



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 09:29 PM UTC

PDB ID : 6VHQ / pdb_00006vhq
Title : Crystal structure of Bacillus subtilis levansucrase (D86A/E342A) in complex with oligosaccharides
Authors : Diaz-Vilchis, A.; Raga-Carbajal, E.; Rojas-Trejo, S.; Olvera, C.; Rudino-Pinera, E.
Deposited on : 2020-01-10
Resolution : 2.05 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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
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

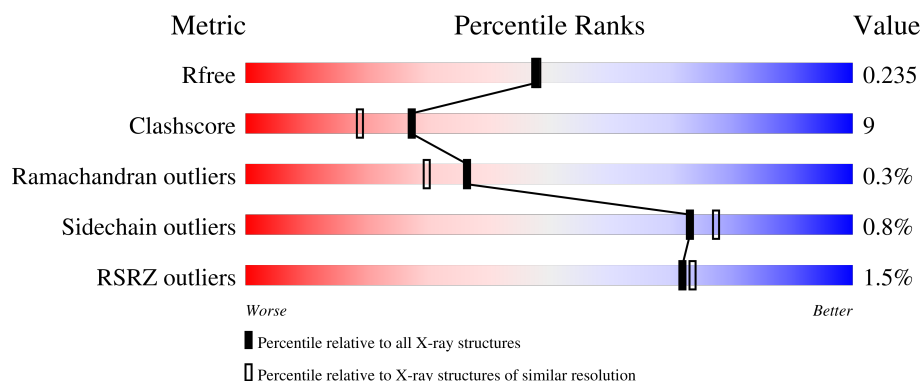
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.05 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	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	466	80% 14% 6%
1	B	466	80% 14% • 6%
2	C	6	100%
2	D	6	83% 17%
2	E	6	83% 17%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
3	F	2	 50%50%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7871 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 Glycoside hydrolase family 68 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	440	Total	C	N	O	S	0	6	0
			3528	2222	588	707	11			
1	B	440	Total	C	N	O	S	0	11	0
			3569	2248	598	712	11			

There are 48 discrepancies between the modelled and reference sequences:

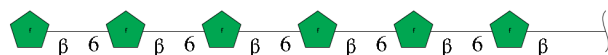
Chain	Residue	Modelled	Actual	Comment	Reference
A	86	ALA	ASP	engineered mutation	UNP A0PFL2
A	342	ALA	GLU	engineered mutation	UNP A0PFL2
A	474	THR	-	expression tag	UNP A0PFL2
A	475	ASP	-	expression tag	UNP A0PFL2
A	476	PRO	-	expression tag	UNP A0PFL2
A	477	ASN	-	expression tag	UNP A0PFL2
A	478	SER	-	expression tag	UNP A0PFL2
A	479	SER	-	expression tag	UNP A0PFL2
A	480	SER	-	expression tag	UNP A0PFL2
A	481	VAL	-	expression tag	UNP A0PFL2
A	482	ASP	-	expression tag	UNP A0PFL2
A	483	LYS	-	expression tag	UNP A0PFL2
A	484	LEU	-	expression tag	UNP A0PFL2
A	485	ALA	-	expression tag	UNP A0PFL2
A	486	ALA	-	expression tag	UNP A0PFL2
A	487	ALA	-	expression tag	UNP A0PFL2
A	488	LEU	-	expression tag	UNP A0PFL2
A	489	GLU	-	expression tag	UNP A0PFL2
A	490	HIS	-	expression tag	UNP A0PFL2
A	491	HIS	-	expression tag	UNP A0PFL2
A	492	HIS	-	expression tag	UNP A0PFL2
A	493	HIS	-	expression tag	UNP A0PFL2
A	494	HIS	-	expression tag	UNP A0PFL2
A	495	HIS	-	expression tag	UNP A0PFL2
B	86	ALA	ASP	engineered mutation	UNP A0PFL2

Continued on next page...

Continued from previous page...

