



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 03:13 AM UTC

PDB ID : 6KAM / pdb_00006kam
Title : Crystal structure of FKRP in complex with Ba ion, CDP-ribitol, and sugar acceptor
Authors : Kuwabara, N.
Deposited on : 2019-06-23
Resolution : 2.46 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

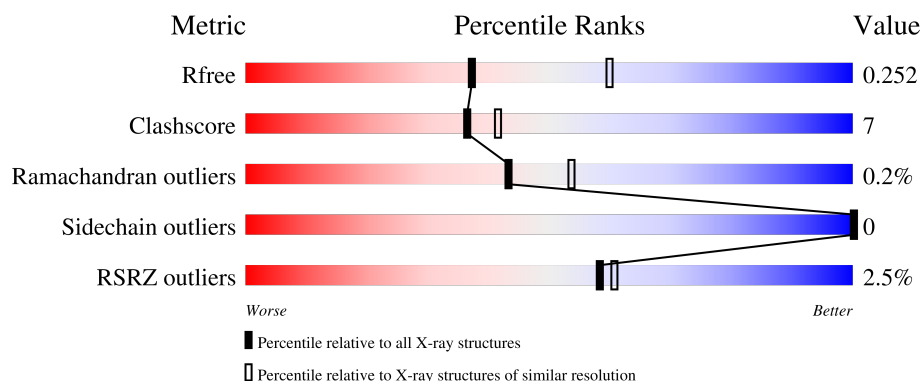
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.46 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	455	3% 84% 13% .
1	B	455	% 83% 15% .
1	C	455	4% 82% 16% .
1	D	455	% 83% 15% .
2	E	3	33% 33% 33%

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
6	RB0	D	505	-	X	-	-
8	PO4	C	507	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 14279 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.

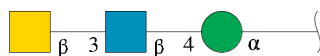
- Molecule 1 is a protein called Fukutin-related protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	446	Total	C	N	O	S	0	2	0
			3444	2211	607	617	9			
1	B	445	Total	C	N	O	S	0	1	0
			3464	2220	616	619	9			
1	C	445	Total	C	N	O	S	0	0	0
			3447	2211	609	618	9			
1	D	443	Total	C	N	O	S	0	0	0
			3441	2206	610	616	9			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	41	GLY	-	expression tag	UNP Q9H9S5
A	42	GLY	-	expression tag	UNP Q9H9S5
A	43	ARG	-	expression tag	UNP Q9H9S5
A	44	PRO	-	expression tag	UNP Q9H9S5
B	41	GLY	-	expression tag	UNP Q9H9S5
B	42	GLY	-	expression tag	UNP Q9H9S5
B	43	ARG	-	expression tag	UNP Q9H9S5
B	44	PRO	-	expression tag	UNP Q9H9S5
C	41	GLY	-	expression tag	UNP Q9H9S5
C	42	GLY	-	expression tag	UNP Q9H9S5
C	43	ARG	-	expression tag	UNP Q9H9S5
C	44	PRO	-	expression tag	UNP Q9H9S5
D	41	GLY	-	expression tag	UNP Q9H9S5
D	42	GLY	-	expression tag	UNP Q9H9S5
D	43	ARG	-	expression tag	UNP Q9H9S5
D	44	PRO	-	expression tag	UNP Q9H9S5

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-alpha-D-mannopyranose.

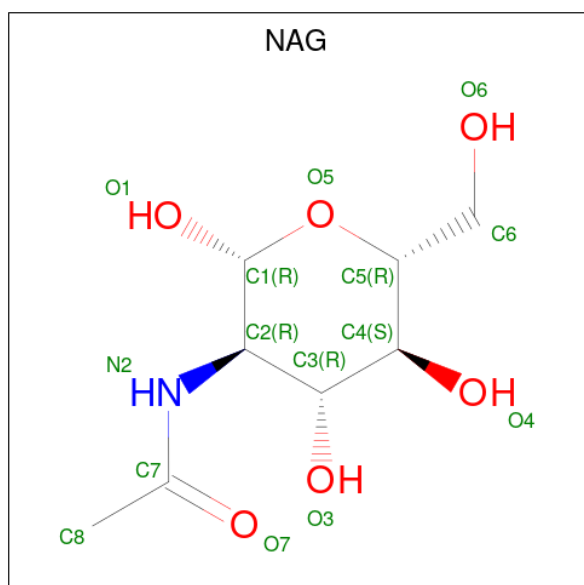


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	3	Total	C	N	O	0	0	0
			40	22	2	16			

- 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	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		
3	D	1	Total	Zn	0	0
			1	1		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆) (labeled as "Ligand of Interest" by depositor).



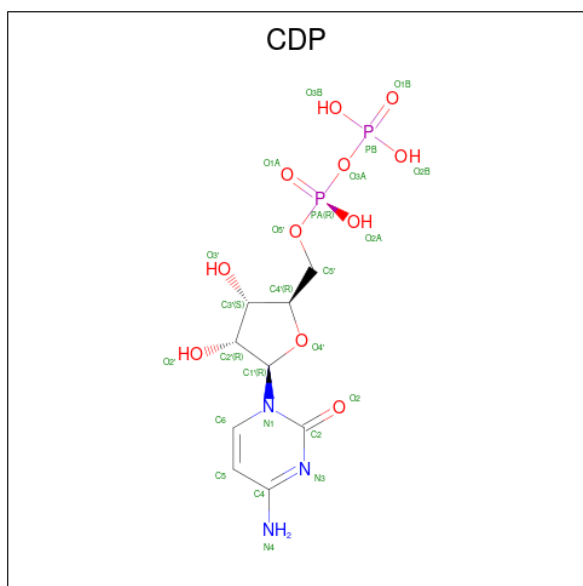
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		

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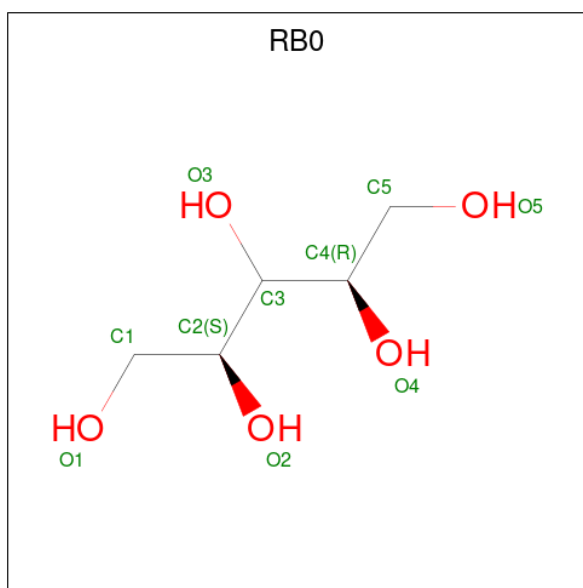
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
	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

- Molecule 5 is CYTIDINE-5'-DIPHOSPHATE (CCD ID: CDP) (formula: $C_9H_{15}N_3O_{11}P_2$) (labeled as "Ligand of Interest" by depositor).



by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			9	5	4		
6	B	1	Total	C	O	0	0
			9	5	4		
6	C	1	Total	C	O	0	0
			9	5	4		
6	D	1	Total	C	O	0	0
			9	5	4		

- Molecule 7 is BARIUM ION (CCD ID: BA) (formula: Ba) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Ba	0	0
			1	1		
7	B	1	Total	Ba	0	0
			1	1		
7	C	1	Total	Ba	0	0
			1	1		
7	D	1	Total	Ba	0	0
			1	1		

- Molecule 8 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P) (labeled as "Ligand of Interest" by depositor).



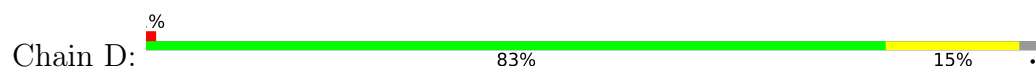
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	C	1	Total	O	P	0	0
			4	3	1		
8	D	1	Total	O	P	0	0
			4	3	1		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	49	Total	O	0	0
			49	49		
9	B	45	Total	O	0	0
			45	45		
9	C	40	Total	O	0	0
			40	40		
9	D	59	Total	O	0	0
			59	59		



- Molecule 1: Fukutin-related protein



- Molecule 2: 2-acetamido-2-deoxy-beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-alpha-D-mannopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.57Å 119.24Å 254.65Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.69 – 2.46 45.69 – 2.46	Depositor EDS
% Data completeness (in resolution range)	99.8 (45.69-2.46) 99.7 (45.69-2.46)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.45Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.208 , 0.249 0.212 , 0.252	Depositor DCC
R_{free} test set	4332 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	55.3	Xtriage
Anisotropy	0.759	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 37.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	14279	wwPDB-VP
Average B, all atoms (Å ²)	75.0	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 23.56 % 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 4.5506e-03. 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, ZN, MAN, NGA, NAG, CDP, BA, RB0

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	0.24	0/3544	0.46	0/4860
1	B	0.27	0/3560	0.52	0/4875
1	C	0.25	0/3543	0.49	0/4854
1	D	0.29	0/3536	0.50	0/4841
All	All	0.26	0/14183	0.49	0/19430

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3444	0	3353	47	0
1	B	3464	0	3400	53	0
1	C	3447	0	3376	62	0
1	D	3441	0	3375	51	0
2	E	40	0	34	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	28	0	26	1	0
4	B	14	0	13	0	0
4	C	28	0	26	0	0
4	D	28	0	26	1	0
5	A	25	0	12	1	0
5	B	25	0	12	0	0
5	C	25	0	12	1	0
5	D	25	0	12	0	0
6	A	9	0	9	0	0
6	B	9	0	9	0	0
6	C	9	0	9	1	0
6	D	9	0	9	1	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
8	C	4	0	0	2	0
8	D	4	0	0	0	0
9	A	49	0	0	0	0
9	B	45	0	0	0	0
9	C	40	0	0	3	0
9	D	59	0	0	4	0
All	All	14279	0	13713	200	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 7.

All (200) 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:D:140:LEU:HD21	1:D:238:LEU:HD13	1.45	0.96
1:A:229:ARG:HG2	1:C:489:LEU:HD22	1.50	0.92
1:D:134:ARG:HH12	1:D:244:ARG:HG3	1.43	0.84
1:B:47:PRO:HA	1:B:48:ARG:HB2	1.65	0.79
1:B:258:GLU:HG2	1:B:259:ARG:N	1.96	0.78
1:B:105:GLN:HE21	1:B:110:ARG:HH12	1.33	0.76
1:A:444:HIS:HA	1:A:447:GLN:HG2	1.68	0.74
1:C:452:LEU:HD23	1:C:453:PRO:HD2	1.68	0.74
1:C:367:ILE:C	1:C:367:ILE:HD12	2.13	0.74
1:D:373:GLY:O	1:D:379:ARG:NH1	2.21	0.73
1:B:255:TRP:HA	1:B:258:GLU:OE2	1.90	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:401:ASP:H	1:B:431:THR:HG21	1.57	0.69
1:B:134:ARG:HG2	1:B:136:GLU:OE2	1.92	0.68
1:D:205:ARG:HA	4:D:503:NAG:H83	1.74	0.68
1:A:327:ARG:HH21	1:A:457:PHE:HB3	1.59	0.67
1:A:303:ASP:OD2	1:A:435:HIS:NE2	2.26	0.67
1:B:47:PRO:HG2	1:B:145:VAL:HG21	1.78	0.66
1:A:488:LEU:O	1:A:489:LEU:HD22	1.96	0.66
1:C:148:LEU:HD23	1:C:156:VAL:HG13	1.76	0.66
1:A:167:ARG:NH1	1:A:195:ASP:OD2	2.28	0.66
1:D:169:LEU:HD11	1:D:195:ASP:HB2	1.76	0.66
1:B:65:GLU:OE1	1:B:134:ARG:NH2	2.29	0.65
1:A:142:GLU:HA	1:A:145:VAL:HG12	1.79	0.65
1:C:439:VAL:HG11	1:C:472:LYS:HB3	1.78	0.64
1:A:447:GLN:HB2	1:A:448:PRO:HD3	1.80	0.63
1:C:295:ARG:NH1	8:C:507:PO4:O4	2.28	0.63
1:B:48:ARG:NH2	1:D:124:GLU:OE1	2.32	0.63
1:A:67:VAL:HG22	1:A:79:VAL:HG11	1.80	0.63
1:D:79:VAL:HG22	1:D:99:VAL:HG22	1.81	0.63
1:C:291:LYS:NZ	9:C:602:HOH:O	2.30	0.63
1:D:148:LEU:HD12	1:D:201:LEU:HD23	1.81	0.63
1:C:452:LEU:HD23	1:C:453:PRO:CD	2.29	0.61
1:D:142:GLU:O	1:D:146:GLU:HG3	2.01	0.60
1:C:376:GLU:OE1	1:C:379:ARG:NH2	2.35	0.60
1:B:142:GLU:O	1:B:145:VAL:HG22	2.01	0.59
1:B:423:ARG:HA	1:B:423:ARG:HE	1.66	0.59
1:B:97:PRO:HB2	1:D:119:THR:HG22	1.85	0.59
1:C:428:THR:HG22	1:C:440:GLU:HB3	1.84	0.59
1:A:303:ASP:OD1	1:A:361:TYR:OH	2.20	0.59
1:C:422:PRO:HB3	1:C:443:GLU:OE2	2.04	0.58
1:D:420:PHE:HB3	1:D:427:MET:HE3	1.85	0.58
1:B:255:TRP:CG	1:B:255:TRP:O	2.56	0.58
1:C:49:VAL:HG22	1:C:125:PHE:HB2	1.85	0.58
1:C:67:VAL:HG23	1:C:79:VAL:HG11	1.84	0.58
1:B:423:ARG:HA	1:B:423:ARG:NE	2.19	0.58
1:C:439:VAL:HG22	1:C:440:GLU:O	2.04	0.57
1:B:97:PRO:HD2	1:D:119:THR:HG21	1.86	0.57
1:B:160:VAL:HG12	1:B:162:THR:HG23	1.85	0.57
1:D:254:ARG:O	1:D:258:GLU:HG3	2.05	0.57
1:C:301:VAL:HG21	1:D:89:PRO:HB3	1.87	0.57
1:A:176:ARG:HG2	1:A:313:TRP:CD2	2.40	0.56
1:C:340:TYR:CZ	1:C:460:GLN:HG3	2.40	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:403:PHE:HB2	1:B:417:LEU:HB2	1.86	0.56
1:D:57:GLU:OE2	1:D:264:ARG:NH2	2.38	0.56
1:C:437:GLN:OE1	1:C:437:GLN:N	2.30	0.56
1:A:400:GLY:C	1:A:431:THR:HG21	2.31	0.56
1:A:166:ALA:HB2	1:A:198:ALA:HB2	1.87	0.55
1:A:148:LEU:HD23	1:A:156:VAL:HG13	1.88	0.55
1:D:134:ARG:NH1	1:D:244:ARG:HG3	2.17	0.55
1:D:140:LEU:HD21	1:D:238:LEU:CD1	2.28	0.55
1:A:437:GLN:NE2	5:A:504:CDP:O1A	2.35	0.55
1:C:141:LEU:O	1:C:145:VAL:HG23	2.06	0.55
1:C:404:ARG:NH2	9:C:605:HOH:O	2.40	0.55
1:C:406:GLN:HG2	1:C:412:HIS:HA	1.88	0.54
1:A:433:LEU:HD13	1:A:434:ASP:N	2.23	0.54
1:B:279:TRP:CD1	1:B:283:ARG:HG2	2.42	0.54
1:A:448:PRO:HD2	1:A:463:ASN:ND2	2.22	0.54
1:C:348:LEU:HD23	1:C:357:ILE:HD11	1.88	0.54
1:D:406:GLN:OE1	1:D:414:HIS:HB3	2.06	0.54
1:A:482:GLN:HA	1:A:491:LEU:HD11	1.90	0.54
1:B:250:THR:HG23	1:B:253:ALA:H	1.71	0.53
1:B:406:GLN:HG2	1:B:412:HIS:HA	1.91	0.53
1:B:62:ALA:HB3	1:B:131:ASP:OD1	2.09	0.53
1:B:68:ASP:O	1:B:72:GLN:HG2	2.09	0.53
1:A:82:ALA:HB1	1:A:104:LEU:HD11	1.91	0.53
1:A:176:ARG:HD3	1:A:313:TRP:CZ2	2.44	0.53
1:D:159:PRO:HD2	1:D:238:LEU:O	2.09	0.53
1:D:305:PRO:HD2	1:D:308:LEU:HD12	1.90	0.53
1:A:140:LEU:HA	1:A:143:ARG:HG3	1.91	0.52
1:A:387:VAL:HG22	1:A:393:VAL:HG22	1.90	0.52
1:A:437:GLN:OE1	1:A:437:GLN:N	2.40	0.52
1:B:47:PRO:N	1:B:48:ARG:HH11	2.08	0.52
1:D:252:HIS:O	1:D:256:LYS:HG3	2.10	0.52
1:A:491:LEU:HD12	1:A:491:LEU:H	1.75	0.52
1:B:258:GLU:N	1:B:258:GLU:OE1	2.40	0.52
1:C:412:HIS:ND1	2:E:3:NGA:O7	2.25	0.51
1:D:234:GLN:NE2	9:D:604:HOH:O	2.42	0.51
1:C:154:ARG:NH2	9:C:607:HOH:O	2.44	0.51
1:C:427:MET:HG2	1:C:443:GLU:HG3	1.93	0.51
1:A:482:GLN:HG3	1:C:176:ARG:HH22	1.76	0.51
1:A:433:LEU:HD12	1:A:435:HIS:CD2	2.46	0.51
1:B:49:VAL:HG12	1:B:125:PHE:HB2	1.94	0.50
1:C:177:GLU:O	1:C:316:PRO:HG3	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:190:ARG:NH2	9:D:604:HOH:O	2.40	0.50
1:C:369:LEU:HD23	1:C:421:TYR:HB3	1.94	0.50
1:D:291:LYS:NZ	9:D:605:HOH:O	2.41	0.50
1:B:140:LEU:HD22	1:B:238:LEU:HD23	1.93	0.50
1:C:367:ILE:C	1:C:367:ILE:CD1	2.85	0.49
1:D:401:ASP:HB2	1:D:431:THR:OG1	2.11	0.49
1:A:137:ALA:HB3	1:A:140:LEU:HD11	1.93	0.49
1:C:406:GLN:HG3	1:C:414:HIS:HB3	1.93	0.49
1:C:452:LEU:HD22	1:C:453:PRO:O	2.12	0.49
1:A:377:GLN:HB3	1:A:394:TRP:CD2	2.46	0.49
1:D:145:VAL:O	1:D:149:ARG:HG3	2.12	0.49
1:A:169:LEU:HD11	1:A:195:ASP:HB2	1.95	0.49
1:D:140:LEU:HA	1:D:140:LEU:HD23	1.37	0.48
1:D:88:TYR:HA	1:D:89:PRO:C	2.38	0.48
1:A:491:LEU:HD13	1:C:176:ARG:HH12	1.76	0.48
1:A:272:LEU:HD13	1:B:108:LEU:HD21	1.96	0.48
1:B:404:ARG:NH2	1:B:416:ASP:OD2	2.45	0.48
1:C:291:LYS:HG3	1:C:408:SER:HB3	1.96	0.48
1:D:375:CYS:HB3	1:D:378:LEU:HD12	1.94	0.48
1:A:403:PHE:HB2	1:A:417:LEU:HB2	1.94	0.48
1:D:107:ALA:HB3	1:D:110:ARG:HG3	1.95	0.48
1:D:159:PRO:HD3	1:D:235:LEU:HD11	1.96	0.48
1:B:134:ARG:NH1	1:B:244:ARG:HG3	2.29	0.48
1:D:441:PHE:HE2	1:D:446:LEU:HD11	1.79	0.48
1:C:67:VAL:HG11	1:C:93:LEU:HD22	1.95	0.48
1:C:340:TYR:CE1	1:C:460:GLN:HG3	2.49	0.48
1:B:170:ALA:HB2	1:B:186:PRO:HD3	1.95	0.47
1:C:387:VAL:HG22	1:C:393:VAL:HG22	1.96	0.47
1:A:199:VAL:HG21	1:A:240:PHE:CE1	2.50	0.47
1:B:369:LEU:O	1:B:372:VAL:HG13	2.15	0.47
1:C:299:THR:HG23	2:E:3:NGA:H3	1.96	0.47
1:B:145:VAL:HG12	1:B:201:LEU:HD11	1.96	0.47
1:C:419:PRO:O	1:C:429:LYS:HB3	2.15	0.47
1:B:98:ASN:OD1	1:D:119:THR:HG23	2.16	0.46
1:B:418:TRP:CZ3	1:B:433:LEU:HD11	2.50	0.46
1:C:363:VAL:HB	1:C:415:VAL:HG22	1.96	0.46
1:D:238:LEU:HD23	1:D:238:LEU:HA	1.65	0.46
1:C:452:LEU:HD23	1:C:453:PRO:N	2.31	0.46
1:D:175:LEU:HD11	1:D:228:LEU:HD11	1.96	0.46
1:A:159:PRO:HD2	1:A:238:LEU:O	2.16	0.46
1:B:141:LEU:O	1:B:145:VAL:HG13	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:171:LEU:HD21	1:B:224:LEU:HD23	1.98	0.46
1:C:88:TYR:HA	1:C:89:PRO:C	2.41	0.46
4:A:502:NAG:H61	1:C:436:ARG:NH1	2.31	0.46
1:A:160:VAL:HG12	1:A:162:THR:HG23	1.98	0.45
1:A:448:PRO:HD2	1:A:463:ASN:HD22	1.79	0.45
1:B:145:VAL:O	1:B:149:ARG:HG3	2.17	0.45
1:C:350:ALA:HA	1:C:354:GLY:HA2	1.98	0.45
1:C:472:LYS:H	1:C:472:LYS:HG2	1.62	0.45
1:D:409:GLU:HG3	9:D:625:HOH:O	2.16	0.45
1:A:108:LEU:HD21	1:B:272:LEU:HD13	1.99	0.44
1:C:110:ARG:O	1:D:211:SER:HB2	2.17	0.44
1:B:429:LYS:NZ	1:B:431:THR:O	2.45	0.44
1:B:291:LYS:HG2	1:B:408:SER:HB3	1.99	0.44
1:C:211:SER:HB2	1:D:110:ARG:O	2.17	0.44
1:B:84:ASP:O	1:B:105:GLN:HB3	2.17	0.44
1:C:134:ARG:HH21	1:C:136:GLU:CD	2.26	0.44
1:D:418:TRP:CD2	6:D:505:RB0:H1A	2.53	0.44
1:D:375:CYS:HB3	1:D:378:LEU:CD1	2.48	0.43
1:D:166:ALA:HB2	1:D:198:ALA:HB2	2.00	0.43
1:D:363:VAL:HB	1:D:415:VAL:HG22	1.99	0.43
1:A:176:ARG:HD3	1:A:313:TRP:CE2	2.54	0.43
1:B:91:LEU:HD23	1:B:91:LEU:HA	1.87	0.43
1:C:377:GLN:HB3	1:C:394:TRP:CD2	2.53	0.43
1:A:142:GLU:CA	1:A:145:VAL:HG12	2.47	0.43
1:B:305:PRO:HD2	1:B:308:LEU:HD12	2.00	0.43
1:C:437:GLN:NE2	5:C:504:CDP:O1A	2.51	0.43
1:D:323:ARG:HG2	1:D:455:ALA:HB1	1.99	0.43
1:A:427:MET:SD	1:A:443:GLU:HG3	2.59	0.43
1:D:238:LEU:CD2	1:D:239:THR:H	2.32	0.43
1:A:367:ILE:O	1:A:419:PRO:HA	2.18	0.43
1:B:265:ARG:HH12	1:B:285:GLU:CD	2.27	0.43
1:C:367:ILE:HD12	1:C:367:ILE:O	2.18	0.42
1:B:134:ARG:NE	1:B:243:ALA:HA	2.34	0.42
1:C:295:ARG:HH22	8:C:507:PO4:P	2.42	0.42
1:A:191:CYS:HB2	1:A:233:VAL:HG13	2.01	0.42
1:B:289:CYS:SG	1:B:295:ARG:HA	2.59	0.42
1:C:56:PHE:HB3	1:C:131:ASP:HB2	2.01	0.42
1:C:160:VAL:HG12	1:C:162:THR:HG23	2.01	0.42
1:B:290:ASN:OD1	1:B:293:THR:HG23	2.20	0.42
1:A:444:HIS:HA	1:A:447:GLN:CG	2.42	0.42
1:C:176:ARG:HD2	1:C:313:TRP:CD2	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:88:TYR:HA	1:B:89:PRO:C	2.45	0.42
1:C:404:ARG:HE	1:C:414:HIS:CD2	2.38	0.42
1:C:438:ASP:OD1	6:C:505:RB0:O1	2.37	0.42
1:A:160:VAL:HG22	1:A:240:PHE:HB2	2.03	0.41
1:C:468:PHE:CE1	1:C:472:LYS:HD3	2.55	0.41
1:B:308:LEU:HD22	1:B:483:TYR:CZ	2.55	0.41
1:C:237:ASP:O	1:C:238:LEU:HD23	2.20	0.41
1:B:255:TRP:HA	1:B:258:GLU:CD	2.46	0.41
1:C:488:LEU:H	1:C:488:LEU:HD22	1.86	0.41
1:D:143:ARG:CZ	1:D:238:LEU:HD11	2.50	0.41
1:D:293:THR:OG1	1:D:294:THR:O	2.29	0.41
1:C:427:MET:CG	1:C:443:GLU:HG3	2.51	0.41
1:D:70:PHE:CE1	1:D:135:ALA:HB2	2.55	0.41
1:C:489:LEU:O	1:C:489:LEU:HG	2.20	0.41
1:D:143:ARG:NH2	1:D:238:LEU:HD11	2.35	0.41
1:A:88:TYR:HA	1:A:89:PRO:C	2.46	0.41
1:A:375:CYS:HB3	1:A:378:LEU:HB2	2.03	0.41
1:B:134:ARG:CZ	1:B:244:ARG:H	2.34	0.41
1:B:48:ARG:CZ	1:D:124:GLU:OE1	2.69	0.40
1:A:166:ALA:CB	1:A:198:ALA:HB2	2.50	0.40
1:B:404:ARG:NE	1:B:416:ASP:OD1	2.53	0.40
1:D:70:PHE:CE1	1:D:141:LEU:HD11	2.57	0.40
1:C:341:TRP:CE2	1:C:366:GLY:HA3	2.56	0.40
1:D:101:LEU:HD21	1:D:103:LEU:HD21	2.04	0.40
1:C:457:PHE:CE2	1:C:459:ALA:HB2	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	446/455 (98%)	432 (97%)	12 (3%)	2 (0%)	30 37

