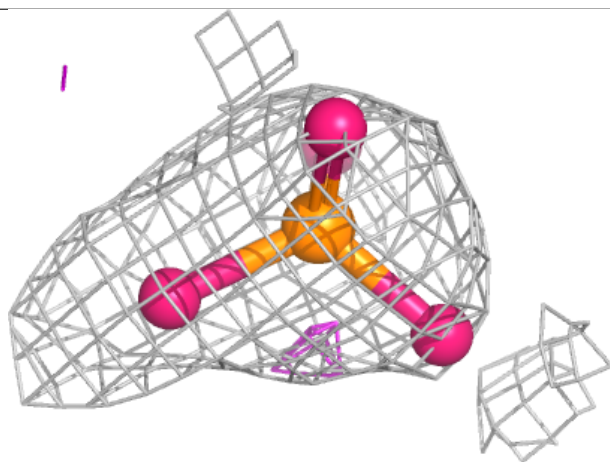
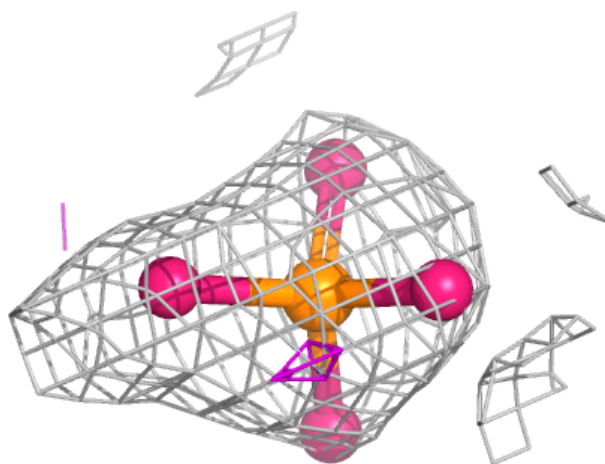
Electron density around CL B 310:

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



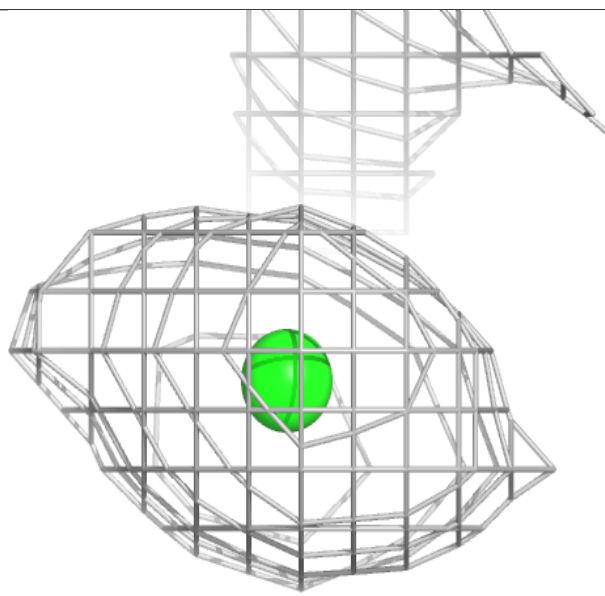
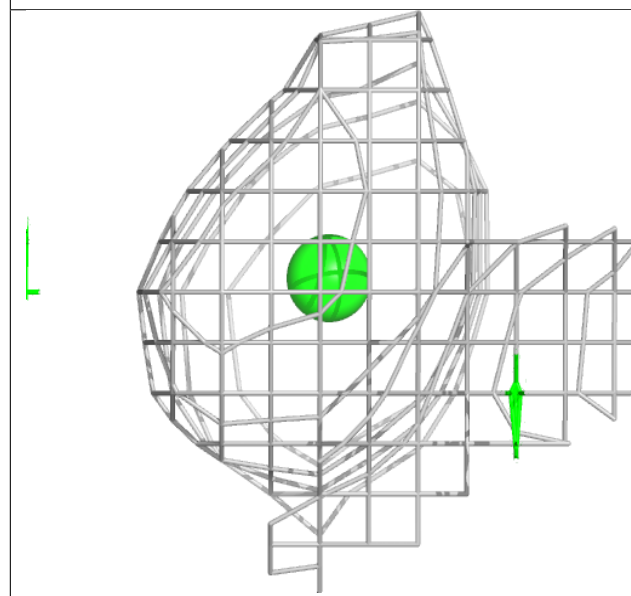
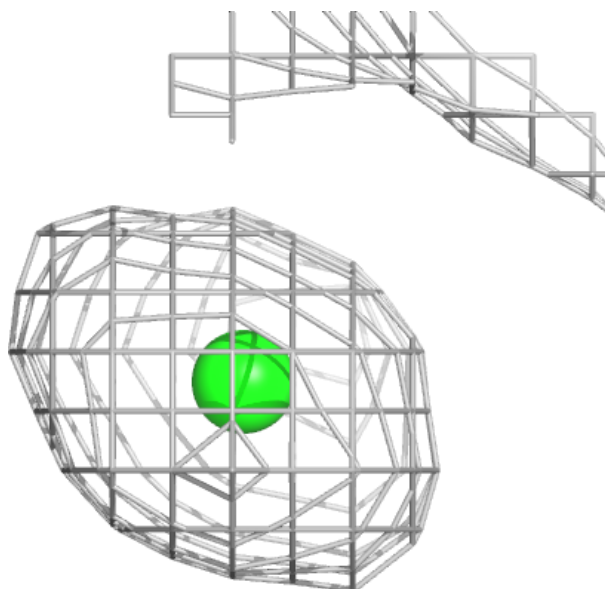
Electron density around PO4 B 320:

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



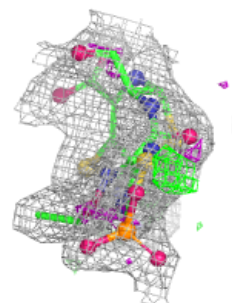
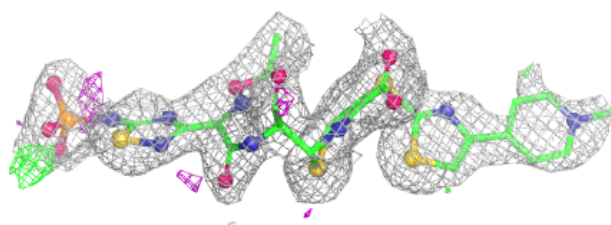
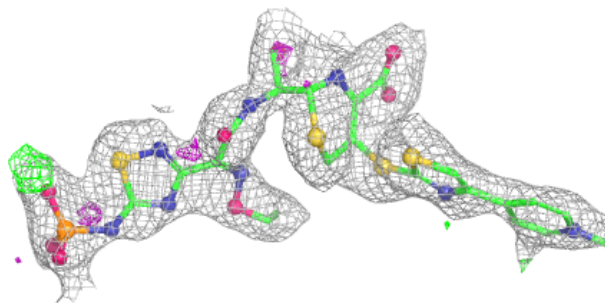
Electron density around CL B 313:

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



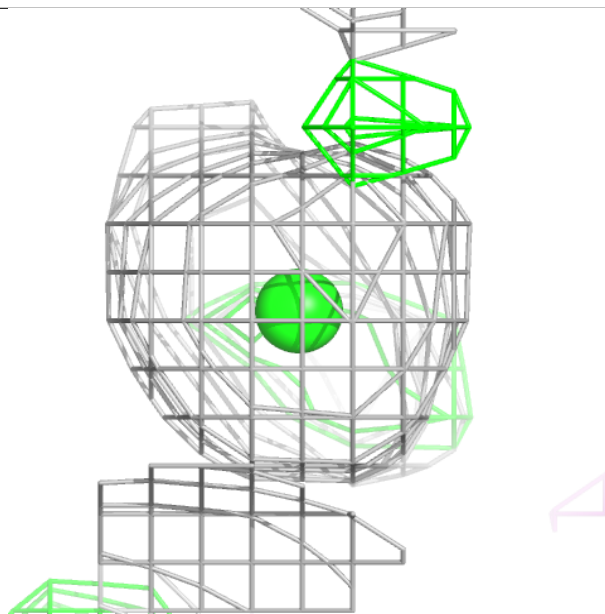
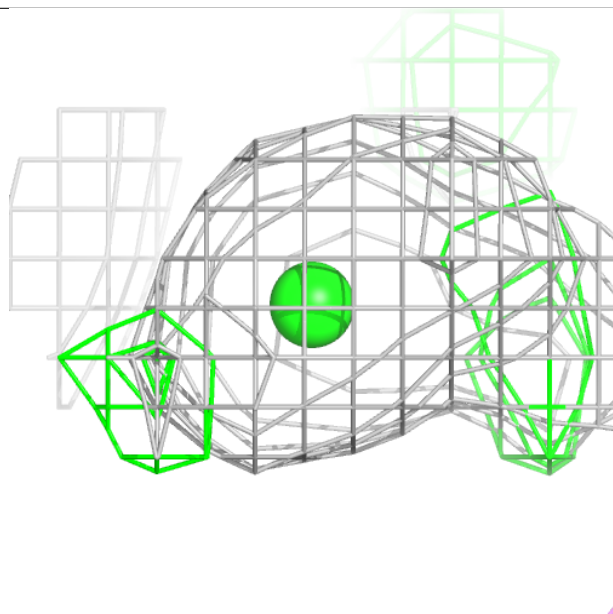
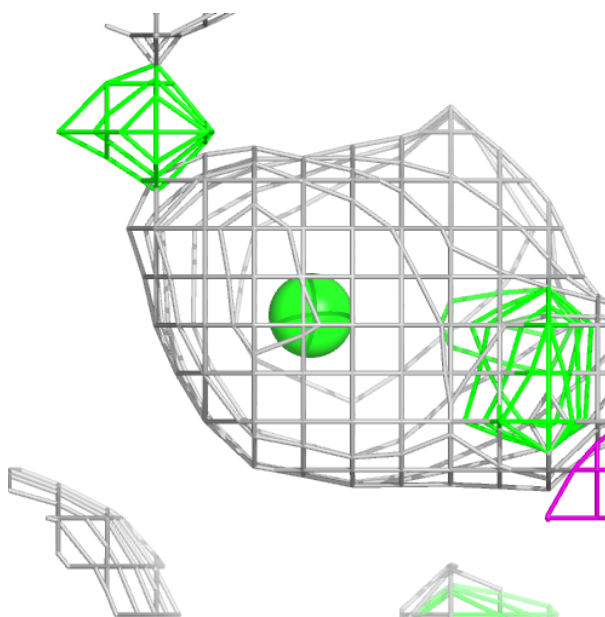
Electron density around A1L2O B 301:

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



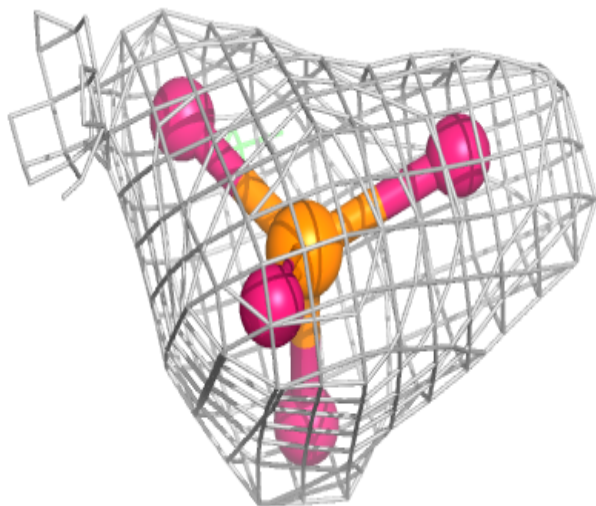
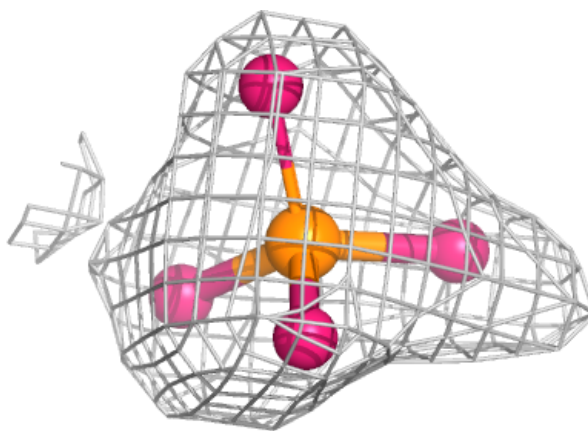
Electron density around CL B 314:

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



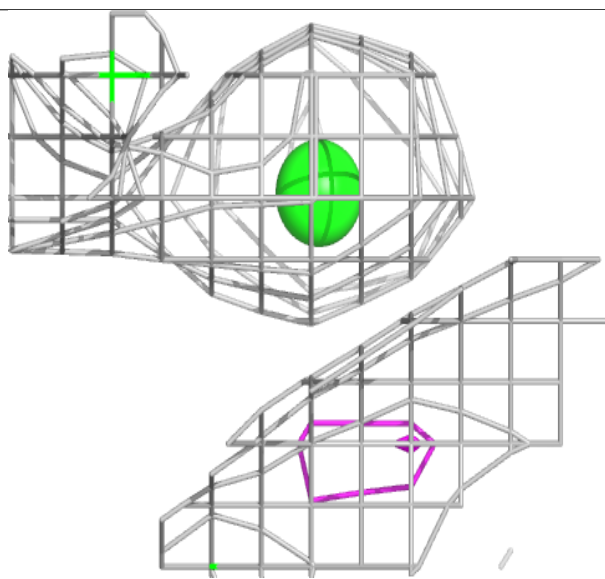
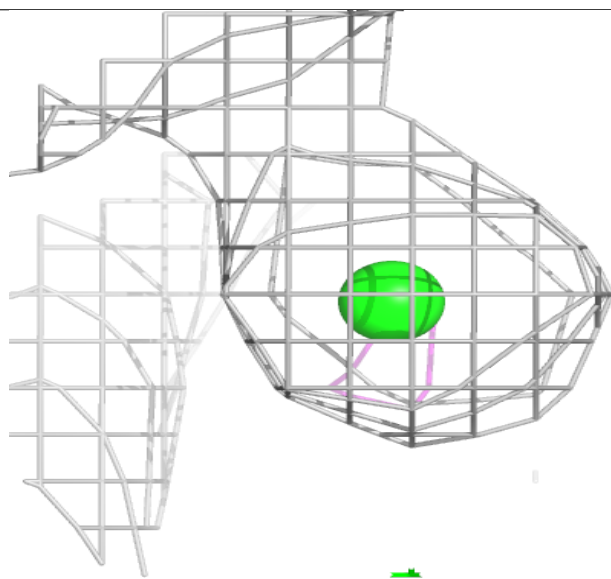
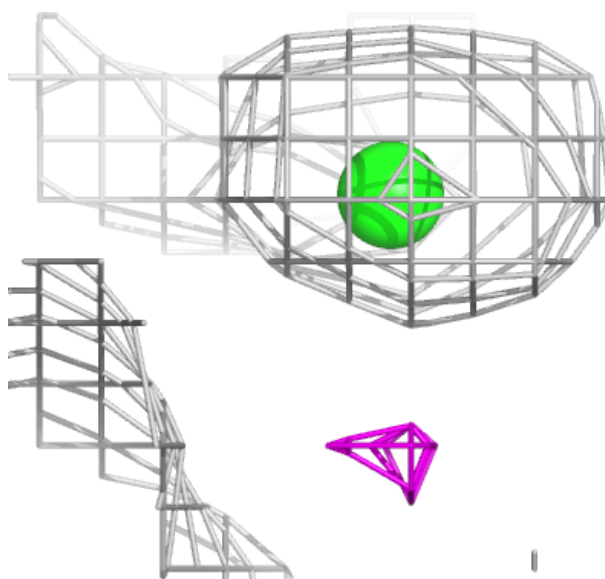
Electron density around PO4 B 319:

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



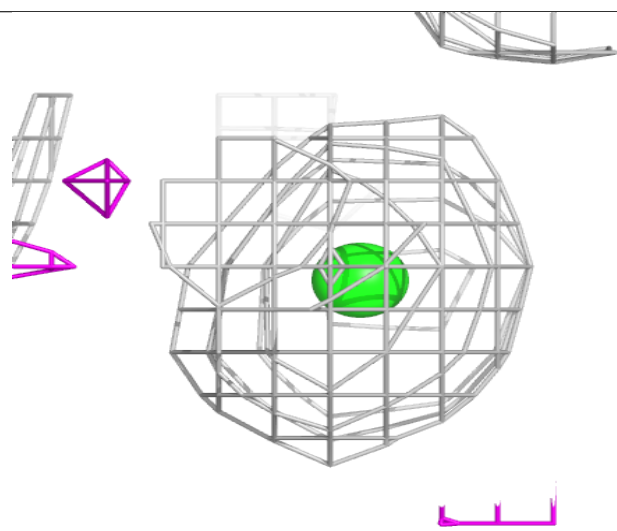
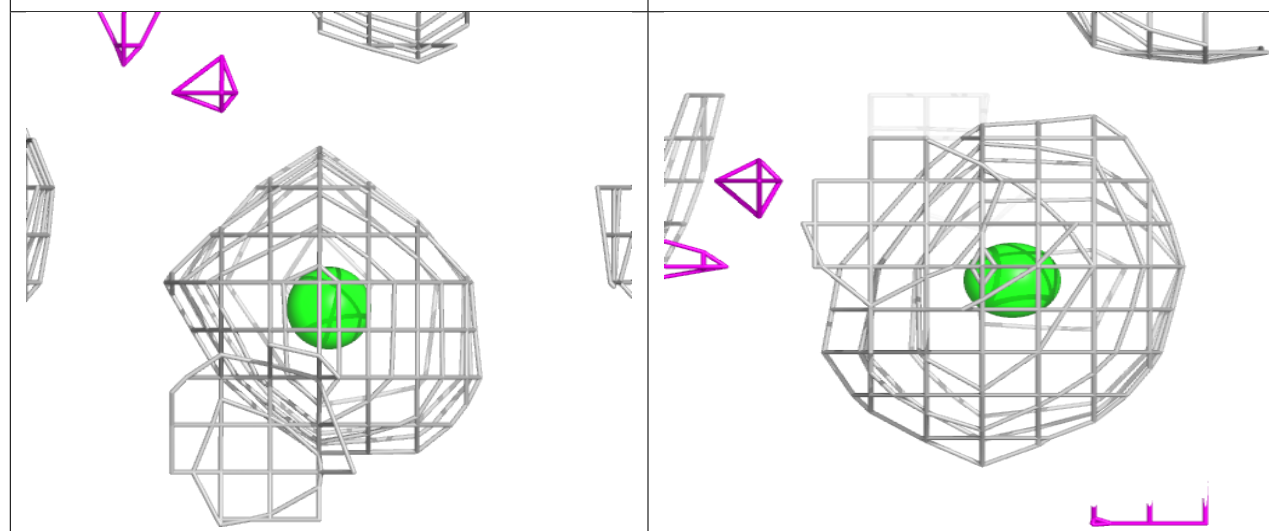
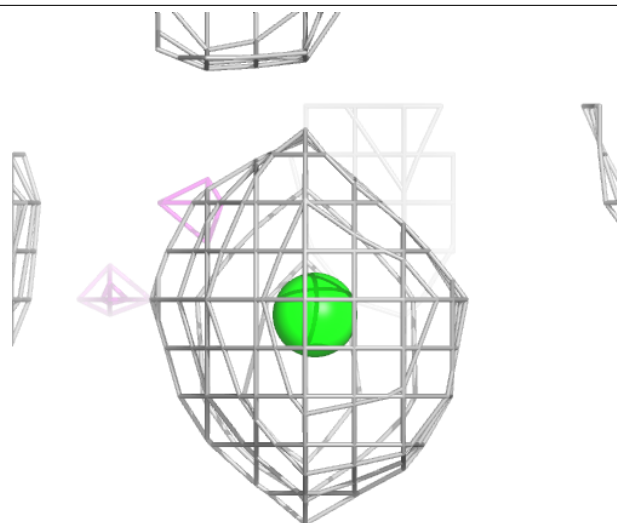
Electron density around CL B 312:

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



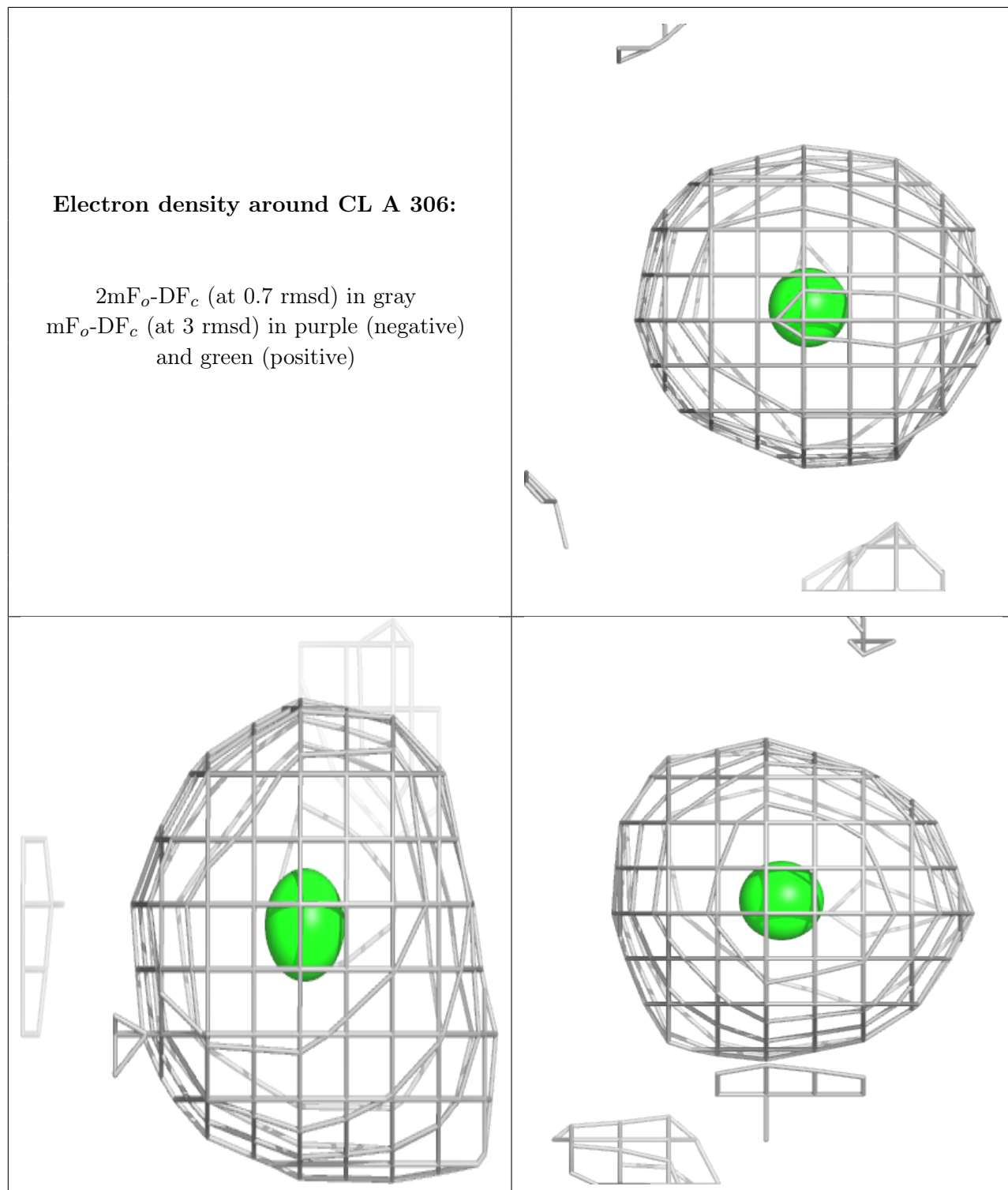
Electron density around CL A 304:

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



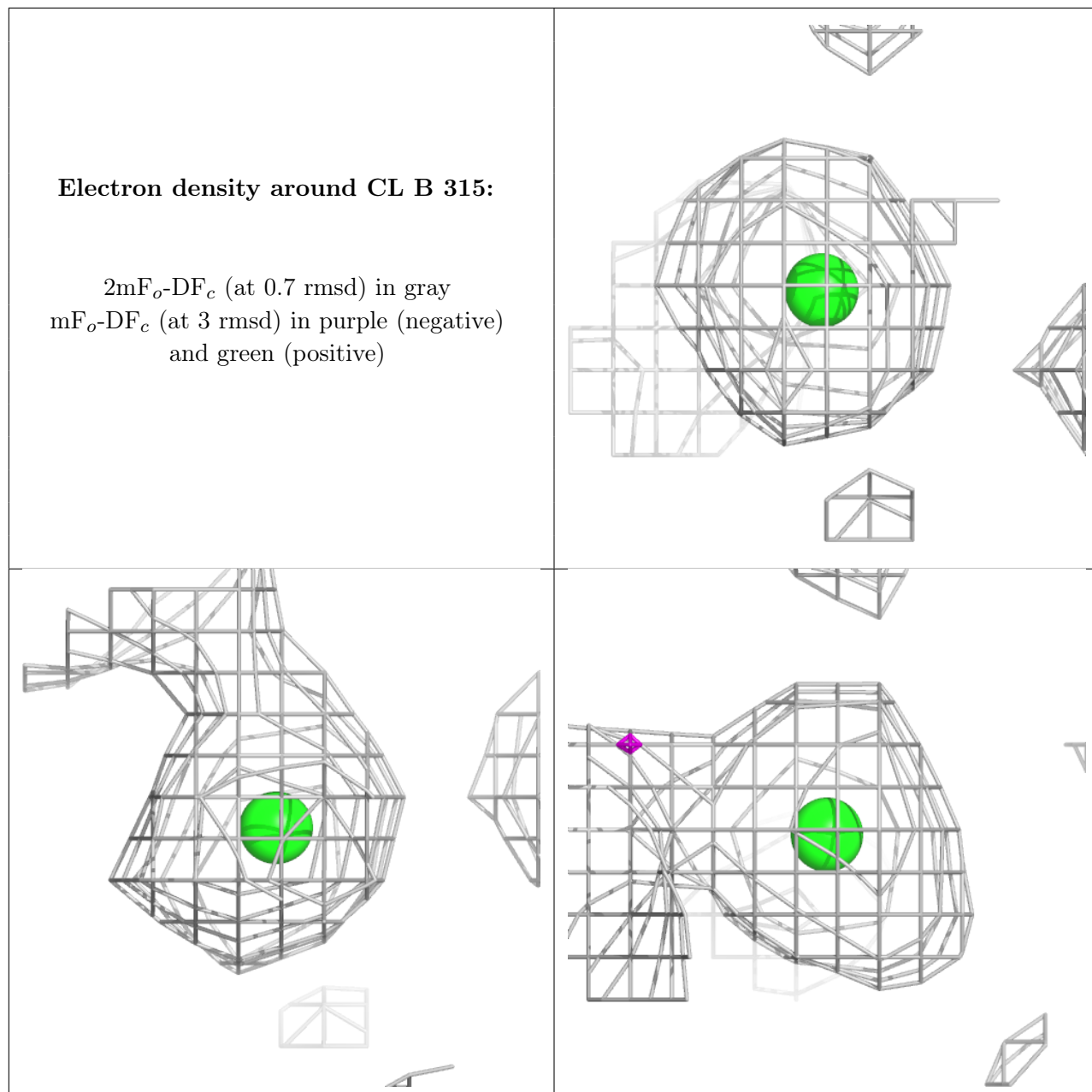
Electron density around CL A 306:

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



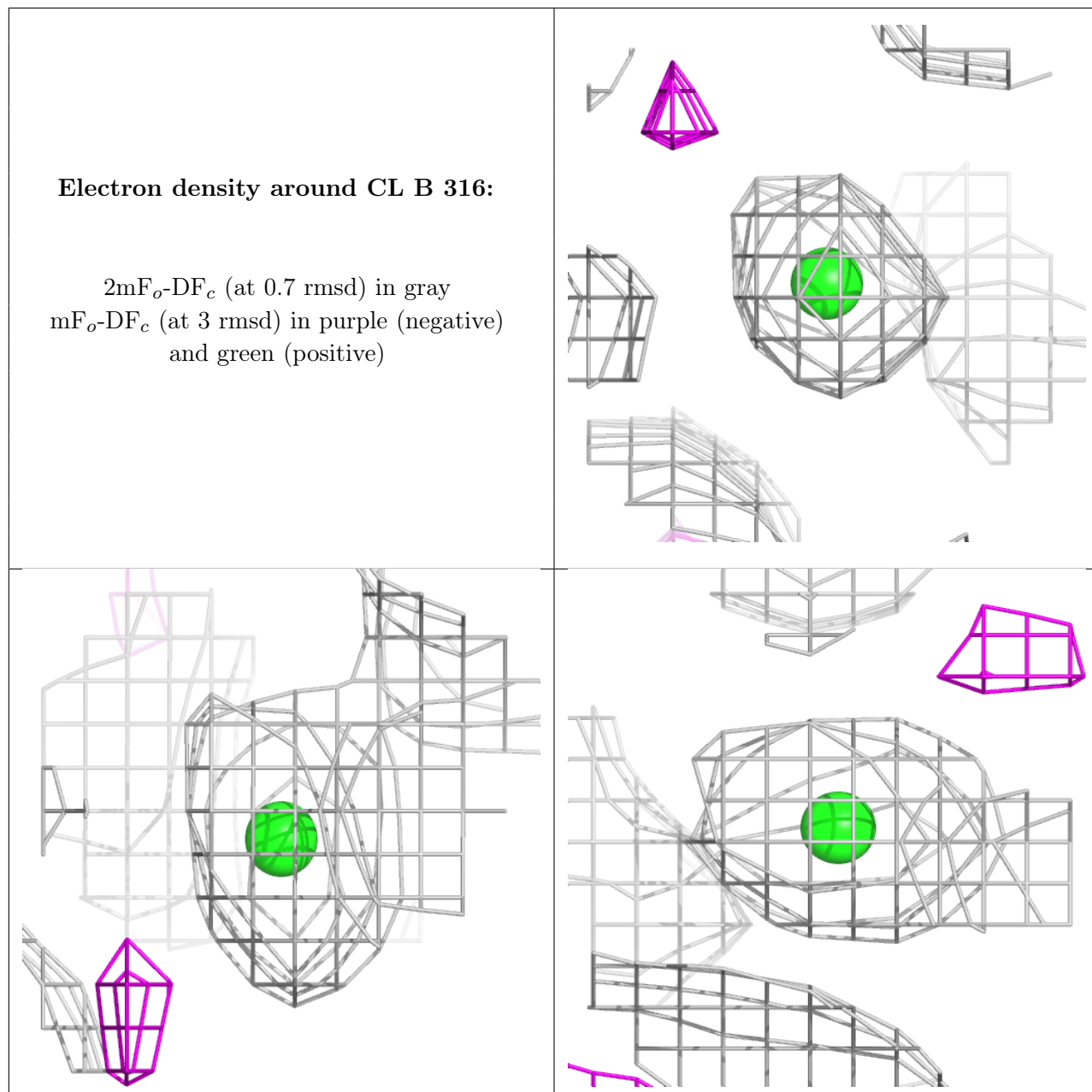
Electron density around CL B 315:

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



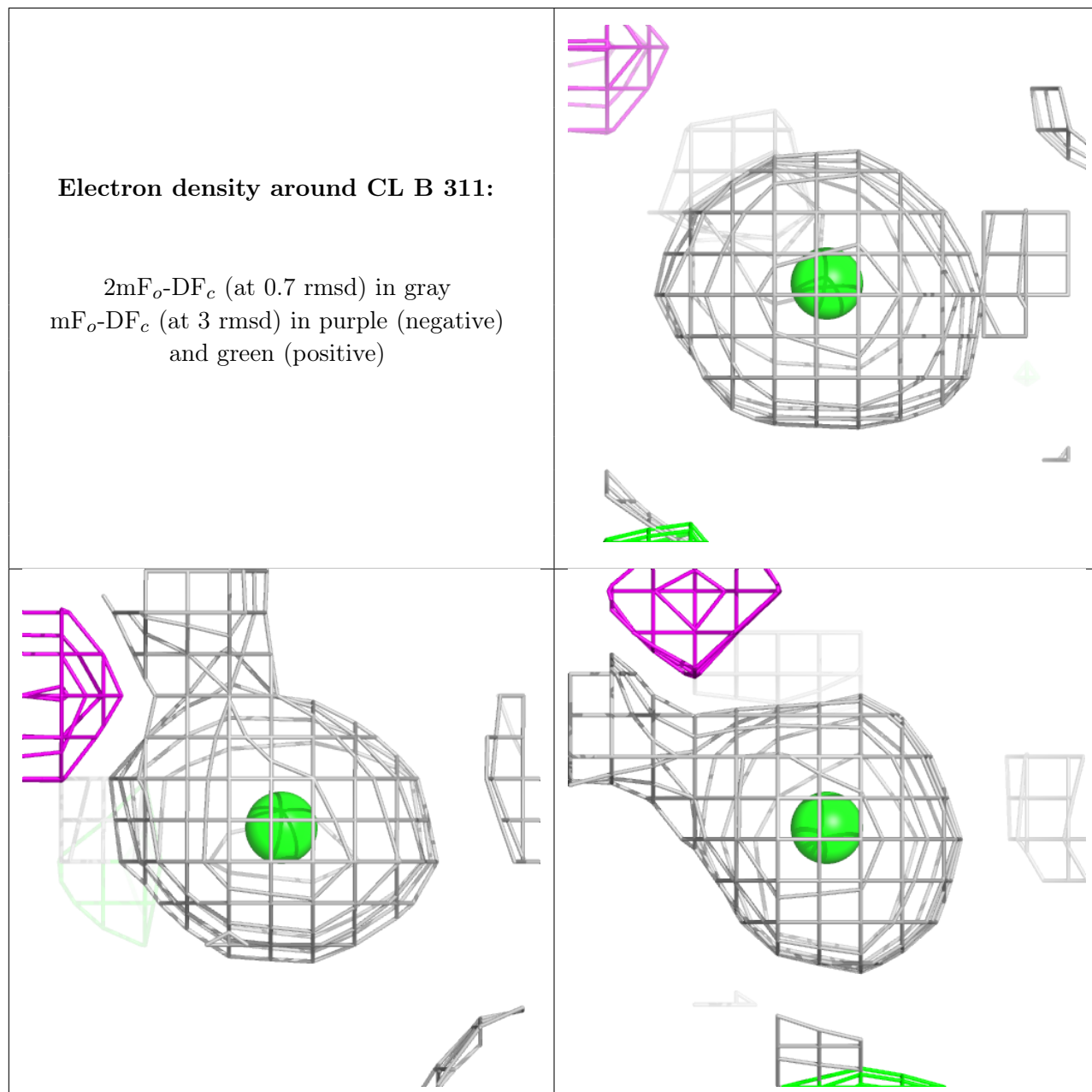
Electron density around CL B 316:

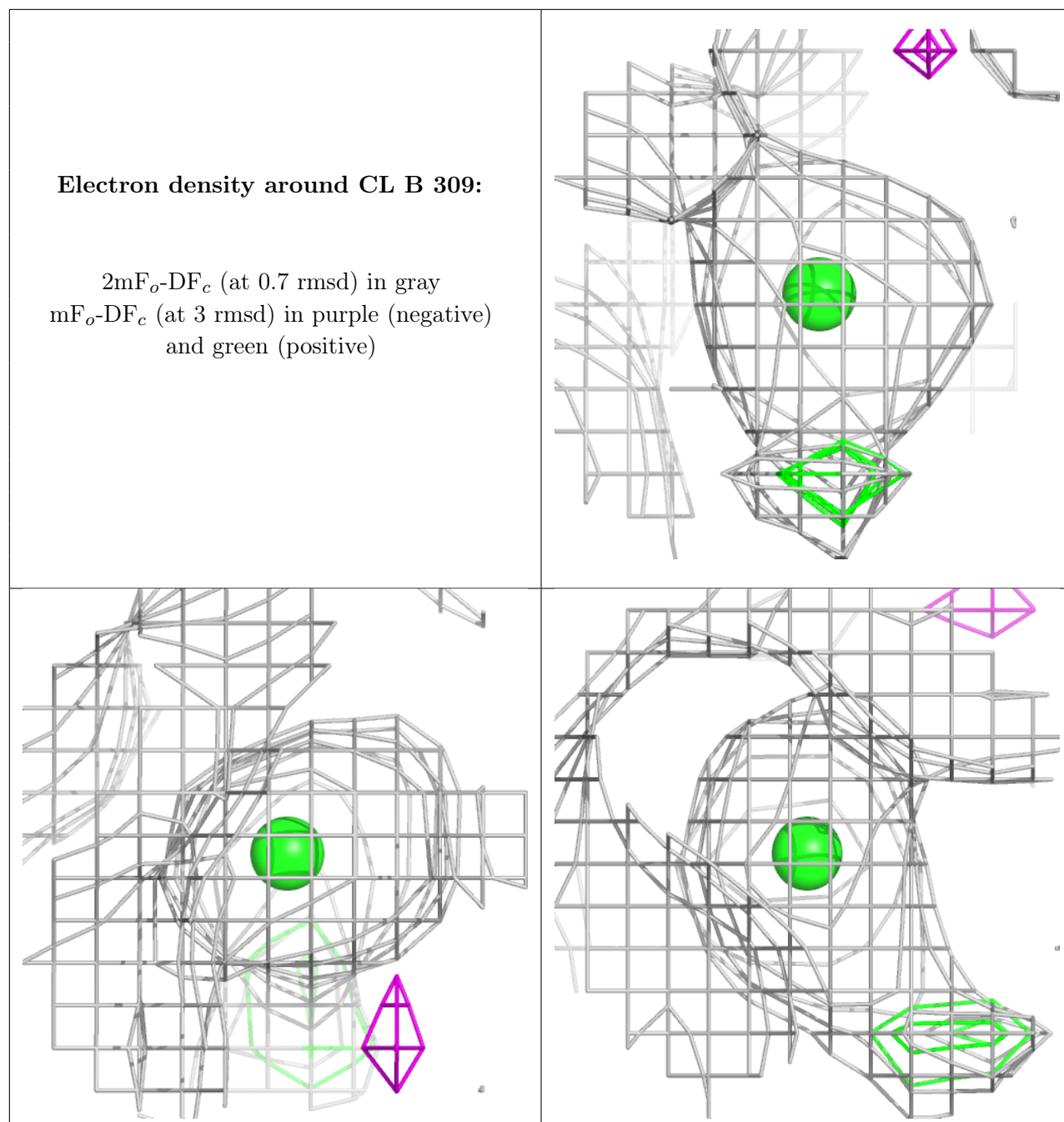
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around CL B 311:

$2mF_o-DF_c$ (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.