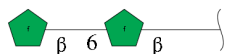
Chain	Residue	Modelled	Actual	Comment	Reference
B	342	ALA	GLU	engineered mutation	UNP A0PFL2
B	474	THR	-	expression tag	UNP A0PFL2
B	475	ASP	-	expression tag	UNP A0PFL2
B	476	PRO	-	expression tag	UNP A0PFL2
B	477	ASN	-	expression tag	UNP A0PFL2
B	478	SER	-	expression tag	UNP A0PFL2
B	479	SER	-	expression tag	UNP A0PFL2
B	480	SER	-	expression tag	UNP A0PFL2
B	481	VAL	-	expression tag	UNP A0PFL2
B	482	ASP	-	expression tag	UNP A0PFL2
B	483	LYS	-	expression tag	UNP A0PFL2
B	484	LEU	-	expression tag	UNP A0PFL2
B	485	ALA	-	expression tag	UNP A0PFL2
B	486	ALA	-	expression tag	UNP A0PFL2
B	487	ALA	-	expression tag	UNP A0PFL2
B	488	LEU	-	expression tag	UNP A0PFL2
B	489	GLU	-	expression tag	UNP A0PFL2
B	490	HIS	-	expression tag	UNP A0PFL2
B	491	HIS	-	expression tag	UNP A0PFL2
B	492	HIS	-	expression tag	UNP A0PFL2
B	493	HIS	-	expression tag	UNP A0PFL2
B	494	HIS	-	expression tag	UNP A0PFL2
B	495	HIS	-	expression tag	UNP A0PFL2

- Molecule 2 is an oligosaccharide called beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	6	Total	C	O	0	0	0
			67	36	31			
2	D	6	Total	C	O	0	0	0
			67	36	31			
2	E	6	Total	C	O	0	0	0
			67	36	31			

- Molecule 3 is an oligosaccharide called beta-D-fructofuranose-(2-6)-beta-D-fructofuranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	F	2	Total	C	O	0	0	0
			22	12	10			

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca	0	0
			1	1		
4	B	1	Total	Ca	0	0
			1	1		

- Molecule 5 is BROMIDE ION (CCD ID: BR) (formula: Br) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	6	Total	Br	0	0
			6	6		
5	B	3	Total	Br	0	0
			3	3		

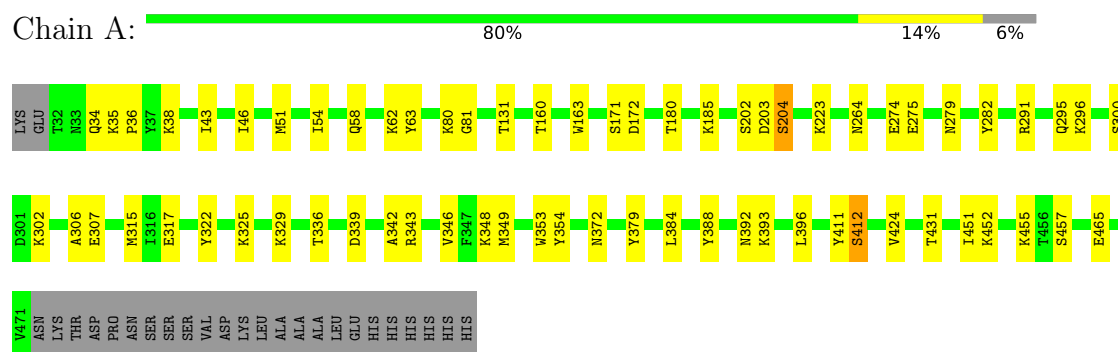
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	290	Total	O	0	0
			290	290		
6	B	250	Total	O	0	0
			250	250		

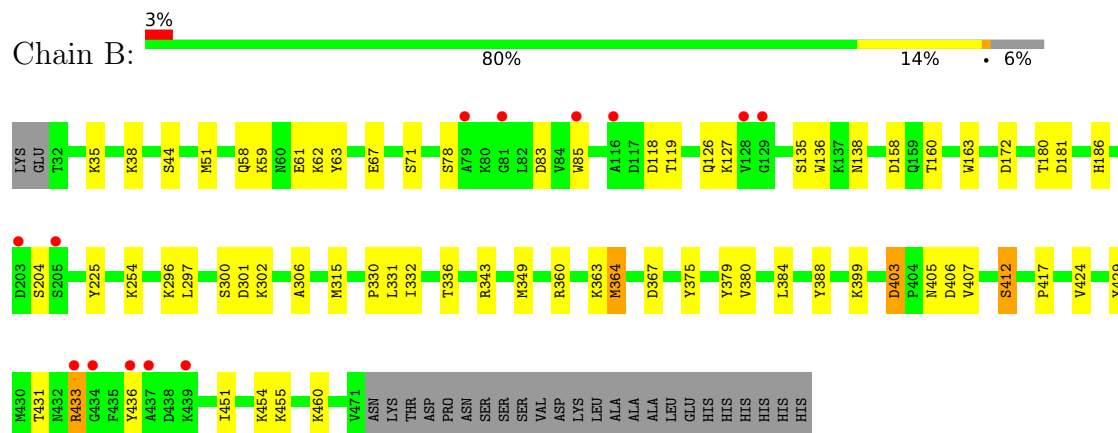
3 Residue-property plots [i](#)