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	444/455 (98%)	426 (96%)	17 (4%)	1 (0%)	43	54
1	C	443/455 (97%)	430 (97%)	13 (3%)	0	100	100
1	D	439/455 (96%)	427 (97%)	12 (3%)	0	100	100
All	All	1772/1820 (97%)	1715 (97%)	54 (3%)	3 (0%)	43	54

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	399	GLU
1	B	399	GLU
1	A	434	ASP

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	344/357 (96%)	344 (100%)	0	100	100
1	B	349/357 (98%)	349 (100%)	0	100	100
1	C	347/357 (97%)	347 (100%)	0	100	100
1	D	347/357 (97%)	347 (100%)	0	100	100
All	All	1387/1428 (97%)	1387 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	374	ASN
1	B	105	GLN
1	B	460	GLN
1	B	464	ASN
1	C	245	GLN
1	C	353	HIS

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Mol	Chain	Res	Type
1	D	61	ASN
1	D	72	GLN
1	D	234	GLN
1	D	460	GLN
1	D	480	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

3 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	MAN	E	1	8,2	12,12,12	0.95	0	17,17,17	1.15	2 (11%)
2	NAG	E	2	2	14,14,15	0.45	0	17,19,21	0.41	0
2	NGA	E	3	8,2	14,14,15	0.64	0	17,19,21	1.22	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MAN	E	1	8,2	-	0/2/22/22	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NGA	E	3	8,2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	3	NGA	O3-C3-C4	-3.19	102.86	110.38
2	E	1	MAN	C3-C4-C5	2.48	114.73	110.23
2	E	1	MAN	O2-C2-C3	-2.17	105.27	110.38

There are no chirality outliers.

All (2) torsion outliers are listed below:

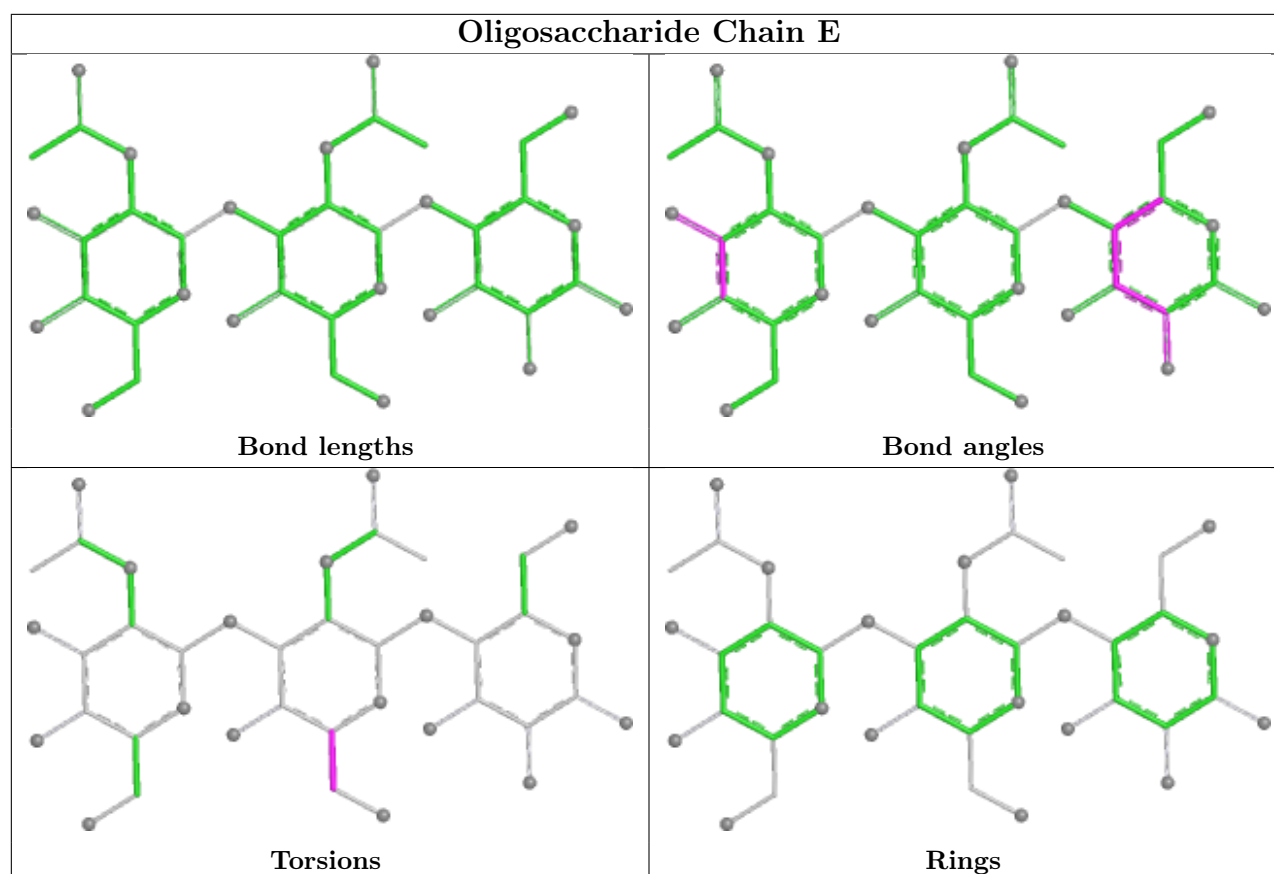
Mol	Chain	Res	Type	Atoms
2	E	2	NAG	O5-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	3	NGA	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry [i](#)

Of 25 ligands modelled in this entry, 8 are monoatomic - leaving 17 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	$\# Z > 2$
8	PO4	C	507	2	0,3,4	-	-	0,3,6	-	-
6	RB0	A	505	5	8,8,9	0.30	0	10,10,11	0.45	0
6	RB0	C	505	5	8,8,9	0.45	0	10,10,11	1.13	0
4	NAG	C	503	1	14,14,15	0.28	0	17,19,21	0.40	0
4	NAG	A	502	1	14,14,15	0.60	1 (7%)	17,19,21	0.75	0
6	RB0	B	504	5	8,8,9	0.65	0	10,10,11	1.16	0
5	CDP	D	504	7,6	25,26,26	0.96	1 (4%)	38,40,40	0.92	1 (2%)
8	PO4	D	507	2	0,3,4	-	-	0,3,6	-	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	CDP	B	503	7,6	25,26,26	0.99	1 (4%)	38,40,40	0.94	2 (5%)
4	NAG	B	502	1	14,14,15	0.40	0	17,19,21	0.43	0
4	NAG	A	503	1	14,14,15	0.36	0	17,19,21	0.57	0
4	NAG	D	502	1	14,14,15	1.11	1 (7%)	17,19,21	1.75	2 (11%)
5	CDP	A	504	7,6	25,26,26	0.99	1 (4%)	38,40,40	0.88	1 (2%)
6	RB0	D	505	5	8,8,9	0.53	0	10,10,11	1.90	2 (20%)
4	NAG	C	502	1	14,14,15	0.28	0	17,19,21	0.54	0
5	CDP	C	504	7,6	25,26,26	1.00	1 (4%)	38,40,40	0.94	1 (2%)
4	NAG	D	503	1	14,14,15	0.43	0	17,19,21	0.76	1 (5%)

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
6	RB0	A	505	5	-	4/10/10/12	-
6	RB0	C	505	5	-	4/10/10/12	-
4	NAG	C	503	1	-	1/6/23/26	0/1/1/1
4	NAG	A	502	1	-	1/6/23/26	0/1/1/1
6	RB0	B	504	5	-	4/10/10/12	-
5	CDP	D	504	7,6	-	3/16/32/32	0/2/2/2
5	CDP	B	503	7,6	-	2/16/32/32	0/2/2/2
4	NAG	B	502	1	-	2/6/23/26	0/1/1/1
4	NAG	A	503	1	-	2/6/23/26	0/1/1/1
4	NAG	D	502	1	-	2/6/23/26	0/1/1/1
5	CDP	A	504	7,6	-	1/16/32/32	0/2/2/2
6	RB0	D	505	5	-	10/10/10/12	-
4	NAG	C	502	1	-	1/6/23/26	0/1/1/1
5	CDP	C	504	7,6	-	5/16/32/32	0/2/2/2
4	NAG	D	503	1	-	2/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	504	CDP	PA-O3A	3.20	1.62	1.59
5	A	504	CDP	PA-O3A	2.98	1.62	1.59
5	B	503	CDP	PA-O3A	2.97	1.62	1.59

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	502	NAG	C1-C2	-2.88	1.48	1.52
5	D	504	CDP	PA-O3A	2.59	1.62	1.59
4	A	502	NAG	O5-C1	-2.12	1.40	1.43

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	502	NAG	C1-O5-C5	6.29	120.61	112.19
6	D	505	RB0	C2-C3-C4	3.18	117.38	112.48
6	D	505	RB0	O3-C3-C2	-3.12	101.84	108.93
5	C	504	CDP	O2B-PB-O3A	3.10	115.03	104.64
5	B	503	CDP	O4'-C1'-N1	2.86	114.84	108.36
5	D	504	CDP	O4'-C1'-N1	2.64	114.34	108.36
4	D	503	NAG	C1-O5-C5	2.55	115.60	112.19
5	A	504	CDP	O2A-PA-O3A	2.17	113.14	107.27
4	D	502	NAG	O3-C3-C2	-2.16	104.92	109.40
5	B	503	CDP	O2A-PA-O3A	2.04	112.79	107.27

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	504	CDP	C5'-O5'-PA-O3A
5	C	504	CDP	C5'-O5'-PA-O1A
5	C	504	CDP	C5'-O5'-PA-O2A
5	D	504	CDP	C5'-O5'-PA-O3A
5	D	504	CDP	C5'-O5'-PA-O2A
6	D	505	RB0	C2-C3-C4-O4
6	D	505	RB0	C2-C3-C4-C5
6	D	505	RB0	O1-C1-C2-O2
6	D	505	RB0	O1-C1-C2-C3
6	D	505	RB0	O2-C2-C3-O3
6	D	505	RB0	O2-C2-C3-C4
6	A	505	RB0	O3-C3-C4-O4
6	B	504	RB0	O3-C3-C4-O4
6	D	505	RB0	O3-C3-C4-O4
6	D	505	RB0	O3-C3-C4-C5
4	D	503	NAG	C8-C7-N2-C2
4	D	503	NAG	O7-C7-N2-C2
6	D	505	RB0	C1-C2-C3-O3
6	A	505	RB0	O3-C3-C4-C5
4	B	502	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
6	B	504	RB0	O3-C3-C4-C5
4	B	502	NAG	C4-C5-C6-O6
4	A	503	NAG	C4-C5-C6-O6
4	D	502	NAG	C4-C5-C6-O6
6	B	504	RB0	C2-C3-C4-C5
4	D	502	NAG	O5-C5-C6-O6
4	C	503	NAG	O5-C5-C6-O6
4	C	502	NAG	O5-C5-C6-O6
4	A	503	NAG	O5-C5-C6-O6
6	A	505	RB0	C2-C3-C4-O4
6	B	504	RB0	C2-C3-C4-O4
5	C	504	CDP	PB-O3A-PA-O2A
5	B	503	CDP	C5'-O5'-PA-O3A
5	D	504	CDP	PB-O3A-PA-O2A
6	C	505	RB0	O3-C3-C4-C5
6	D	505	RB0	C1-C2-C3-C4
6	C	505	RB0	O2-C2-C3-O3
6	C	505	RB0	O3-C3-C4-O4
6	C	505	RB0	C1-C2-C3-O3
5	B	503	CDP	PB-O3A-PA-O2A
4	A	502	NAG	C1-C2-N2-C7
6	A	505	RB0	C2-C3-C4-C5
5	A	504	CDP	O4'-C4'-C5'-O5'
5	C	504	CDP	O4'-C4'-C5'-O5'

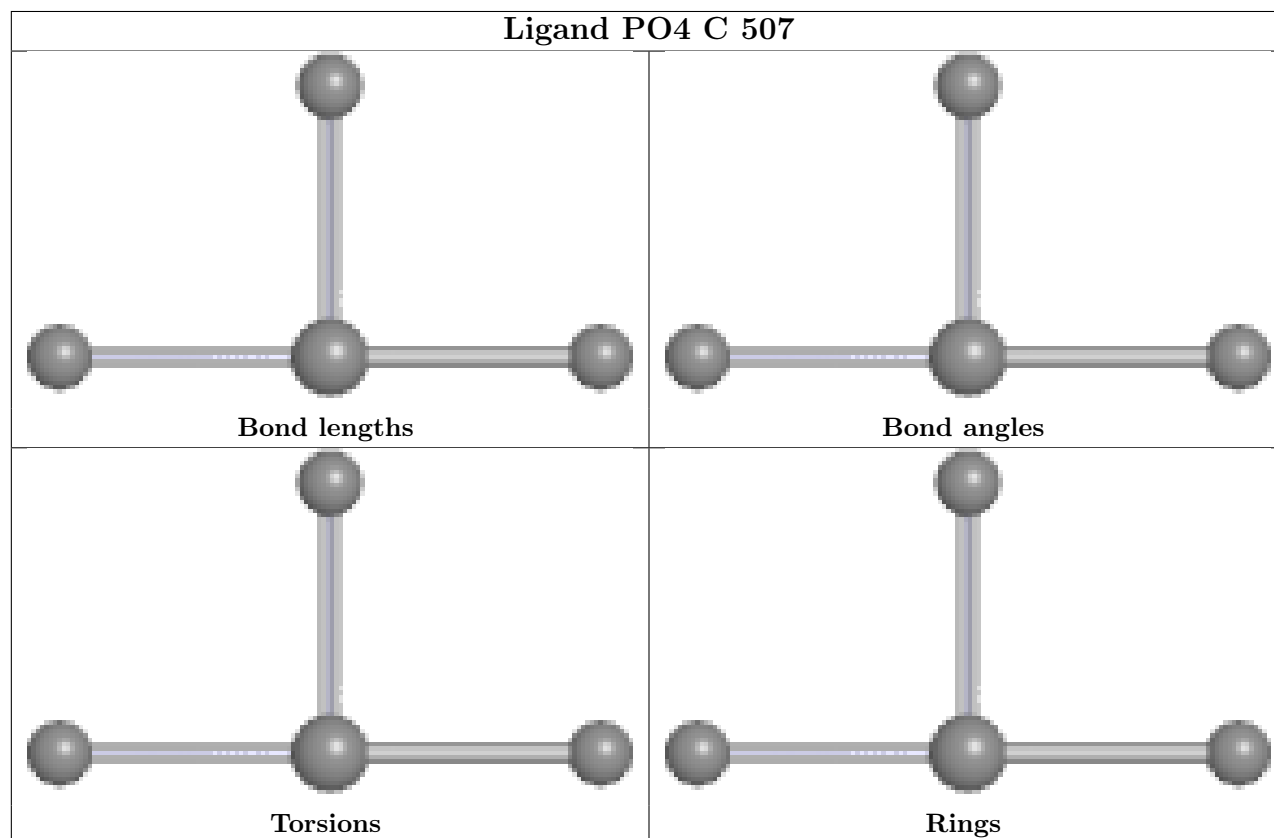
There are no ring outliers.