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

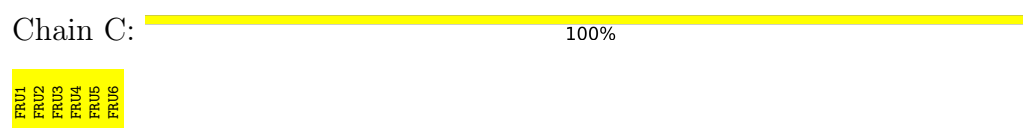
- Molecule 1: Glycoside hydrolase family 68 protein



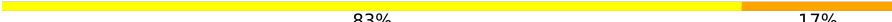
- Molecule 1: Glycoside hydrolase family 68 protein



- Molecule 2: beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose



- Molecule 2: beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose

Chain D:  83% 17%

FRU1
FRU2
FRU3
FRU4
FRU5
FRU6

- Molecule 2: beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose-(2-6)-beta-D-fructofuranose

Chain E:  83% 17%

FRU1
FRU2
FRU3
FRU4
FRU5
FRU6

- Molecule 3: beta-D-fructofuranose-(2-6)-beta-D-fructofuranose

Chain F:  50% 50%

FRU1
FRU2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	69.31Å 78.67Å 78.73Å 90.00° 93.94° 90.00°	Depositor
Resolution (Å)	35.20 – 2.05 35.20 – 2.05	Depositor EDS
% Data completeness (in resolution range)	99.2 (35.20-2.05) 99.2 (35.20-2.05)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.92 (at 2.05Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.167 , 0.235 0.168 , 0.235	Depositor DCC
R_{free} test set	2606 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	25.1	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 63.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7871	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FRU, BR, CA

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	0.69	0/3605	0.83	0/4872
1	B	0.70	0/3649	0.82	2/4932 (0.0%)
All	All	0.70	0/7254	0.82	2/9804 (0.0%)

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	B	0	1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	364	MET	CG-SD-CE	6.42	115.02	100.90
1	B	433	ARG	NE-CZ-NH1	-6.19	115.31	121.50

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	375	TYR	Peptide

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	3528	0	3387	50	0
1	B	3569	0	3417	81	0
2	C	67	0	57	0	0
2	D	67	0	57	1	0
2	E	67	0	58	1	0
3	F	22	0	18	1	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	6	0	0	0	0
5	B	3	0	0	2	0
6	A	290	0	0	9	0
6	B	250	0	0	14	0
All	All	7871	0	6994	130	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 9.

The worst 5 of 130 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:85[A]:TRP:CD1	1:B:433:ARG:HD3	1.62	1.35
5:B:502:BR:BR	6:B:831:HOH:O	2.19	1.12
1:B:85[A]:TRP:NE1	1:B:433:ARG:CD	2.13	1.11
1:B:85[A]:TRP:CE2	1:B:433:ARG:HD2	1.92	1.05
1:B:85[A]:TRP:CE2	1:B:433:ARG:CD	2.44	1.01

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	444/466 (95%)	424 (96%)	19 (4%)	1 (0%)	43	37
1	B	448/466 (96%)	433 (97%)	13 (3%)	2 (0%)	30	22
All	All	892/932 (96%)	857 (96%)	32 (4%)	3 (0%)	36	30

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	204	SER
1	A	431	THR
1	B	431	THR

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	389/406 (96%)	386 (99%)	3 (1%)	73	77
1	B	393/406 (97%)	390 (99%)	3 (1%)	73	77
All	All	782/812 (96%)	776 (99%)	6 (1%)	73	77

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	181	ASP
1	B	403	ASP
1	B	412	SER
1	A	204	SER
1	A	171	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 13 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	126	GLN
1	B	152	ASN
1	B	450	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	382	ASN
1	B	405	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 ⓘ

20 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	FRU	C	1	2	11,12,12	4.36	7 (63%)	10,18,18	1.82	2 (20%)
2	FRU	C	2	2	11,11,12	4.43	8 (72%)	15,15,18	1.19	2 (13%)
2	FRU	C	3	2	11,11,12	4.30	8 (72%)	15,15,18	2.57	8 (53%)
2	FRU	C	4	2	11,11,12	4.47	8 (72%)	15,15,18	1.26	2 (13%)
2	FRU	C	5	2	11,11,12	4.31	8 (72%)	15,15,18	1.87	4 (26%)
2	FRU	C	6	2	11,11,12	4.14	8 (72%)	15,15,18	1.35	2 (13%)
2	FRU	D	1	2	11,12,12	4.46	7 (63%)	10,18,18	4.96	1 (10%)
2	FRU	D	2	2	11,11,12	4.52	8 (72%)	15,15,18	1.33	1 (6%)
2	FRU	D	3	2	11,11,12	4.52	8 (72%)	15,15,18	1.18	1 (6%)
2	FRU	D	4	2	11,11,12	4.10	8 (72%)	15,15,18	2.23	5 (33%)
2	FRU	D	5	2	11,11,12	4.34	8 (72%)	15,15,18	1.72	2 (13%)
2	FRU	D	6	2	11,11,12	4.51	8 (72%)	15,15,18	0.96	1 (6%)
2	FRU	E	1	2	11,12,12	4.28	7 (63%)	10,18,18	3.64	2 (20%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	FRU	E	2	2	11,11,12	4.51	8 (72%)	15,15,18	1.50	3 (20%)
2	FRU	E	3	2	11,11,12	4.53	8 (72%)	15,15,18	1.76	4 (26%)
2	FRU	E	4	2	11,11,12	4.52	8 (72%)	15,15,18	2.74	3 (20%)
2	FRU	E	5	2	11,11,12	4.76	7 (63%)	15,15,18	1.51	3 (20%)
2	FRU	E	6	2	11,11,12	4.27	8 (72%)	15,15,18	1.85	6 (40%)
3	FRU	F	1	3	11,12,12	4.17	7 (63%)	10,18,18	1.07	2 (20%)
3	FRU	F	2	3	10,10,12	5.19	8 (80%)	13,14,18	2.00	4 (30%)

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	FRU	C	1	2	-	3/5/24/24	0/1/1/1
2	FRU	C	2	2	-	1/4/20/24	0/1/1/1
2	FRU	C	3	2	-	1/4/20/24	0/1/1/1
2	FRU	C	4	2	-	1/4/20/24	0/1/1/1
2	FRU	C	5	2	-	2/4/20/24	0/1/1/1
2	FRU	C	6	2	-	0/4/20/24	0/1/1/1
2	FRU	D	1	2	-	3/5/24/24	0/1/1/1
2	FRU	D	2	2	-	2/4/20/24	0/1/1/1
2	FRU	D	3	2	-	2/4/20/24	0/1/1/1
2	FRU	D	4	2	-	0/4/20/24	0/1/1/1
2	FRU	D	5	2	-	2/4/20/24	0/1/1/1
2	FRU	D	6	2	-	0/4/20/24	0/1/1/1
2	FRU	E	1	2	-	3/5/24/24	0/1/1/1
2	FRU	E	2	2	-	2/4/20/24	0/1/1/1
2	FRU	E	3	2	-	4/4/20/24	0/1/1/1
2	FRU	E	4	2	-	4/4/20/24	0/1/1/1
2	FRU	E	5	2	-	3/4/20/24	0/1/1/1
2	FRU	E	6	2	-	0/4/20/24	0/1/1/1
3	FRU	F	1	3	-	0/5/24/24	0/1/1/1
3	FRU	F	2	3	-	2/2/18/24	0/1/1/1