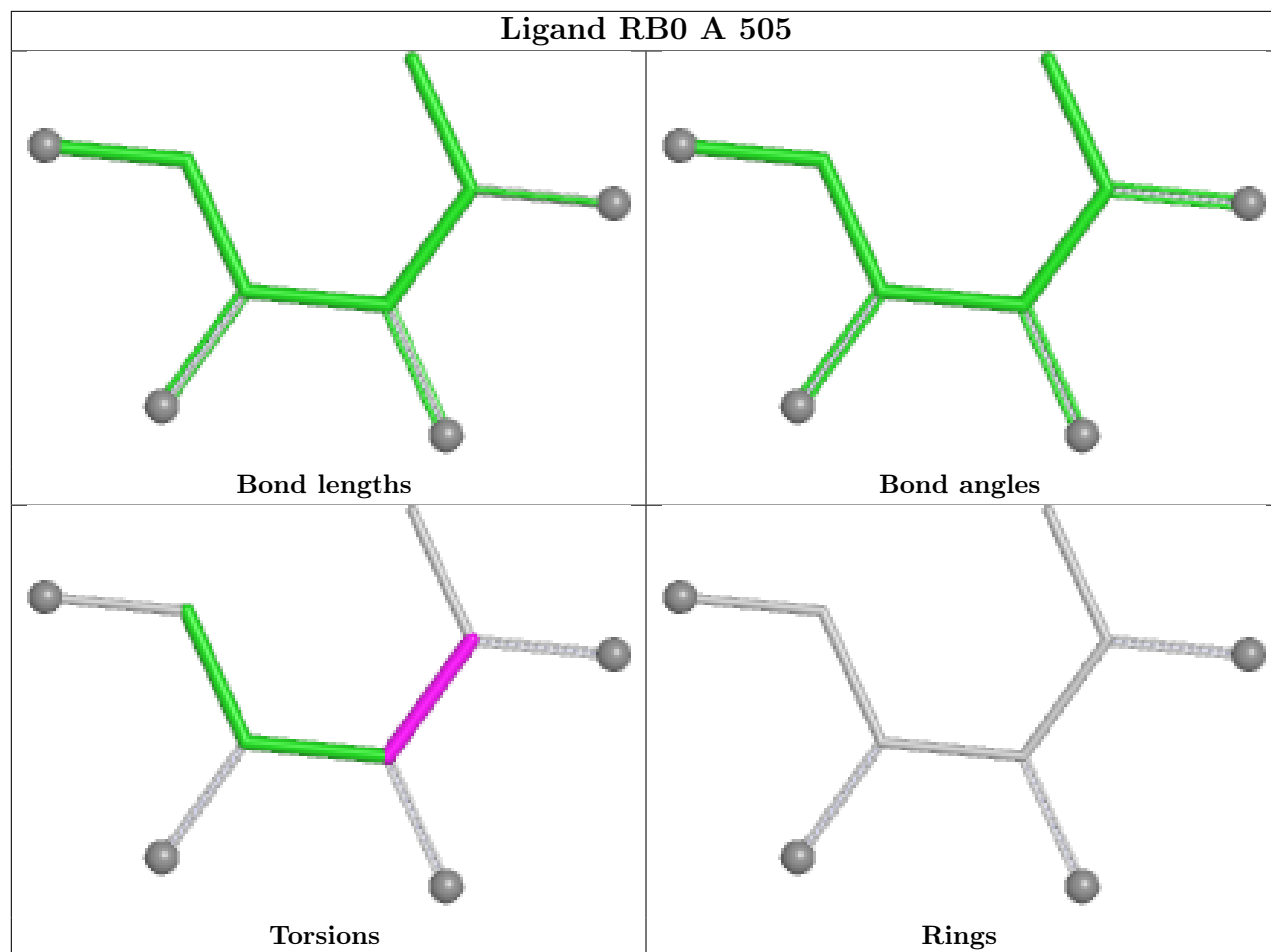
7 monomers are involved in 8 short contacts:

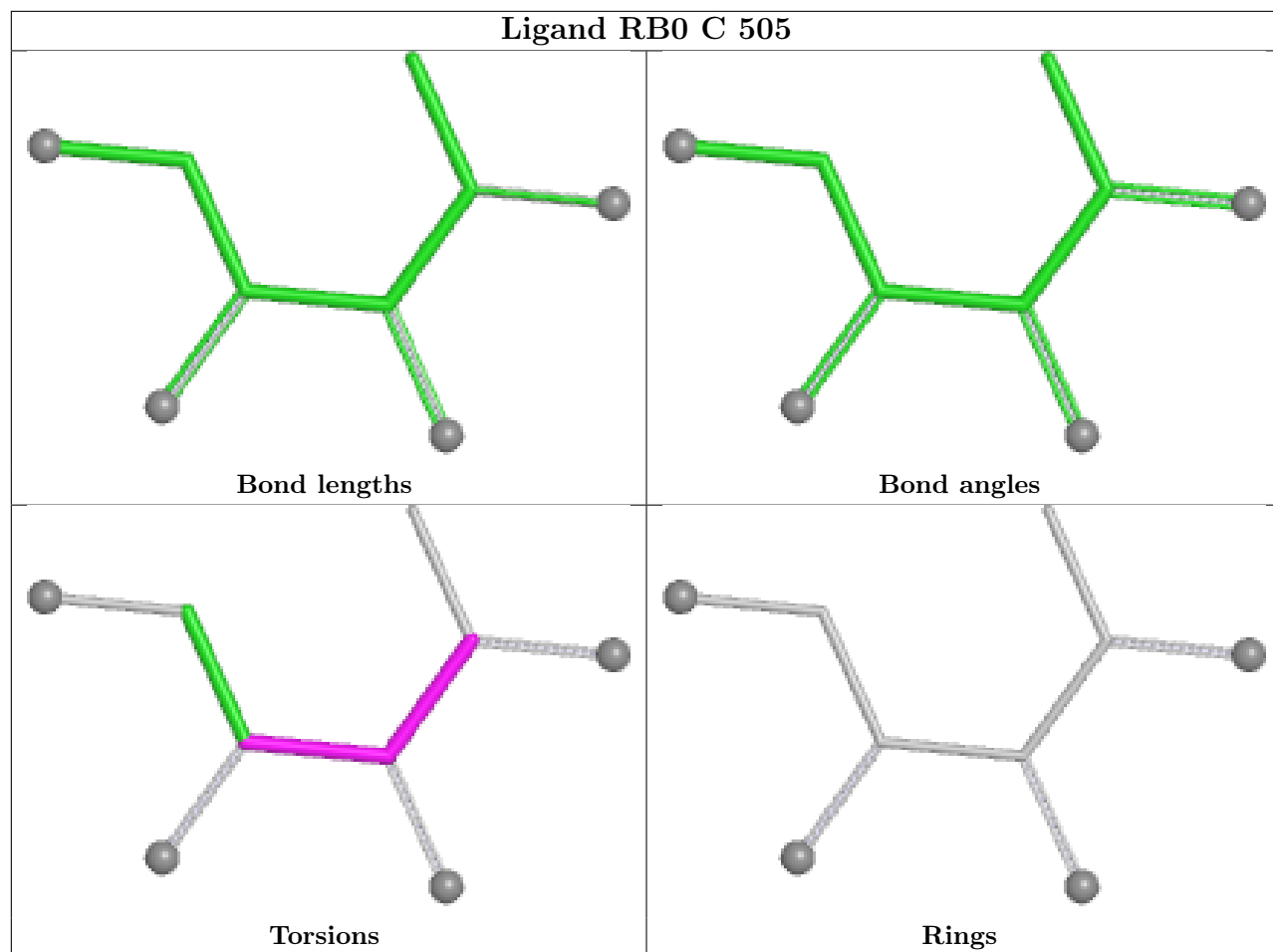
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	C	507	PO4	2	0
6	C	505	RB0	1	0
4	A	502	NAG	1	0
5	A	504	CDP	1	0
6	D	505	RB0	1	0
5	C	504	CDP	1	0
4	D	503	NAG	1	0

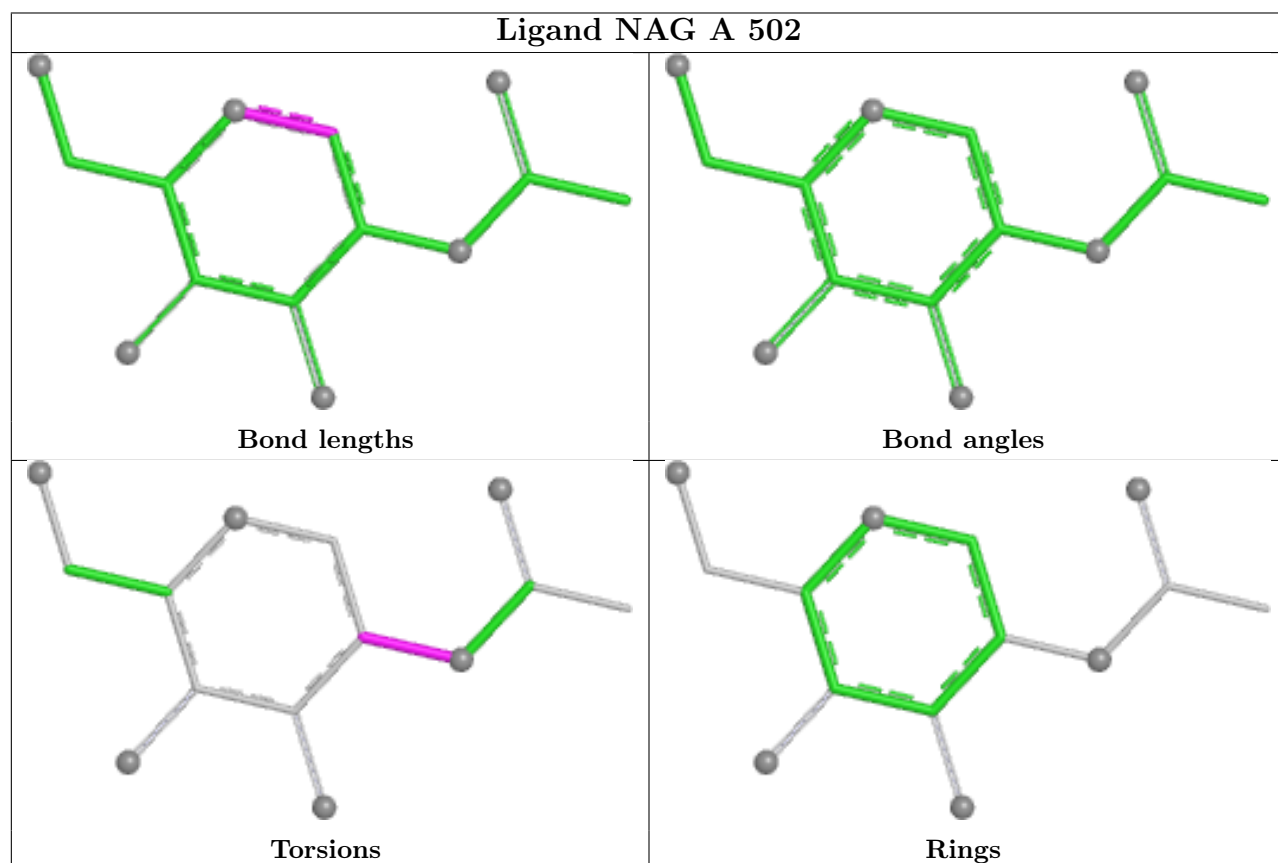
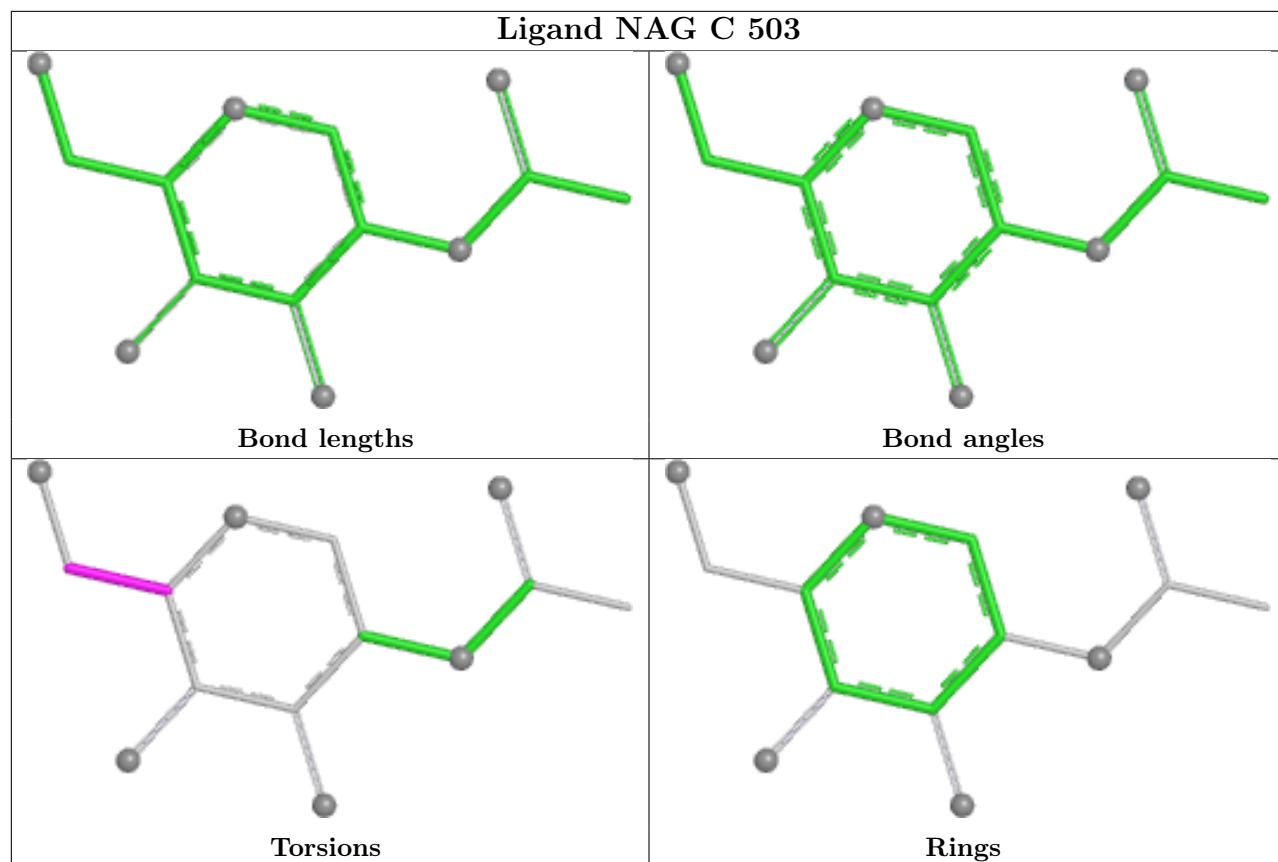
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

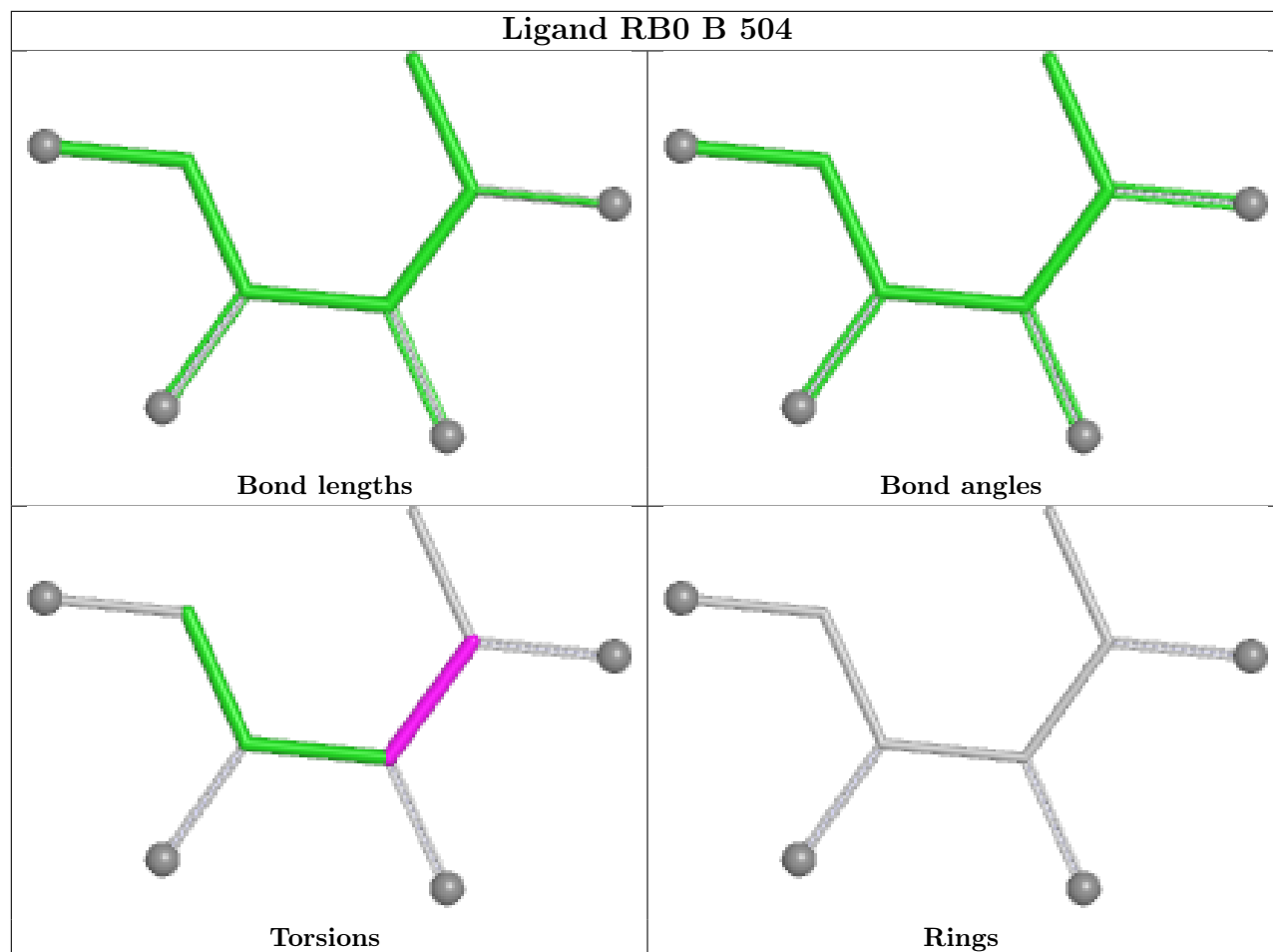




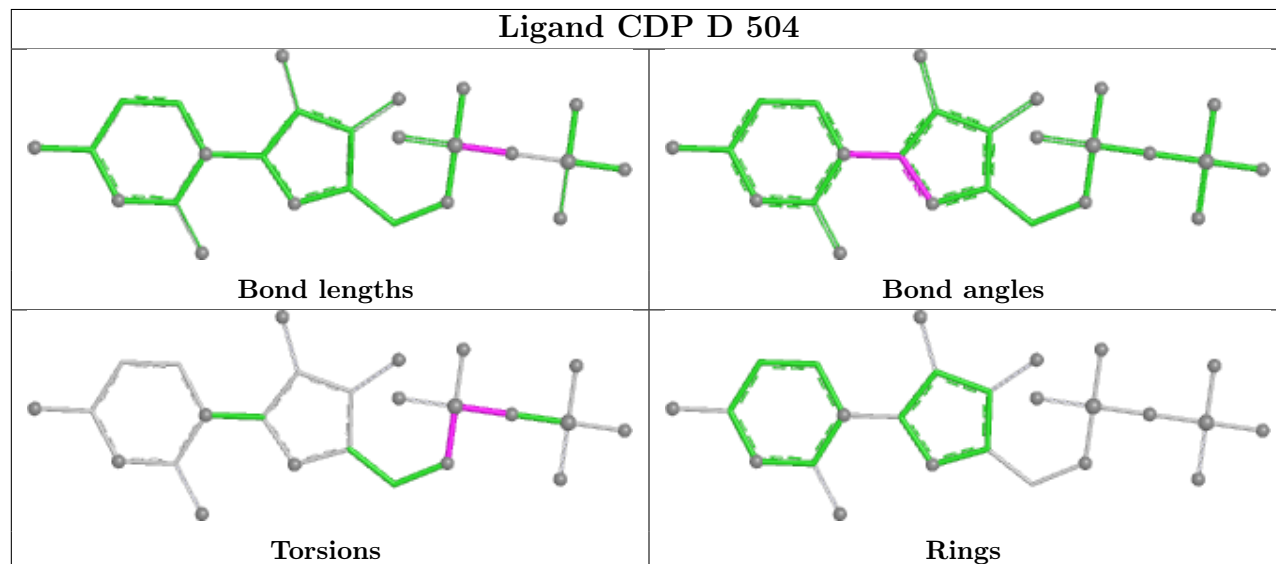


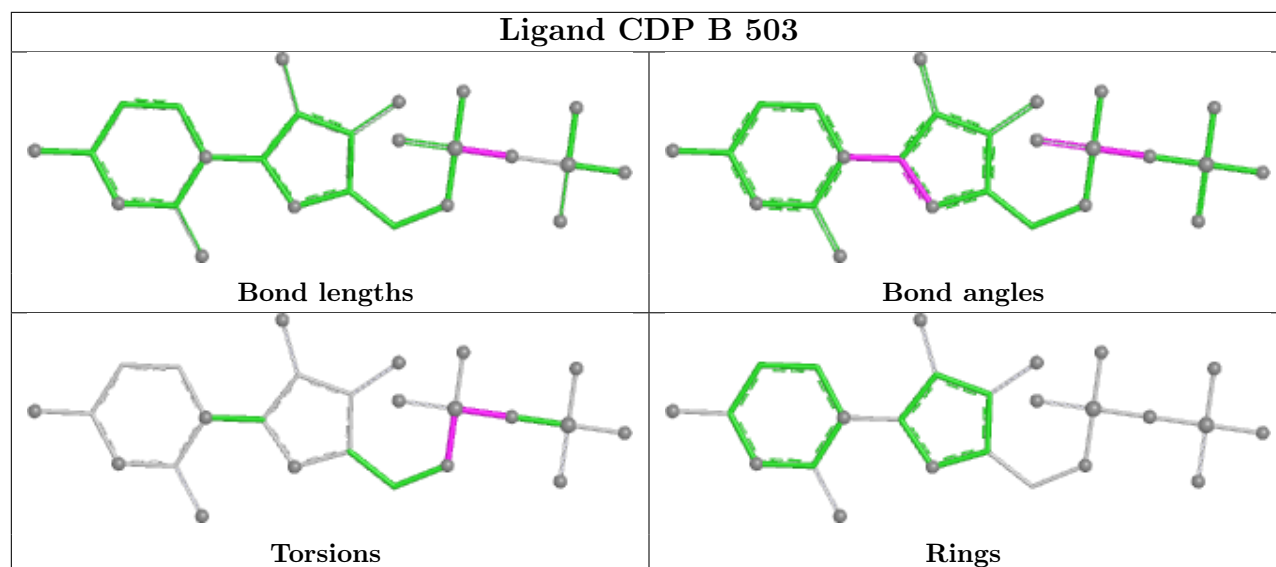
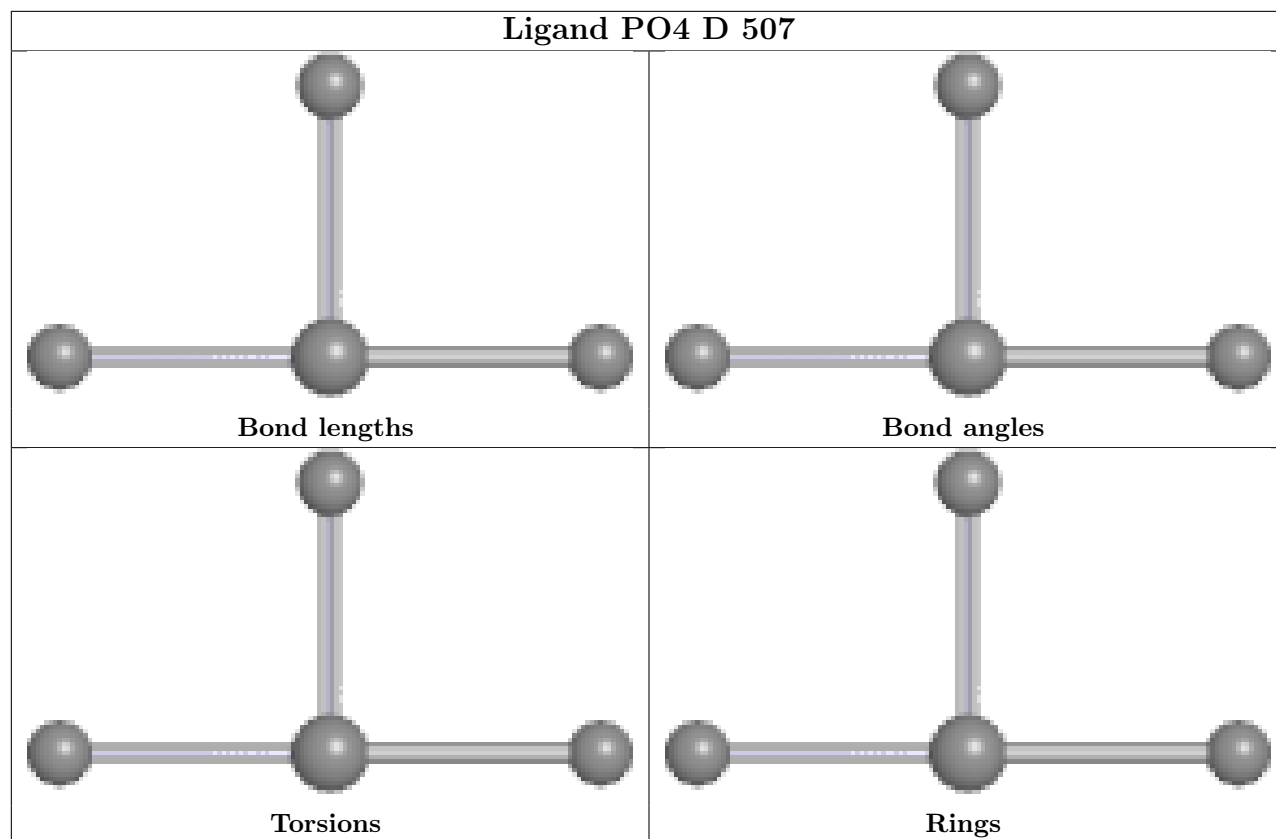


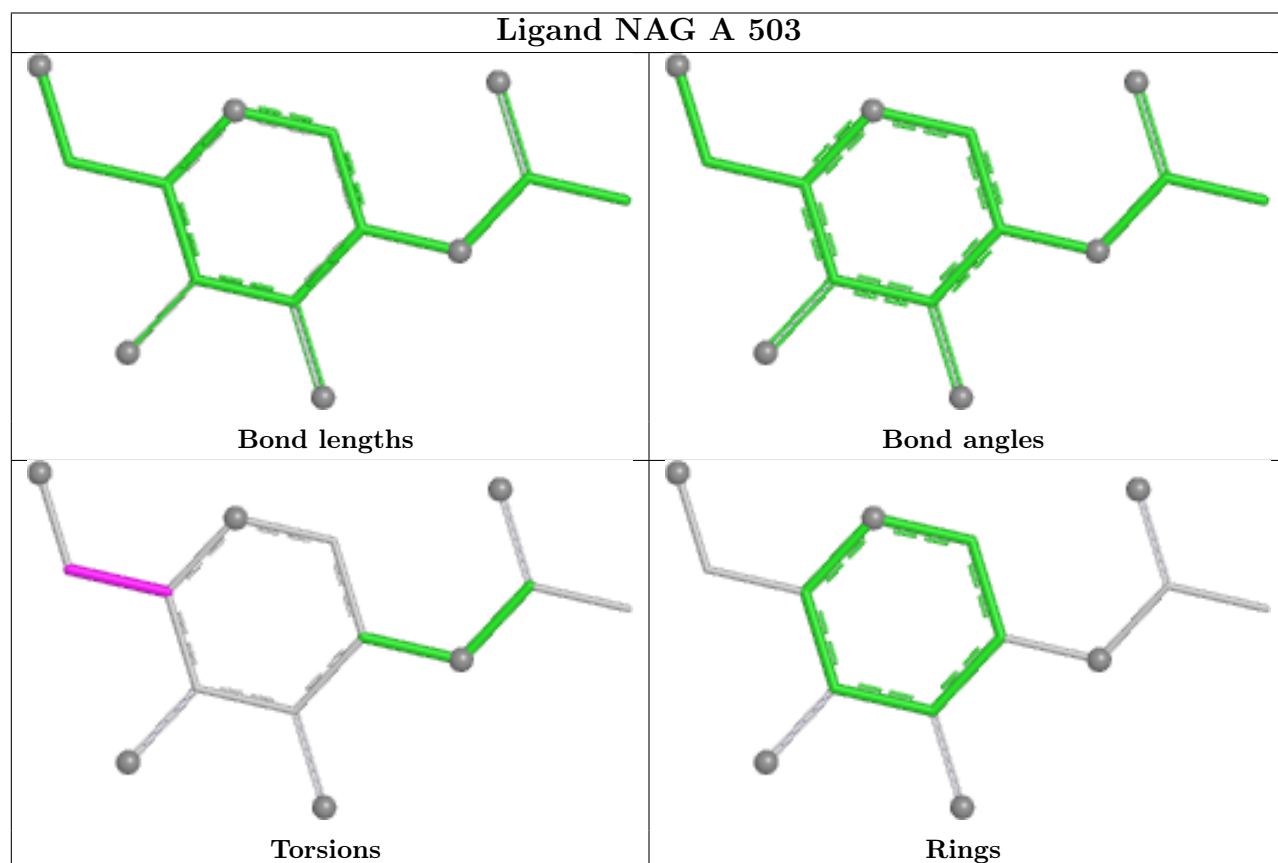
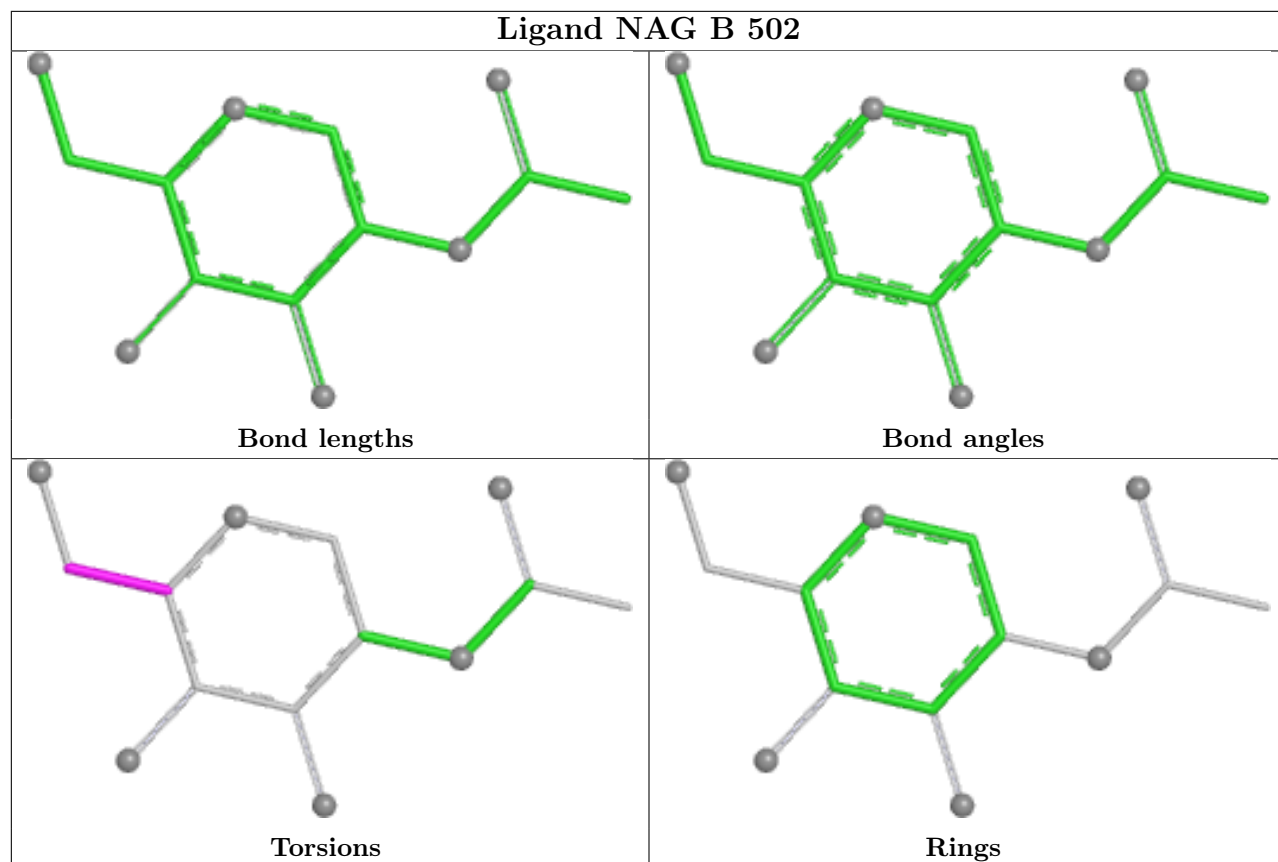
Ligand RB0 B 504

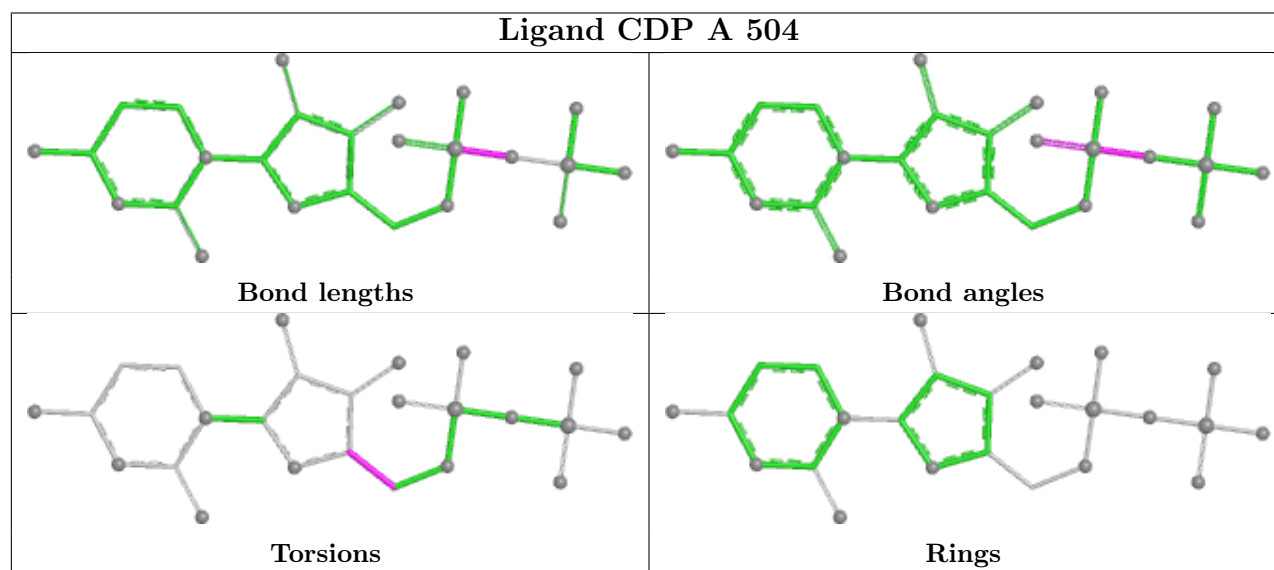
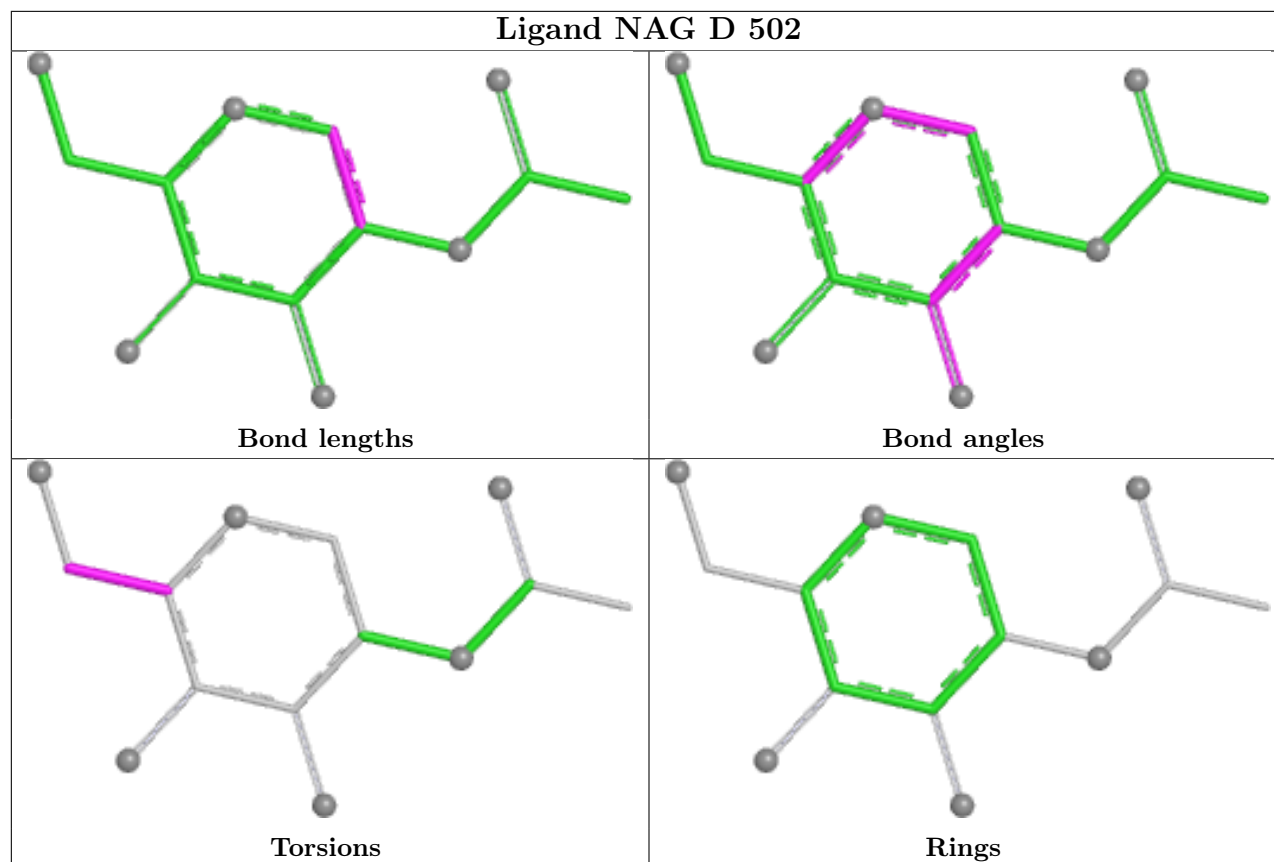