The worst 5 of 155 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	5	FRU	C4-C3	-11.22	1.22	1.53
3	F	2	FRU	C4-C3	-10.61	1.24	1.53
2	E	3	FRU	C4-C3	-10.47	1.25	1.53
2	E	2	FRU	C4-C3	-10.36	1.25	1.53
2	C	5	FRU	C4-C3	-10.36	1.25	1.53

The worst 5 of 58 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	FRU	O6-C6-C5	15.17	162.99	111.33
2	E	1	FRU	O6-C6-C5	11.13	149.22	111.33
2	E	4	FRU	O6-C6-C5	9.33	143.10	111.33
2	C	3	FRU	O5-C2-C1	6.51	122.98	109.22
3	F	2	FRU	C6-C5-C4	-5.60	109.81	115.70

There are no chirality outliers.

5 of 35 torsion outliers are listed below:

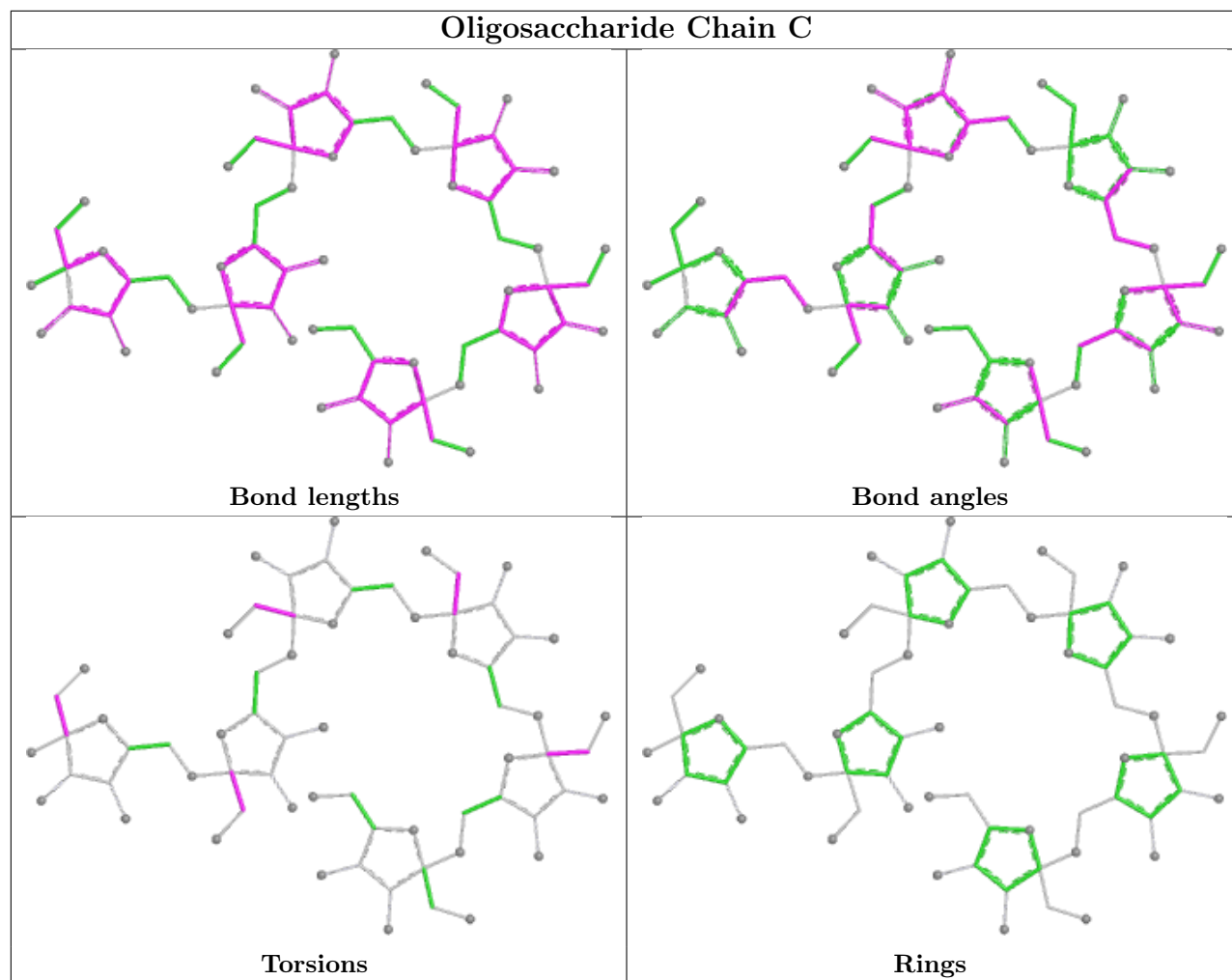
Mol	Chain	Res	Type	Atoms
2	C	1	FRU	O1-C1-C2-C3
2	C	1	FRU	O1-C1-C2-O2
2	D	1	FRU	O1-C1-C2-C3
2	E	1	FRU	O1-C1-C2-C3
2	E	1	FRU	O1-C1-C2-O2