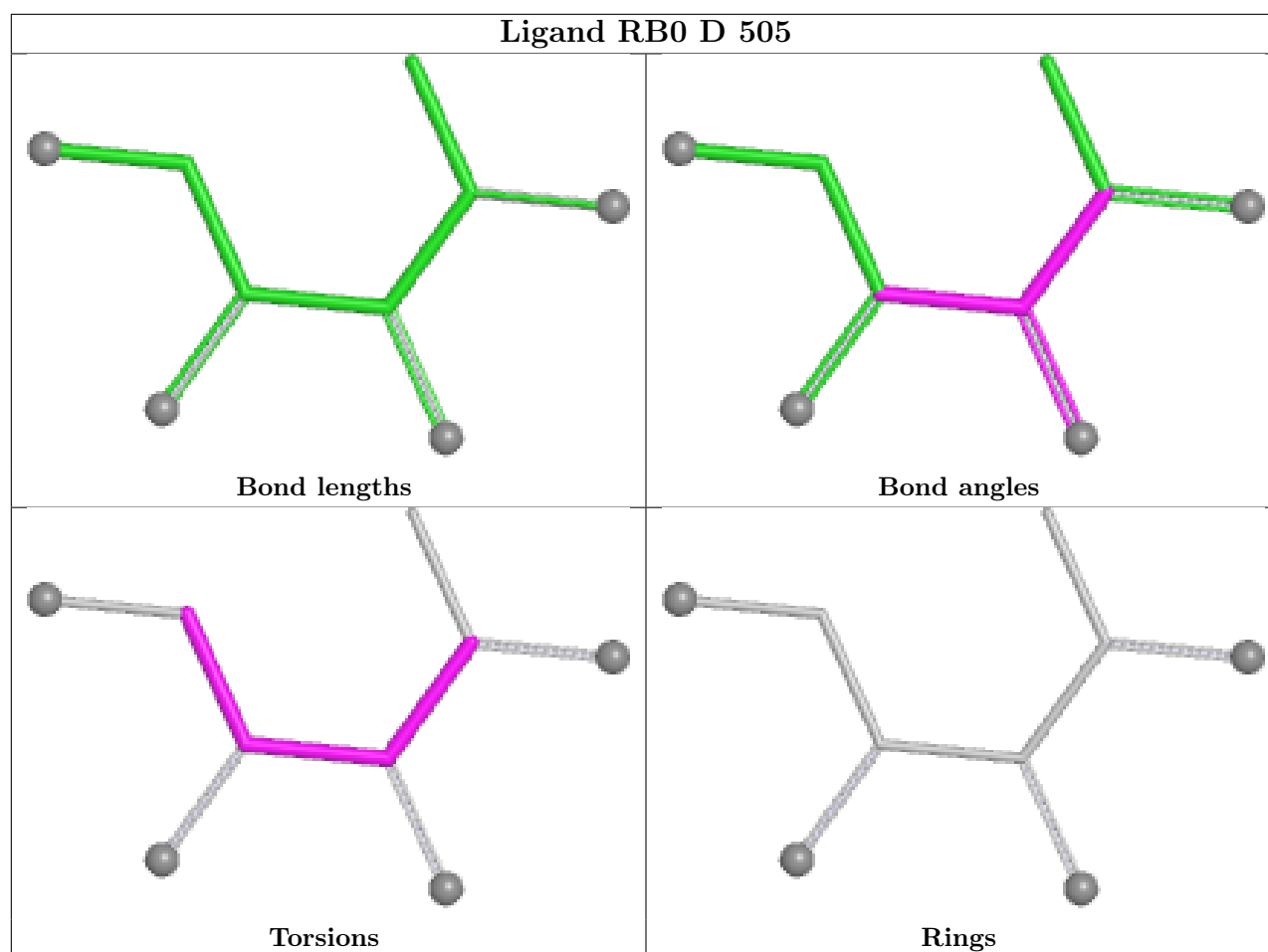
Ligand CDP D 504

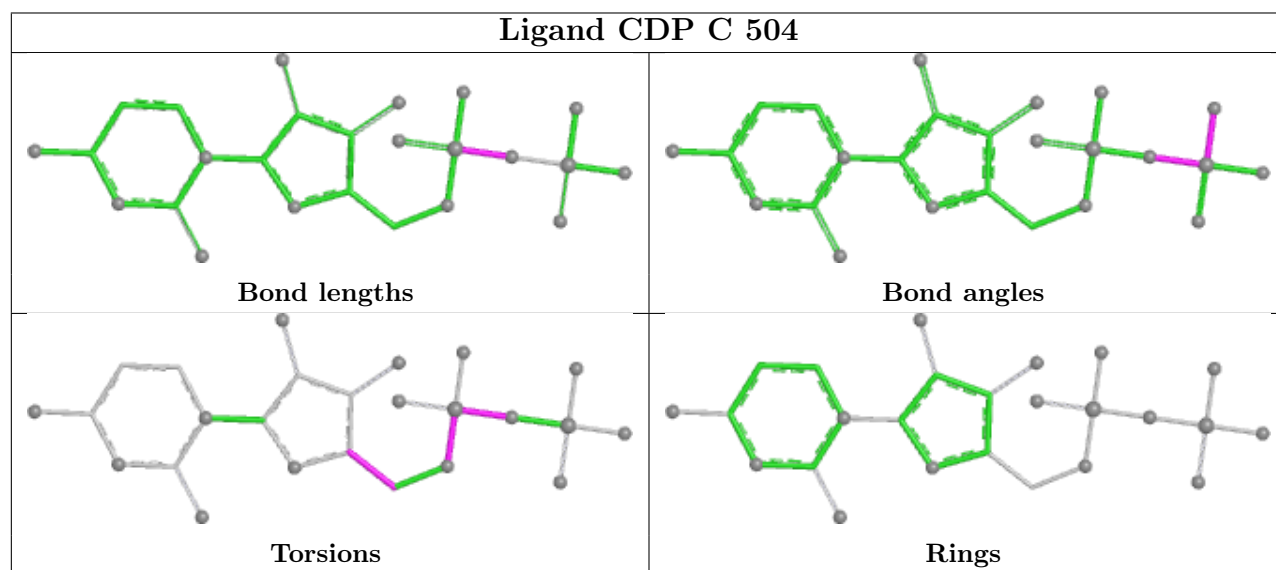
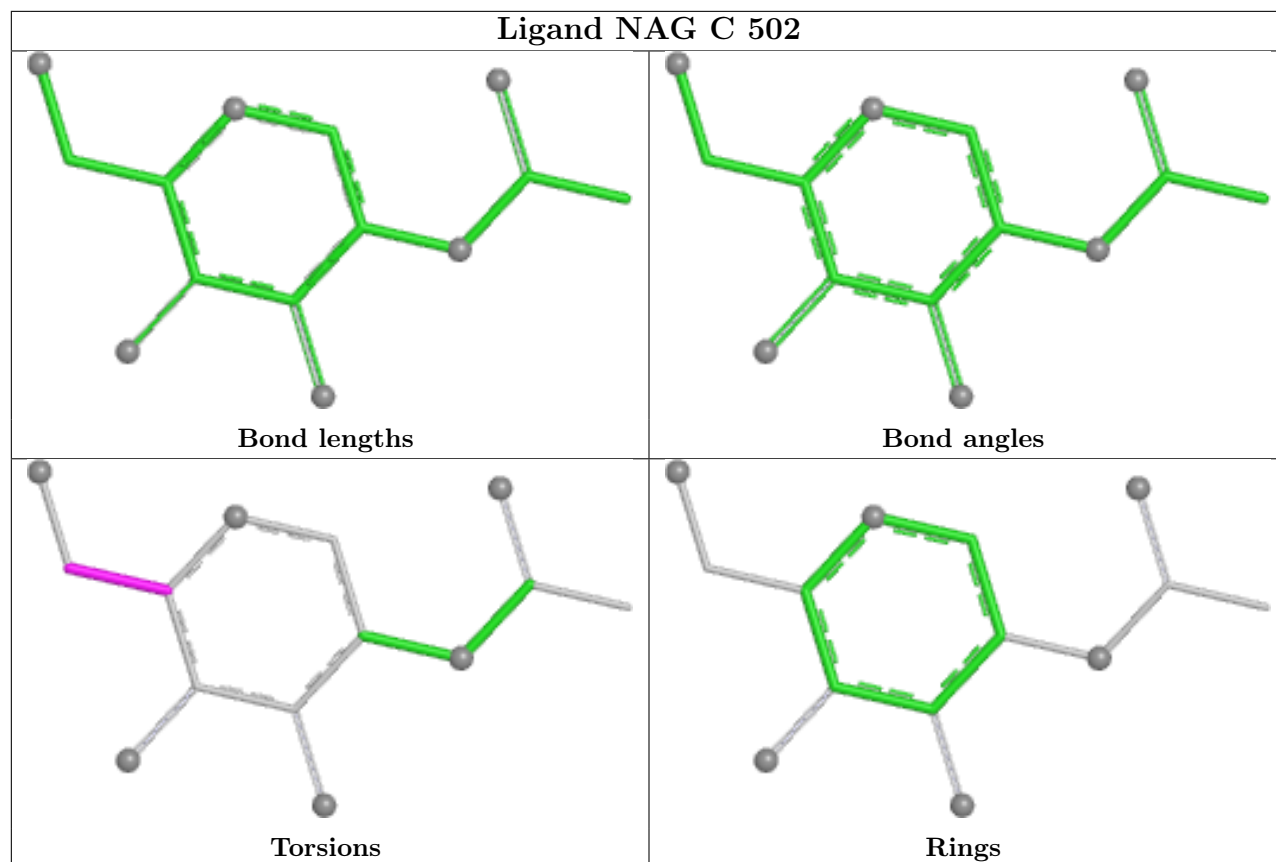


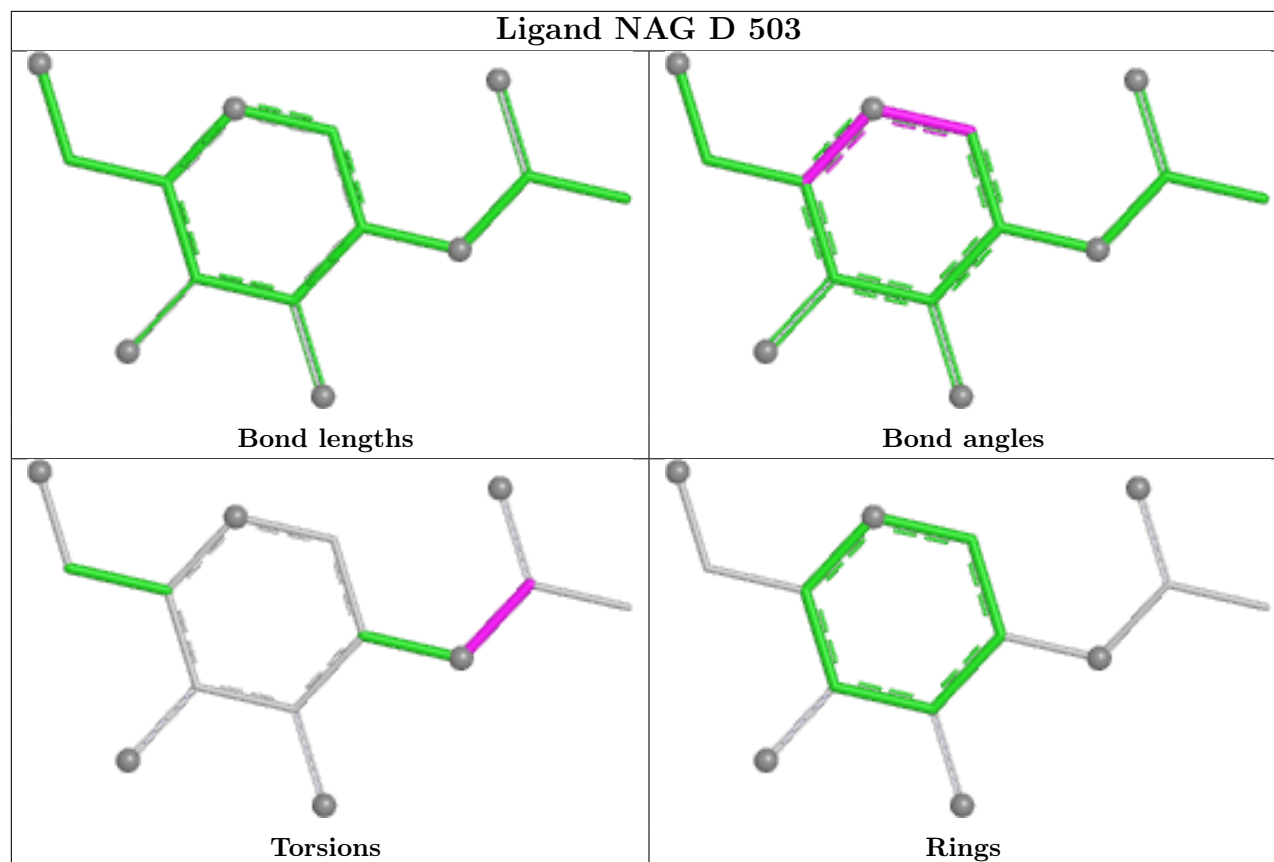












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	446/455 (98%)	0.24	14 (3%) 51 52	34, 73, 114, 151	2 (0%)
1	B	445/455 (97%)	0.18	4 (0%) 81 82	30, 68, 107, 153	1 (0%)
1	C	445/455 (97%)	0.32	20 (4%) 38 37	48, 74, 117, 152	0
1	D	443/455 (97%)	0.04	6 (1%) 73 75	45, 67, 100, 135	0
All	All	1779/1820 (97%)	0.19	44 (2%) 58 61	30, 70, 110, 153	3 (0%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	483	TYR	7.1
1	C	487	ALA	5.7
1	D	47	PRO	5.2
1	C	491	LEU	4.7
1	B	47	PRO	4.4
1	C	486	PRO	4.3
1	A	398	VAL	4.3
1	A	488	LEU	4.0
1	C	484	PRO	4.0
1	C	47	PRO	3.8
1	D	400	GLY	3.7
1	D	397	ALA	3.6
1	C	488	LEU	3.6
1	A	487	ALA	3.5
1	A	433	LEU	3.5
1	C	398	VAL	3.3
1	C	168	CYS	3.3
1	A	430	ASP	3.1
1	C	489	LEU	3.1
1	C	146	GLU	3.0
1	A	491	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
1	D	238	LEU	2.8
1	C	399	GLU	2.8
1	C	137	ALA	2.8
1	A	399	GLU	2.7
1	A	489	LEU	2.6
1	C	439	VAL	2.6
1	A	370	GLU	2.6
1	C	490	SER	2.6
1	A	400	GLY	2.6
1	B	255	TRP	2.5
1	C	138	PRO	2.5
1	C	379	ARG	2.5
1	A	137	ALA	2.5
1	D	68	ASP	2.4
1	C	396	LYS	2.4
1	D	378	LEU	2.3
1	A	439	VAL	2.3
1	A	46	GLY	2.3
1	C	482	GLN	2.3
1	A	168	CYS	2.3
1	B	138	PRO	2.3
1	C	378	LEU	2.2
1	B	48	ARG	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.

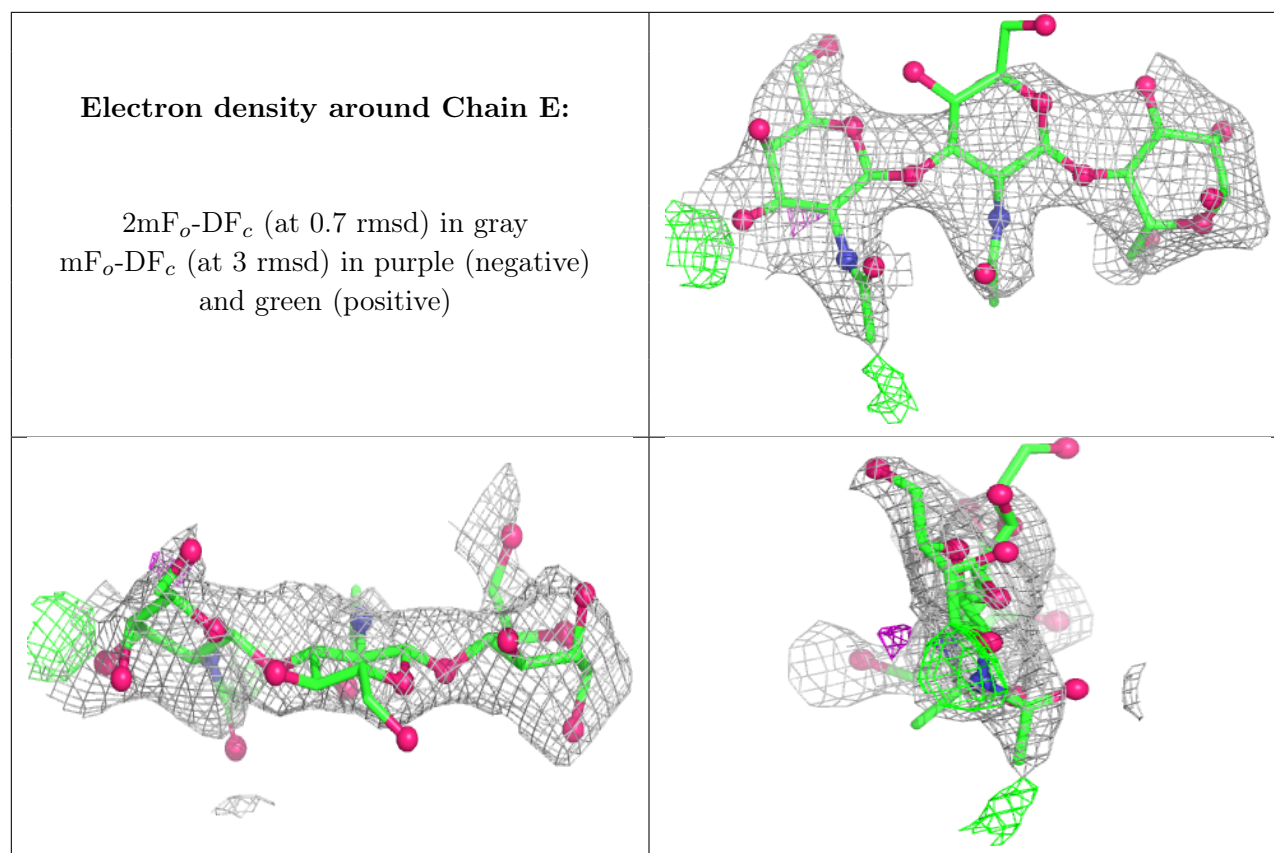
6.3 Carbohydrates [i](#)

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
2	MAN	E	1	12/12	0.51	0.12	149,153,158,161	0
2	NAG	E	2	14/15	0.71	0.15	153,161,168,171	0
2	NGA	E	3	14/15	0.73	0.16	133,142,154,162	0

The following is a graphical depiction of the model fit to experimental electron density for oligosac-

charide. Each fit is shown from different orientation to approximate a three-dimensional view.



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
8	PO4	D	507	4/5	0.44	0.11	154,155,155,155	0
4	NAG	D	503	14/15	0.60	0.17	80,85,87,88	0
4	NAG	D	502	14/15	0.62	0.16	82,89,97,100	0
4	NAG	A	503	14/15	0.63	0.16	77,81,86,86	0
4	NAG	C	503	14/15	0.65	0.18	123,130,138,140	0
4	NAG	B	502	14/15	0.73	0.13	80,89,98,103	0
8	PO4	C	507	4/5	0.77	0.26	167,168,170,170	0
4	NAG	A	502	14/15	0.80	0.12	76,79,87,91	0
6	RB0	A	505	9/10	0.82	0.17	101,105,108,109	0
6	RB0	D	505	9/10	0.83	0.20	83,85,87,88	0
6	RB0	C	505	9/10	0.84	0.15	93,96,99,100	0
7	BA	A	506	1/1	0.84	0.16	252,252,252,252	0

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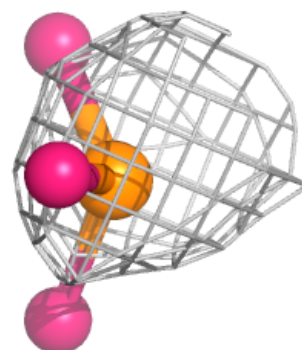
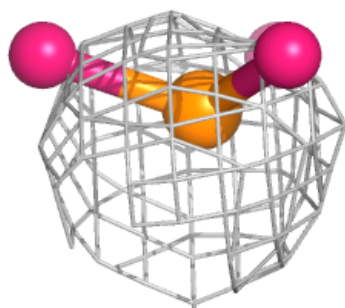
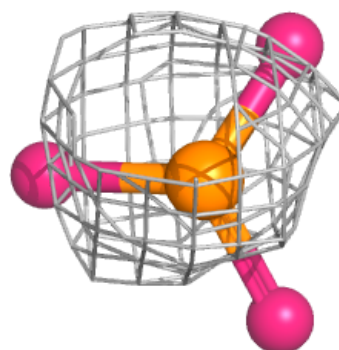
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	RB0	B	504	9/10	0.86	0.21	86,94,98,101	0
4	NAG	C	502	14/15	0.87	0.12	79,87,100,101	0
7	BA	D	506	1/1	0.90	0.14	213,213,213,213	0
7	BA	B	505	1/1	0.92	0.14	177,177,177,177	0
5	CDP	A	504	25/25	0.92	0.09	77,82,96,98	0
5	CDP	C	504	25/25	0.93	0.09	70,75,99,99	0
7	BA	C	506	1/1	0.94	0.12	224,224,224,224	0
5	CDP	D	504	25/25	0.94	0.08	56,59,77,82	0
5	CDP	B	503	25/25	0.96	0.06	52,56,66,77	0
3	ZN	B	501	1/1	0.99	0.02	59,59,59,59	0
3	ZN	C	501	1/1	0.99	0.04	60,60,60,60	0
3	ZN	D	501	1/1	0.99	0.02	56,56,56,56	0
3	ZN	A	501	1/1	0.99	0.03	57,57,57,57	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.