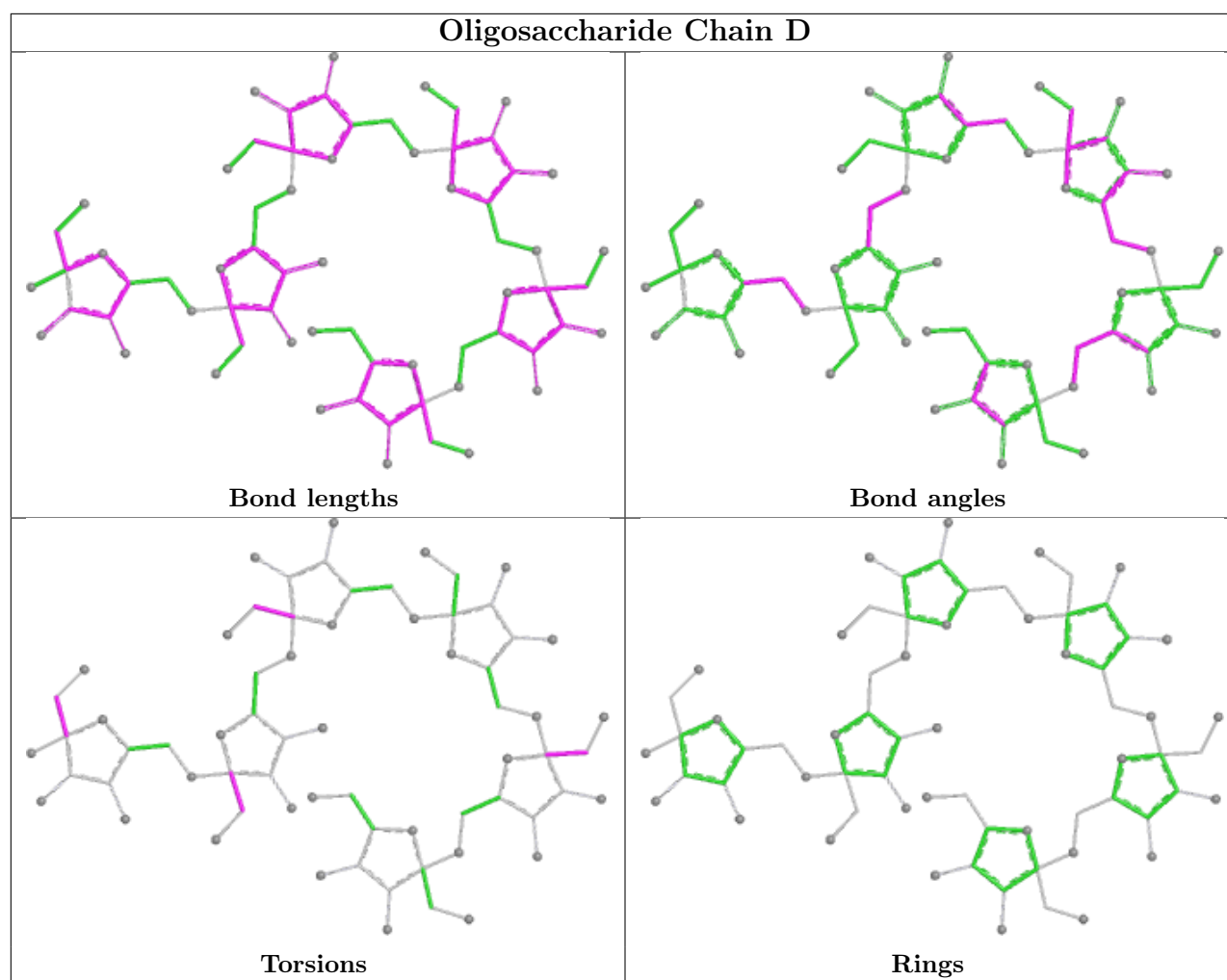
There are no ring outliers.