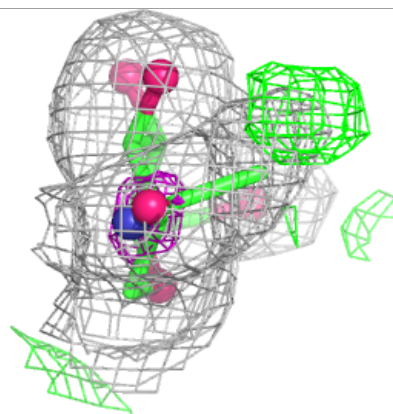
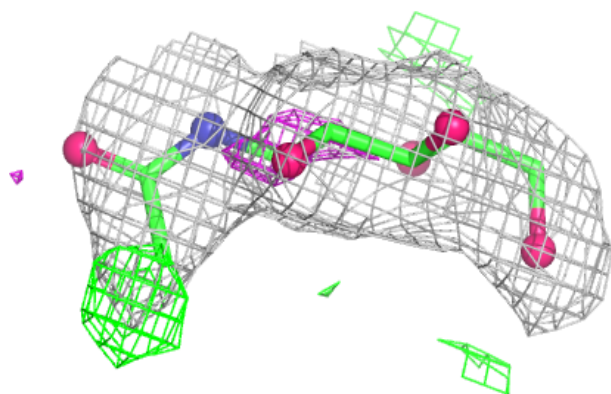
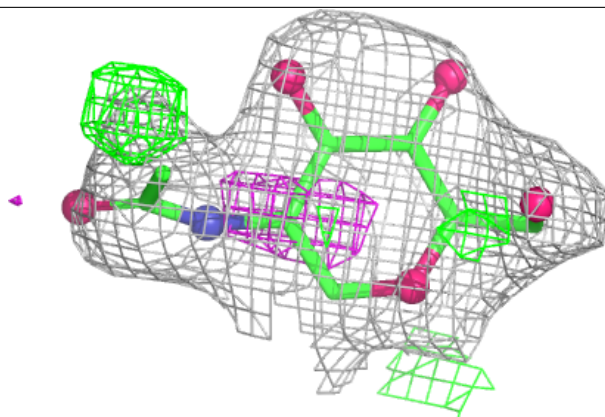
Electron density around PO4 D 507:

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

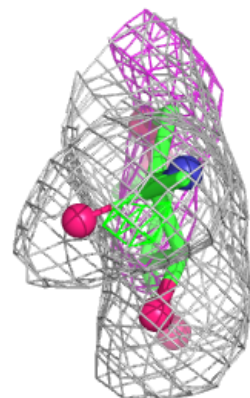
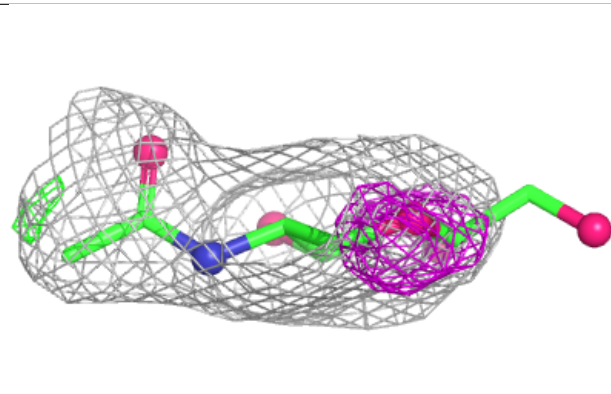
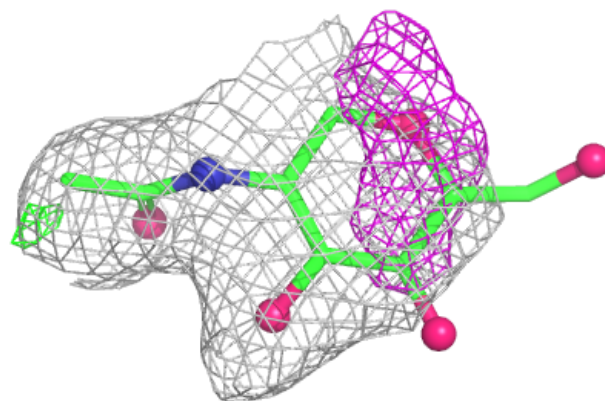


Electron density around NAG D 503:

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

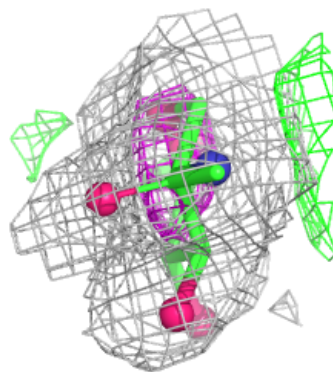
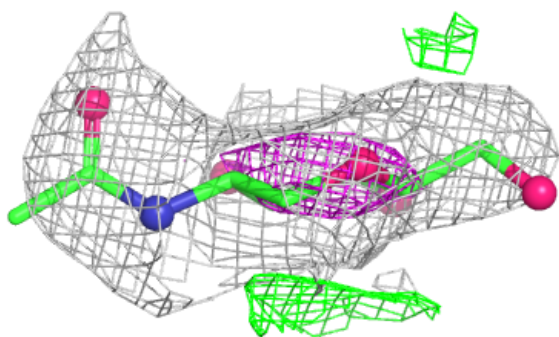
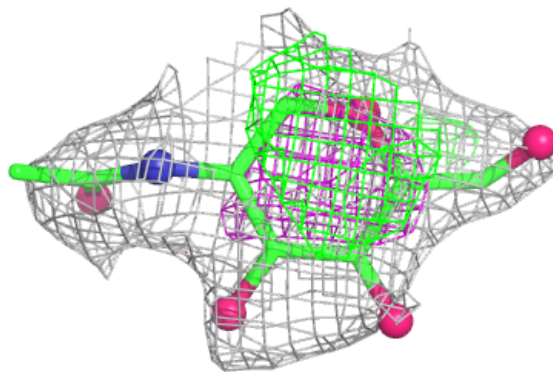
**Electron density around NAG D 502:**

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and green (positive)

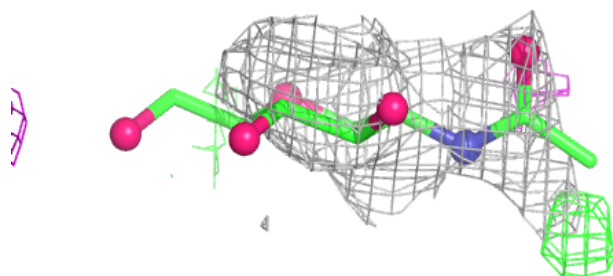
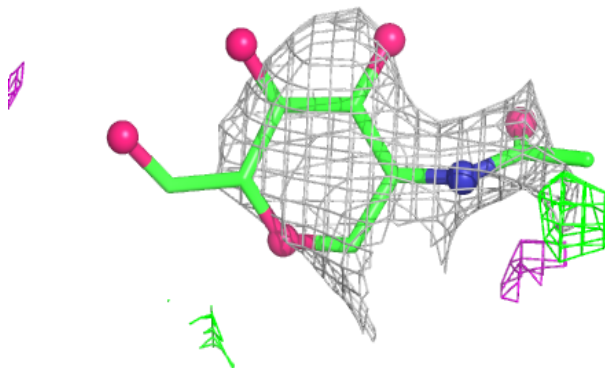


Electron density around NAG A 503:

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

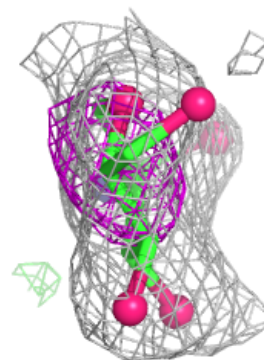
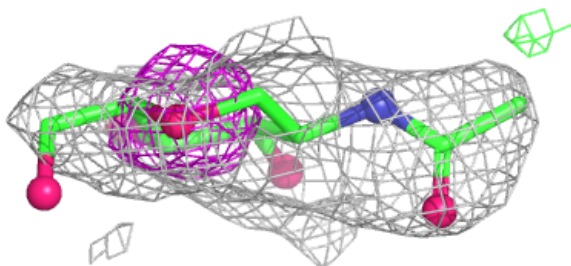
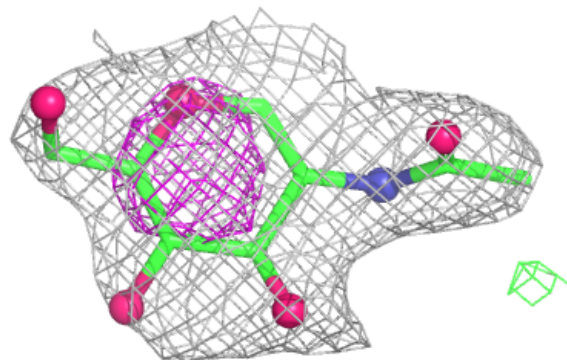
**Electron density around NAG C 503:**

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



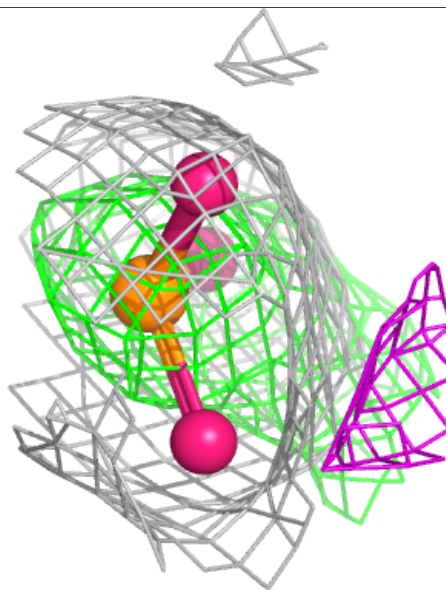
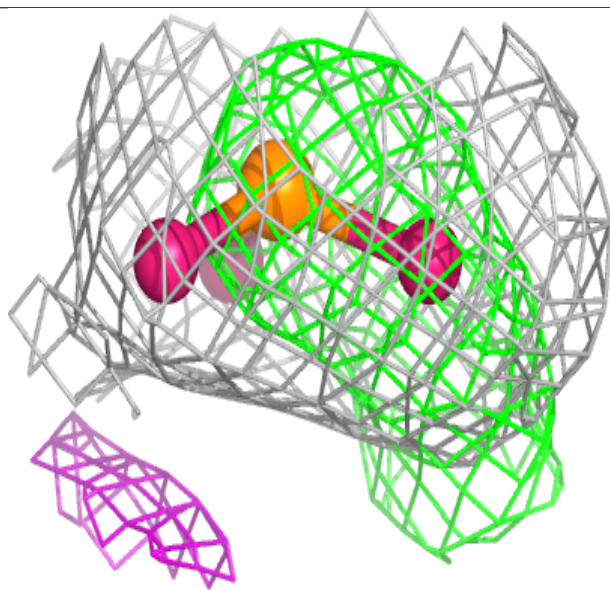
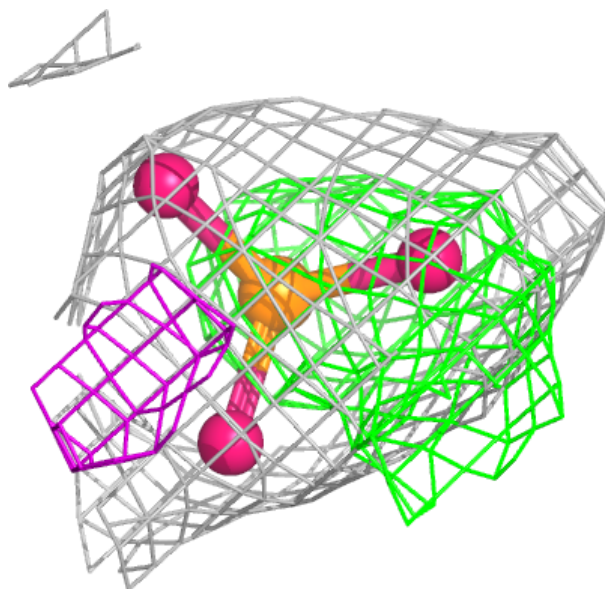
Electron density around NAG B 502:

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



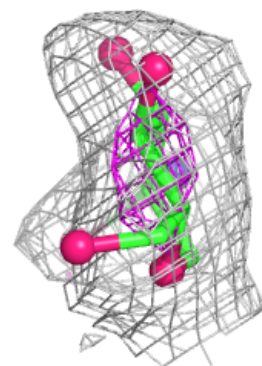
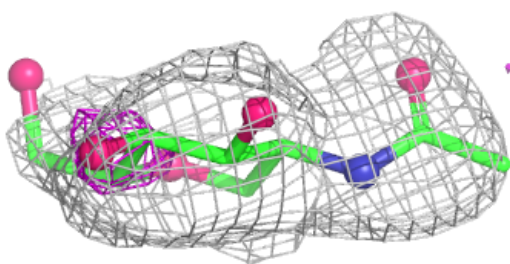
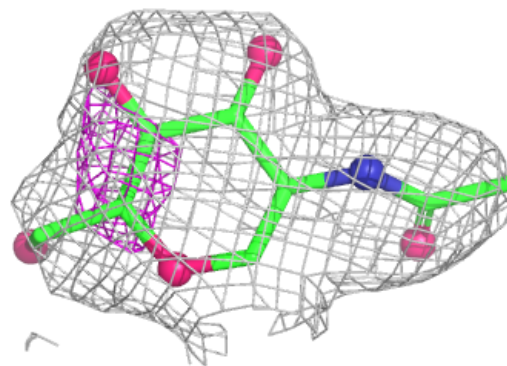
Electron density around PO4 C 507:

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

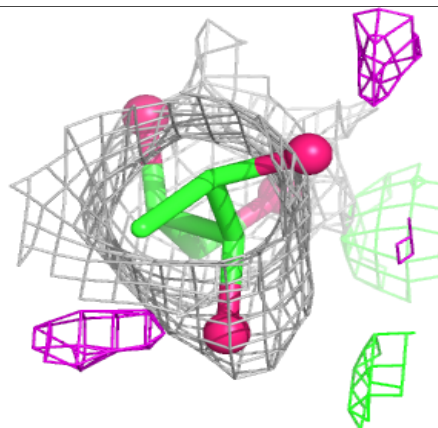
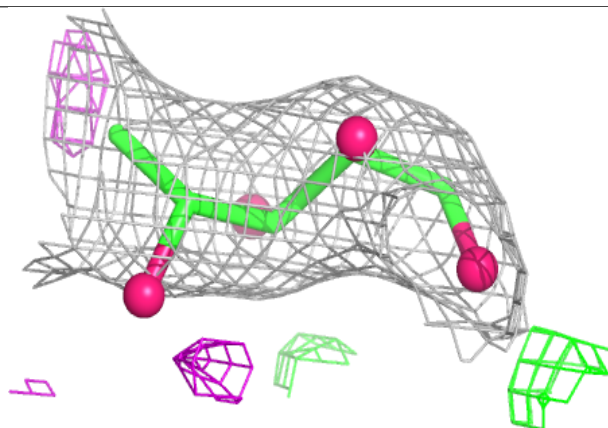
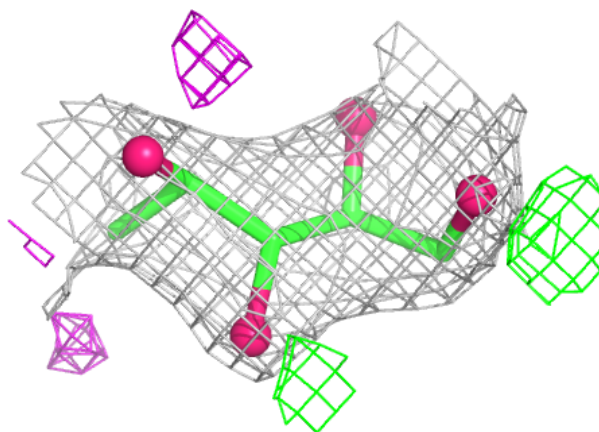


Electron density around NAG A 502:

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

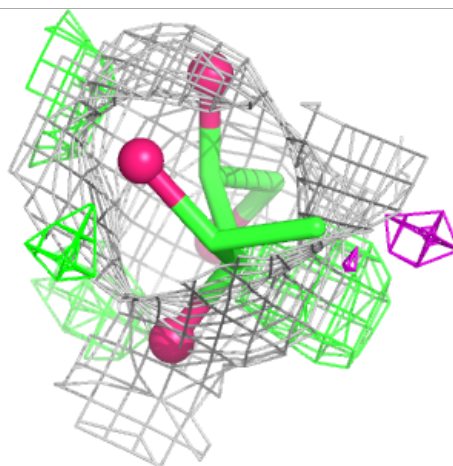
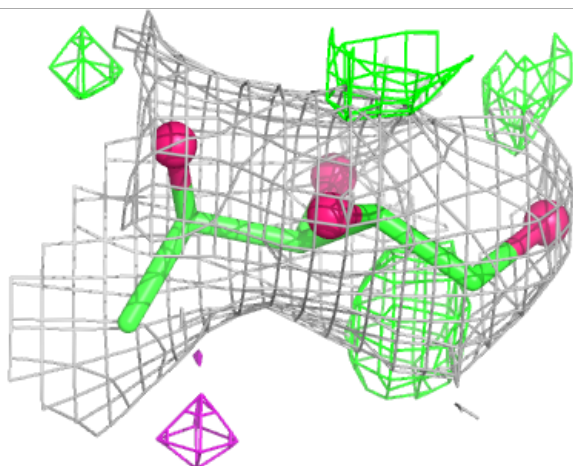
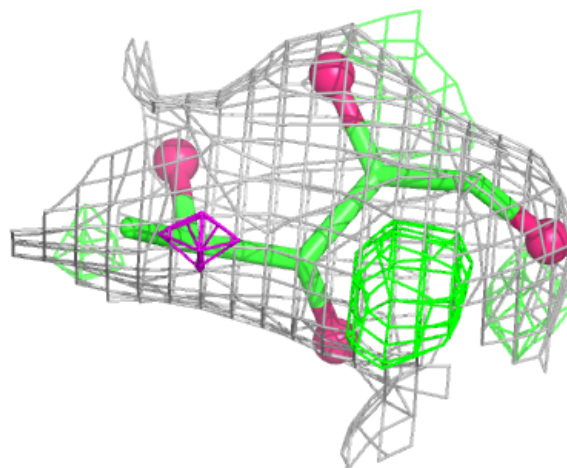
**Electron density around RB0 A 505:**

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



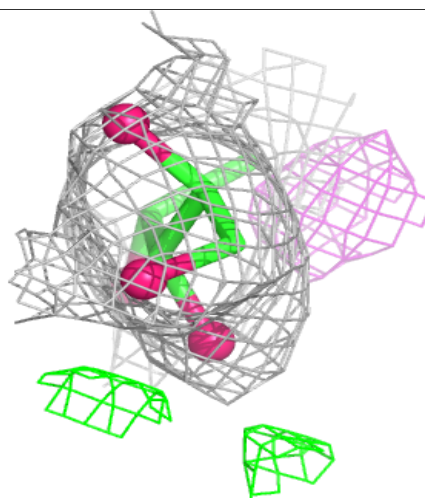
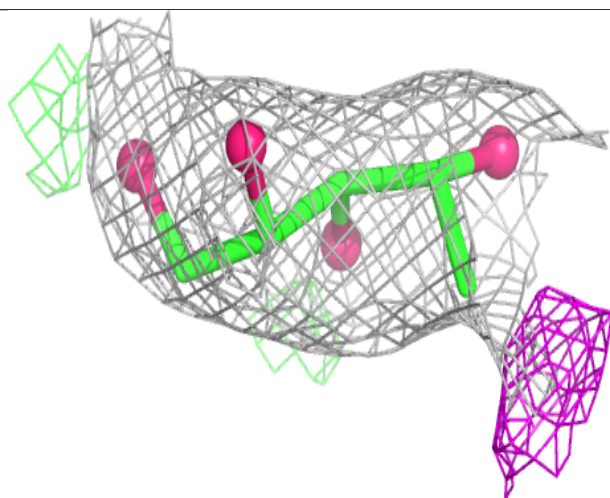
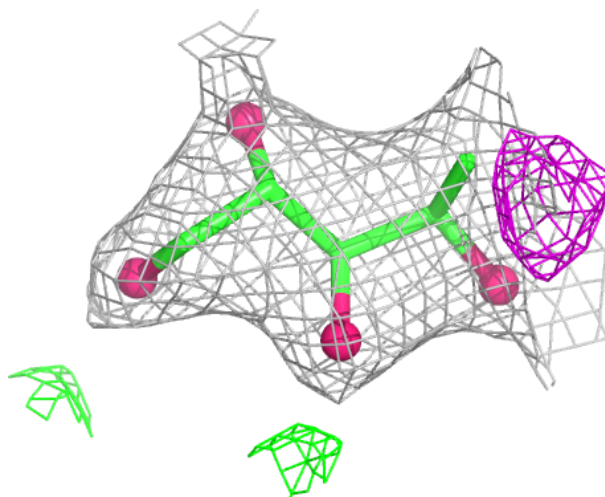
Electron density around RB0 D 505:

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



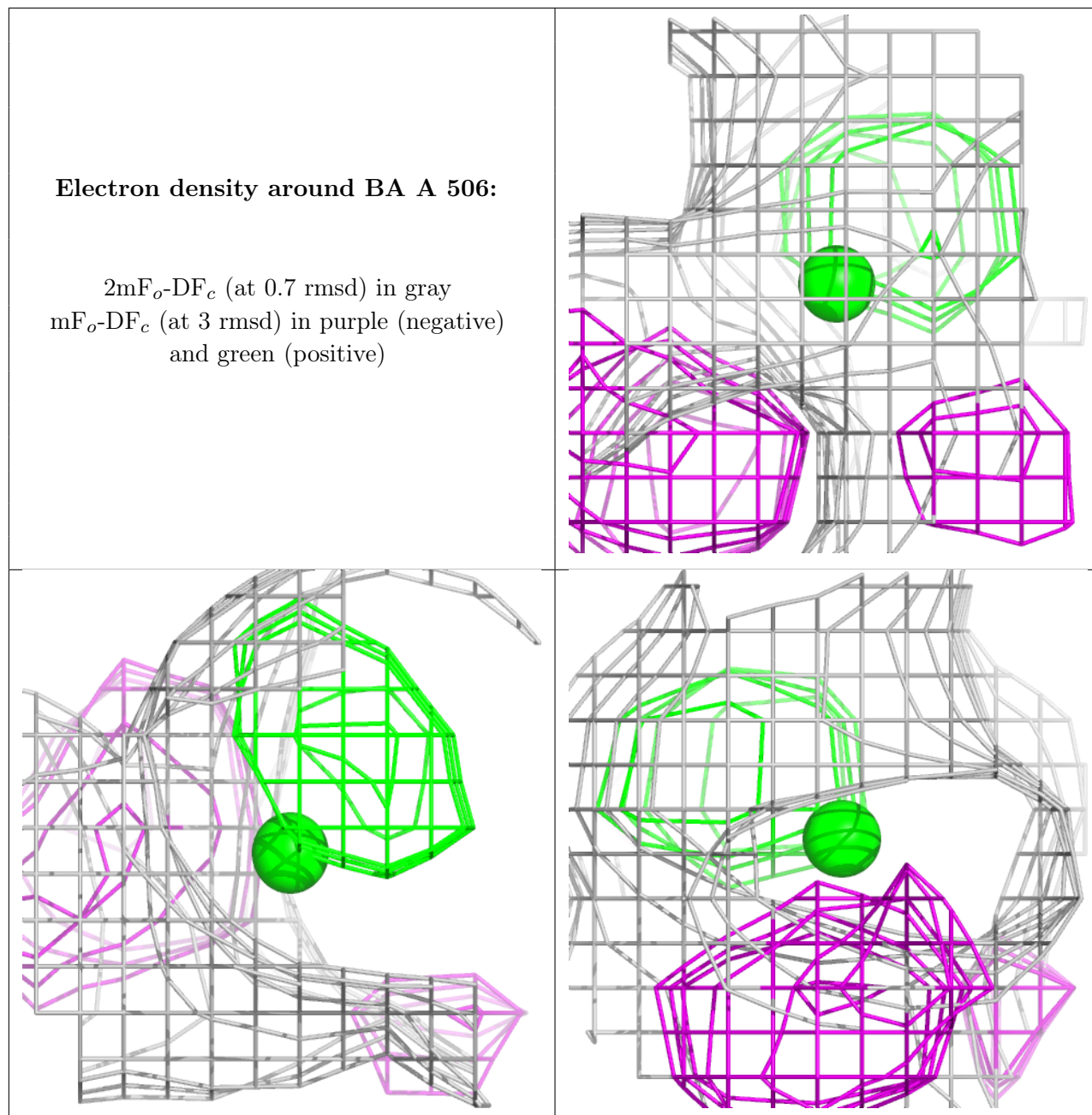
Electron density around RB0 C 505:

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



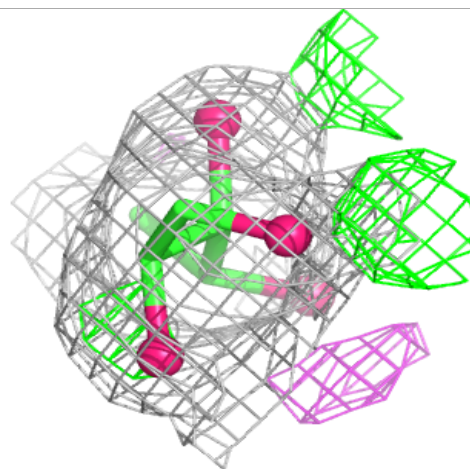
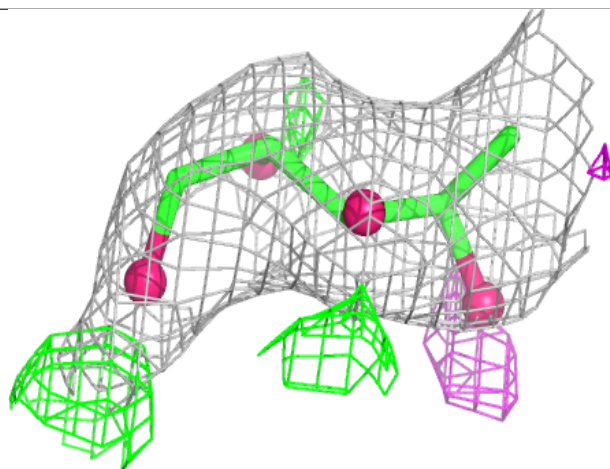
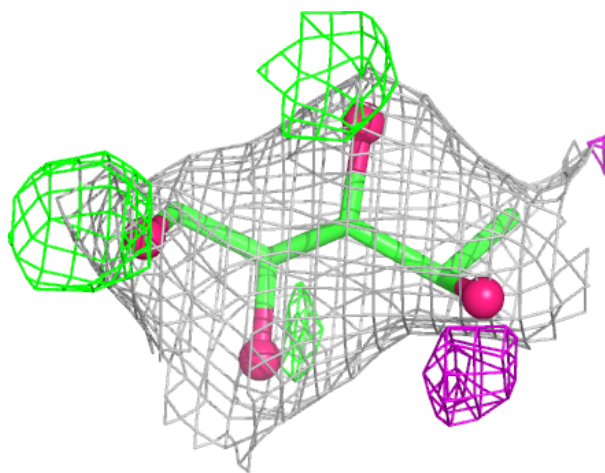
Electron density around BA A 506:

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



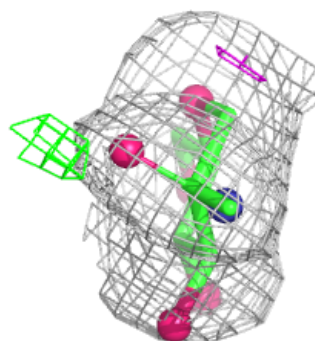
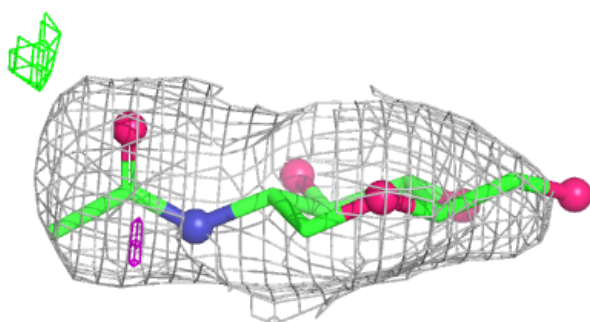
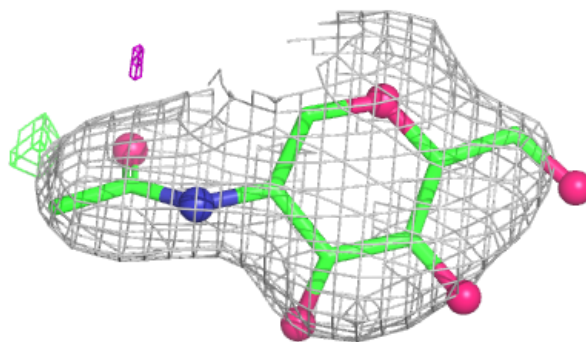
Electron density around RB0 B 504:

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



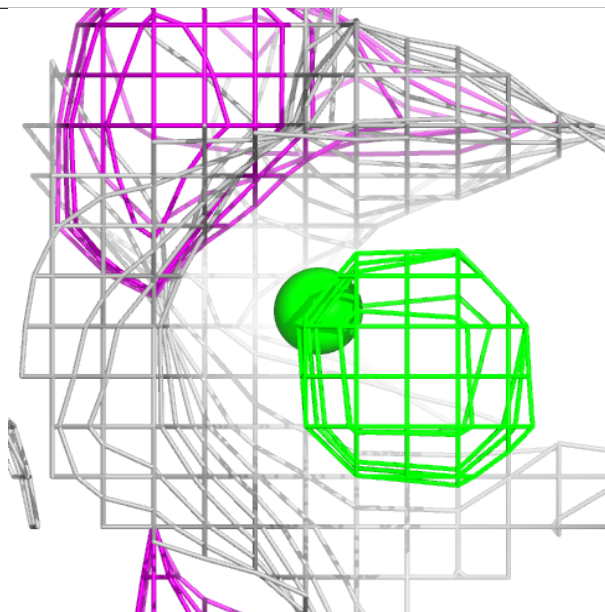
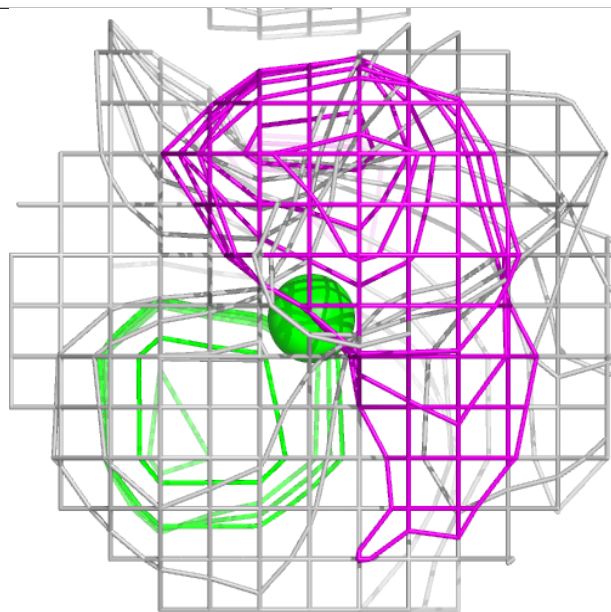
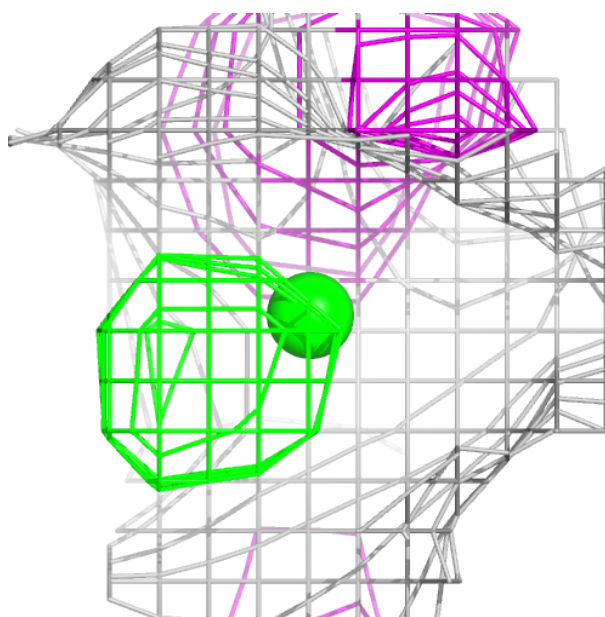
Electron density around NAG C 502:

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



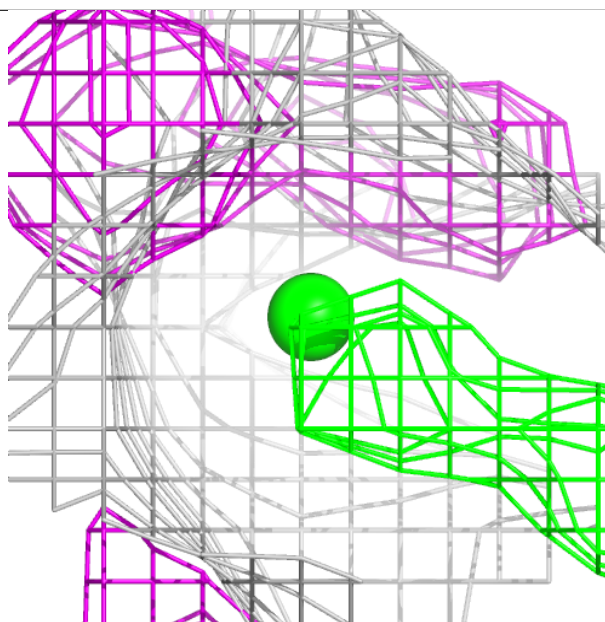
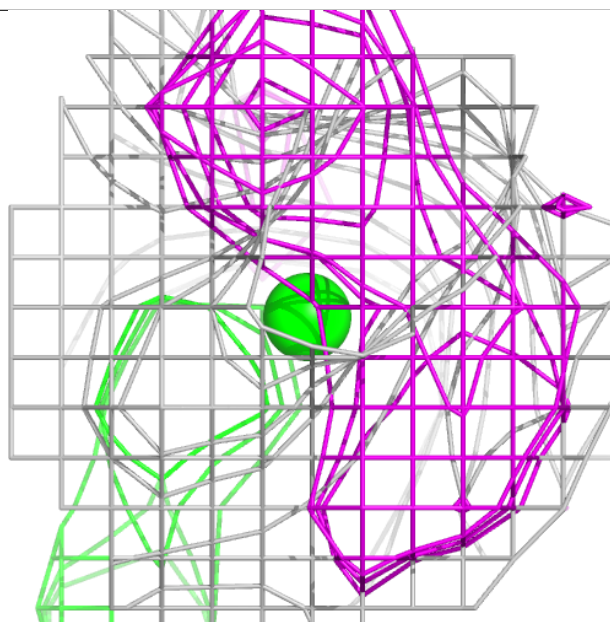
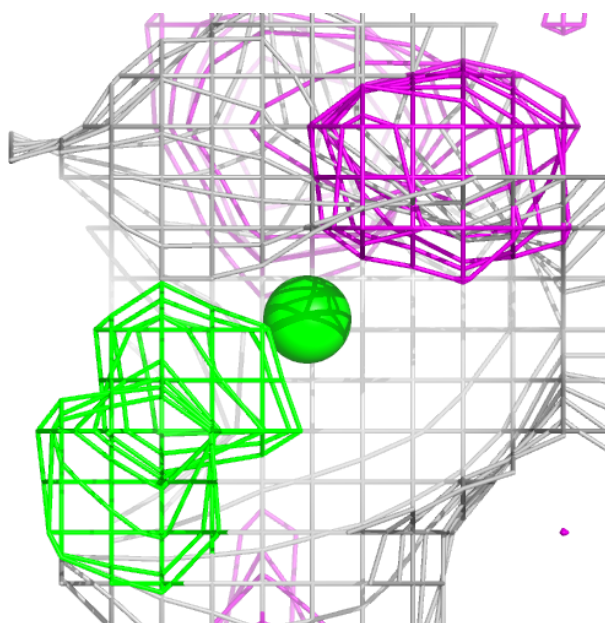
Electron density around BA D 506:

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



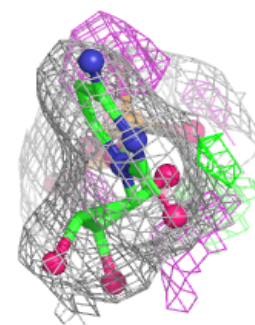
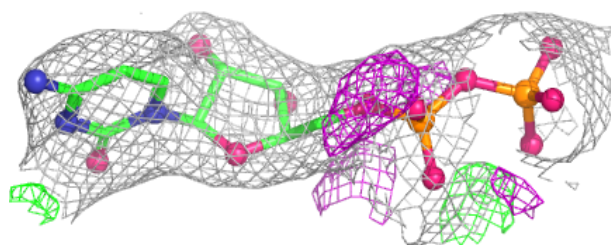
Electron density around BA B 505:

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

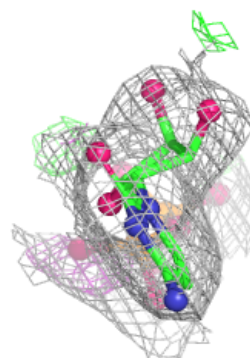
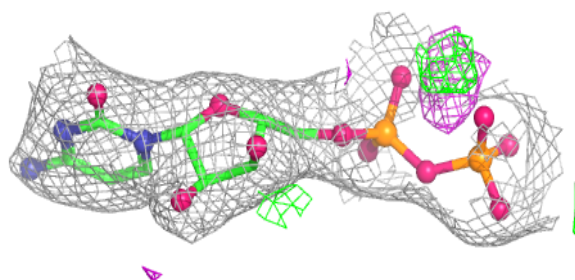
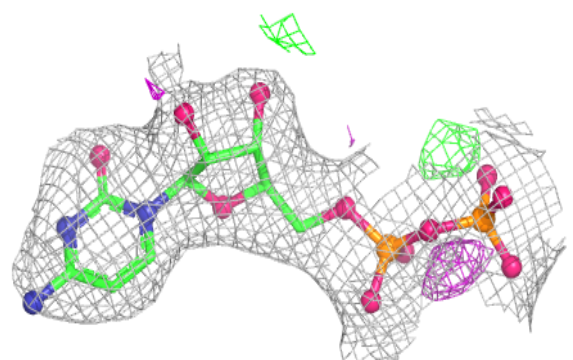


Electron density around CDP A 504:

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

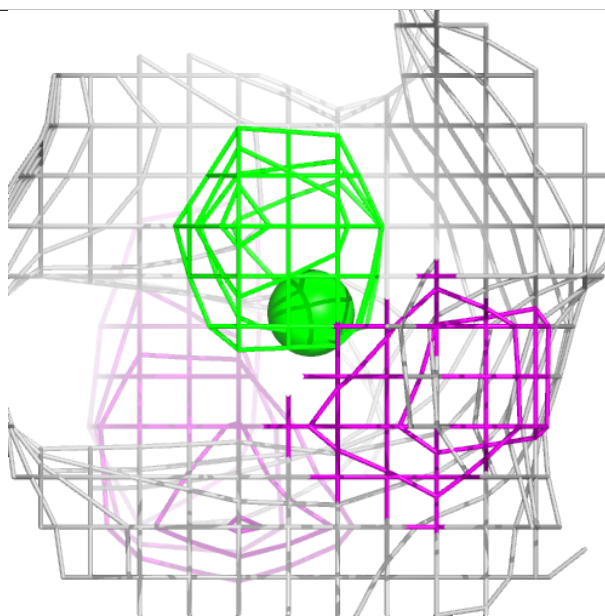
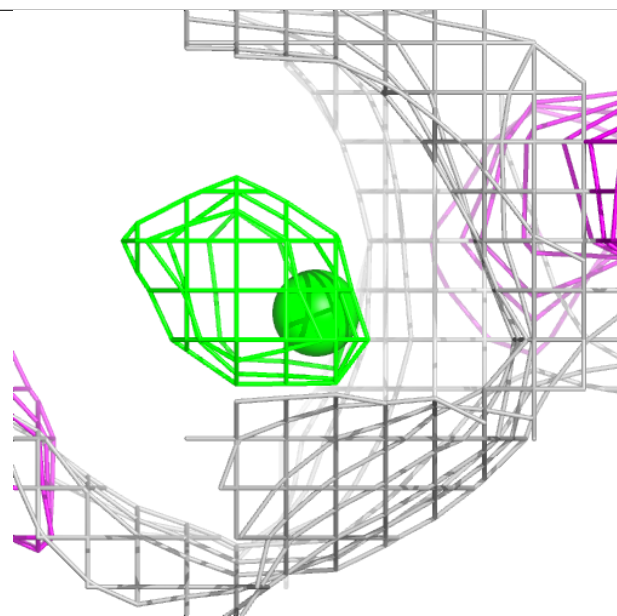
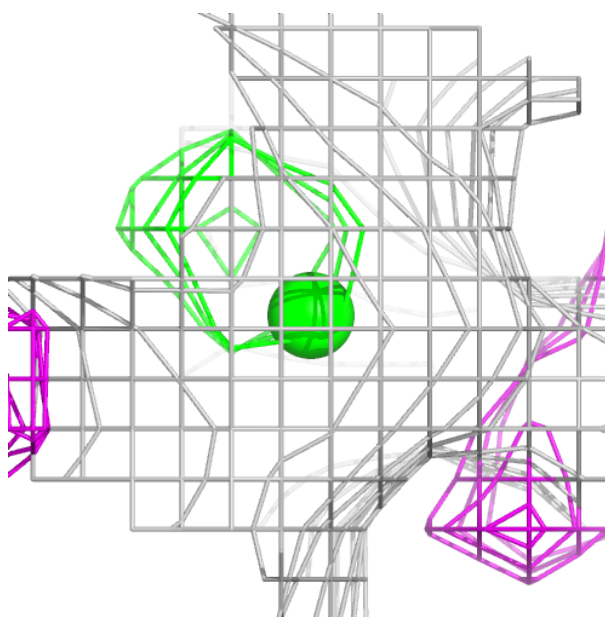
**Electron density around CDP C 504:**

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



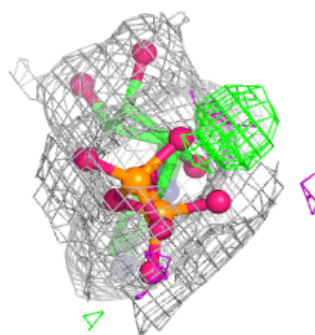
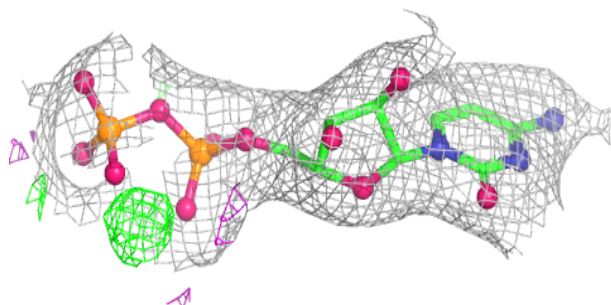
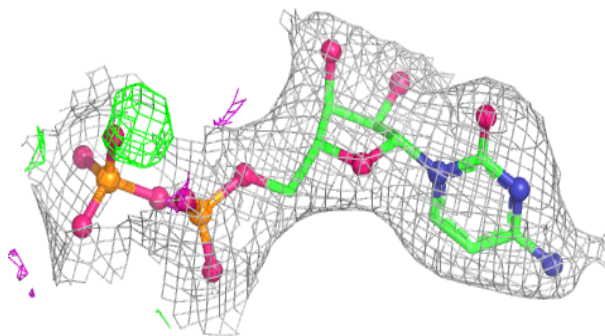
Electron density around BA C 506:

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

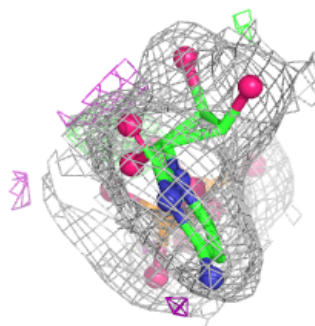
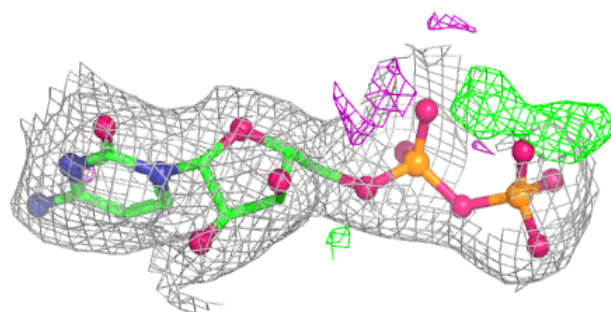
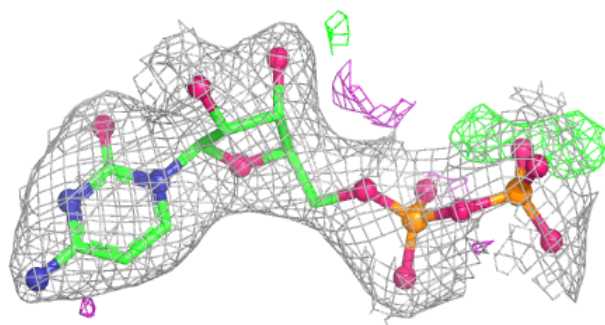


Electron density around CDP D 504:

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

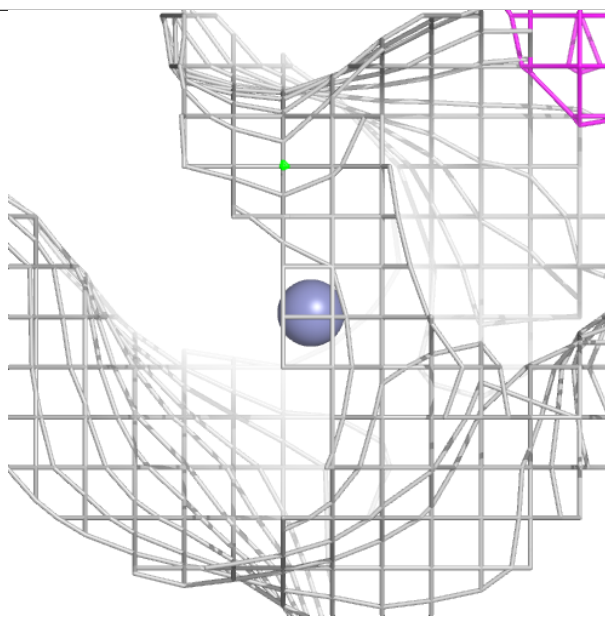
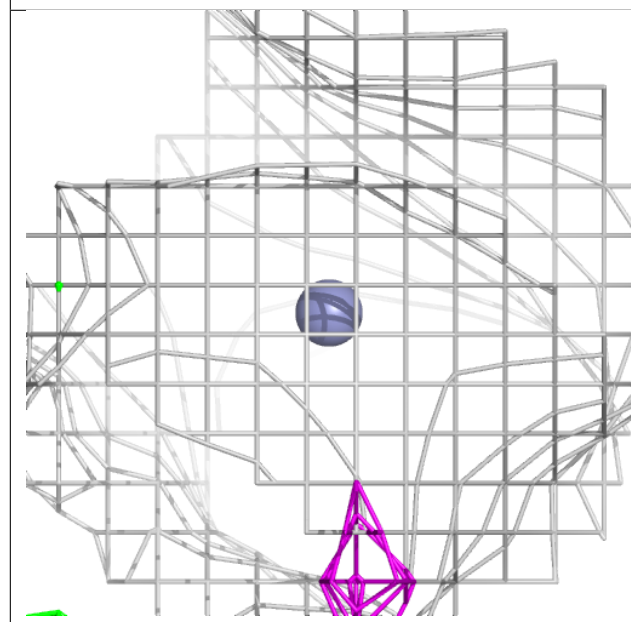
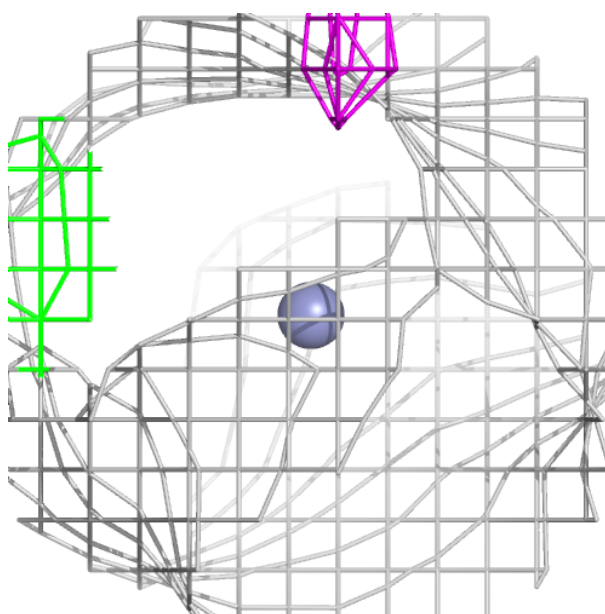
**Electron density around CDP B 503:**

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



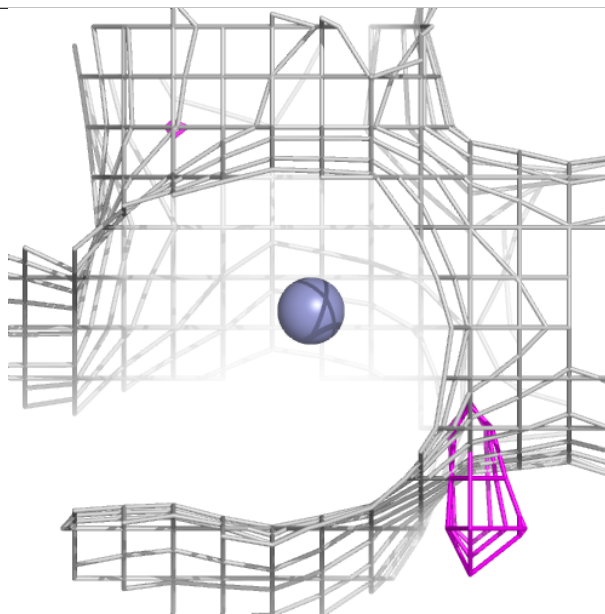
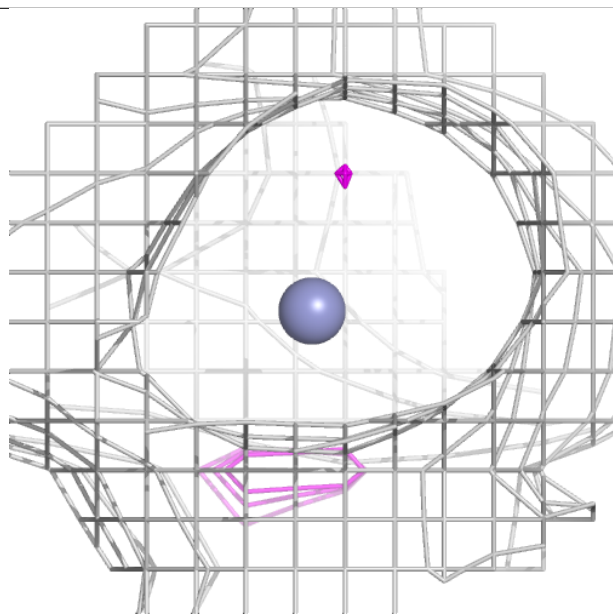
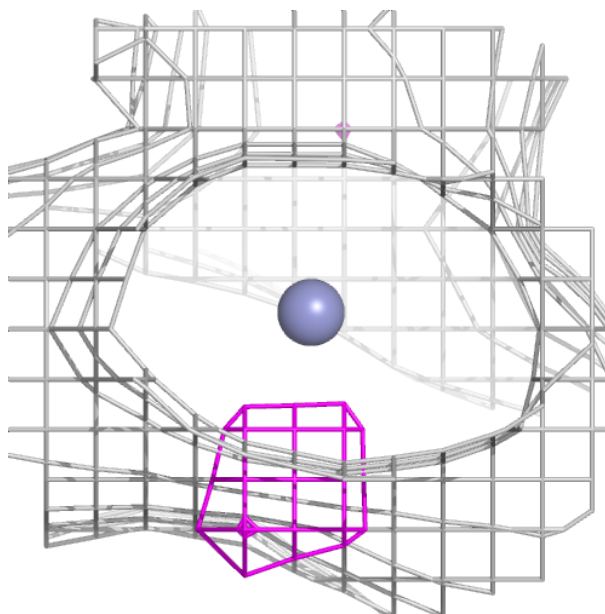
Electron density around ZN B 501:

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



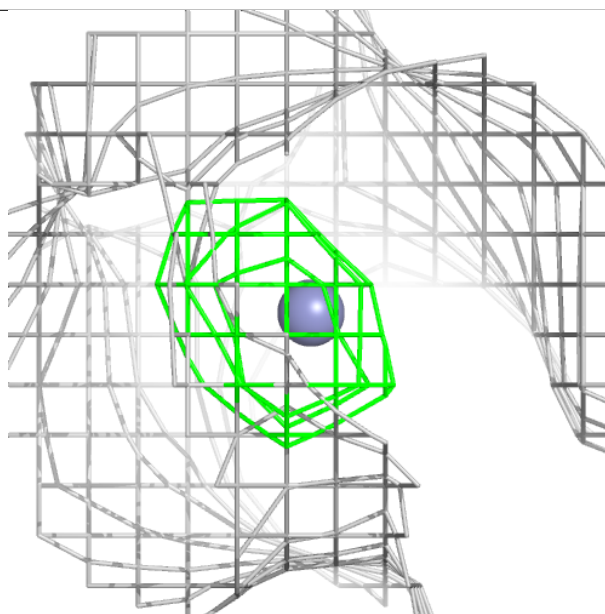
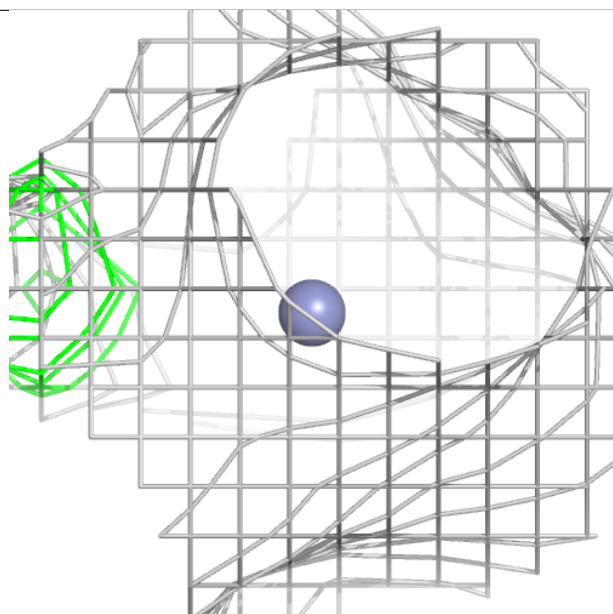
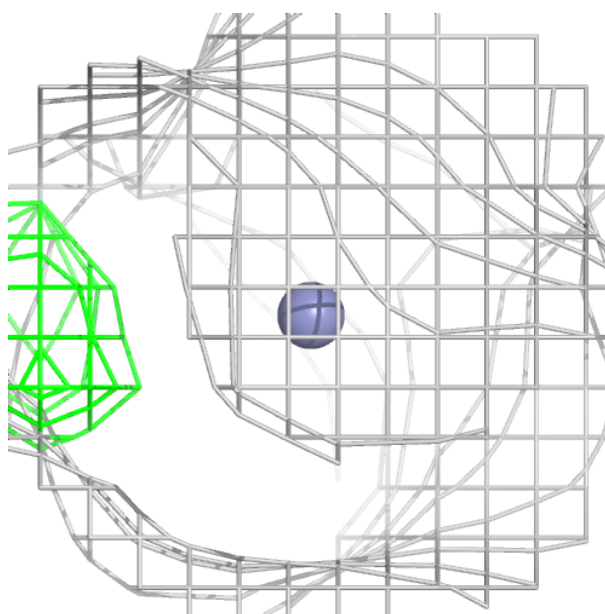
Electron density around ZN C 501:

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



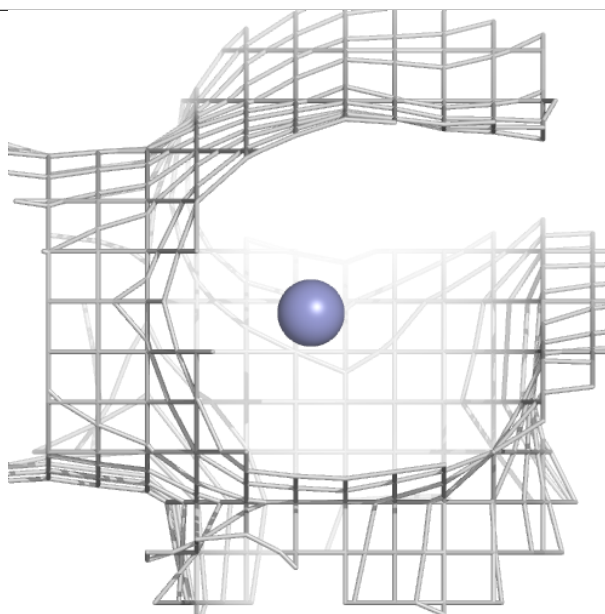
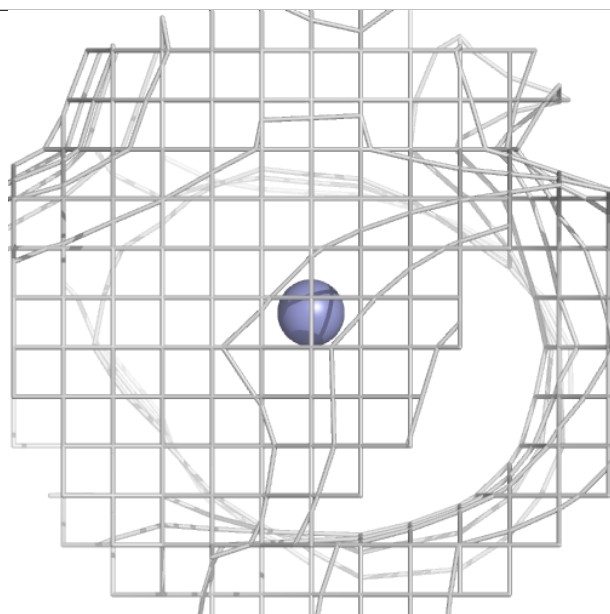
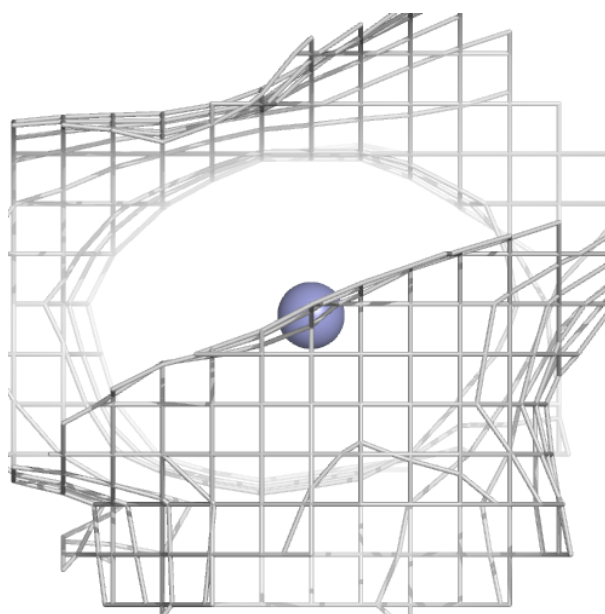
Electron density around ZN D 501:

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



Electron density around ZN A 501:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.