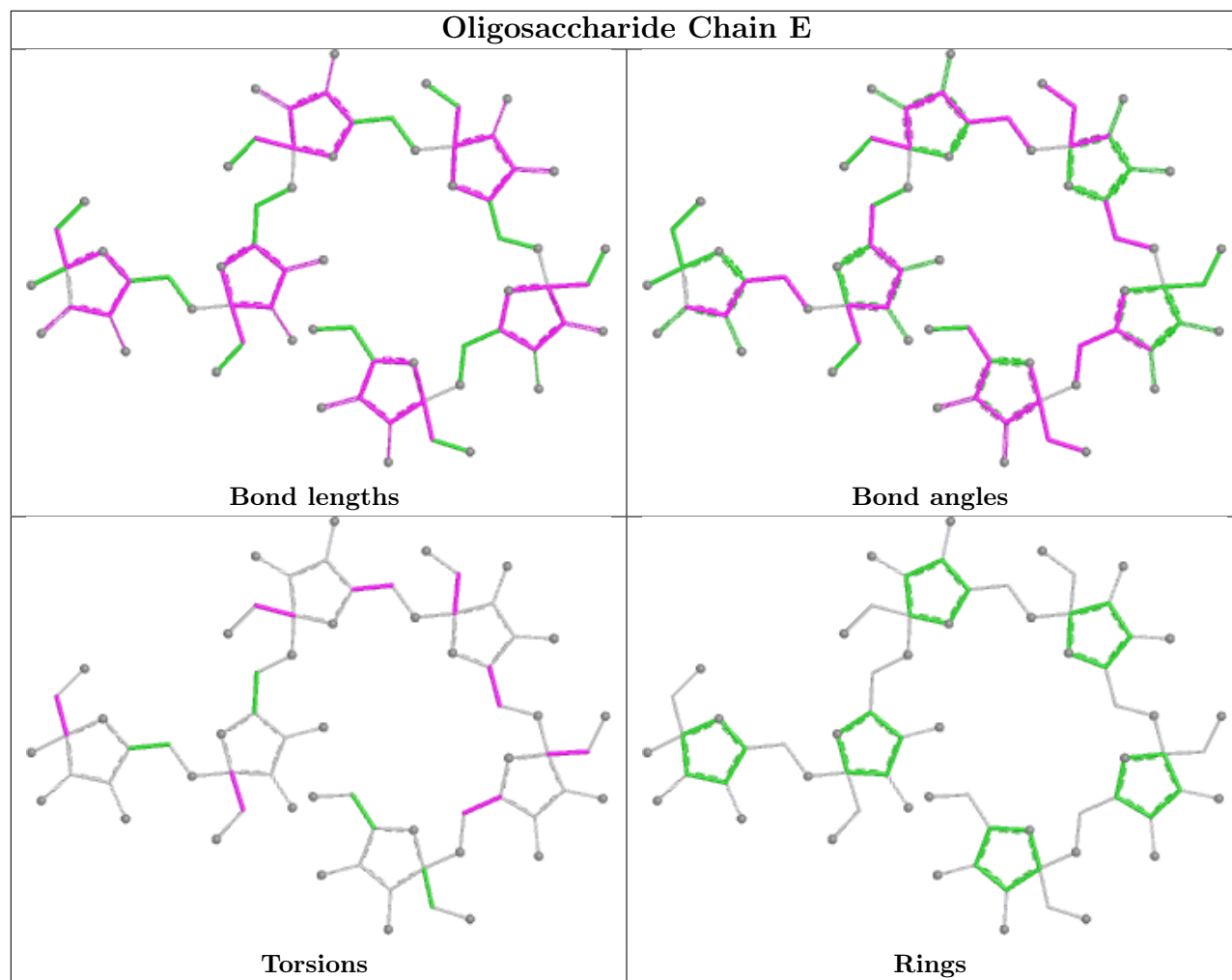
3 monomers are involved in 3 short contacts:

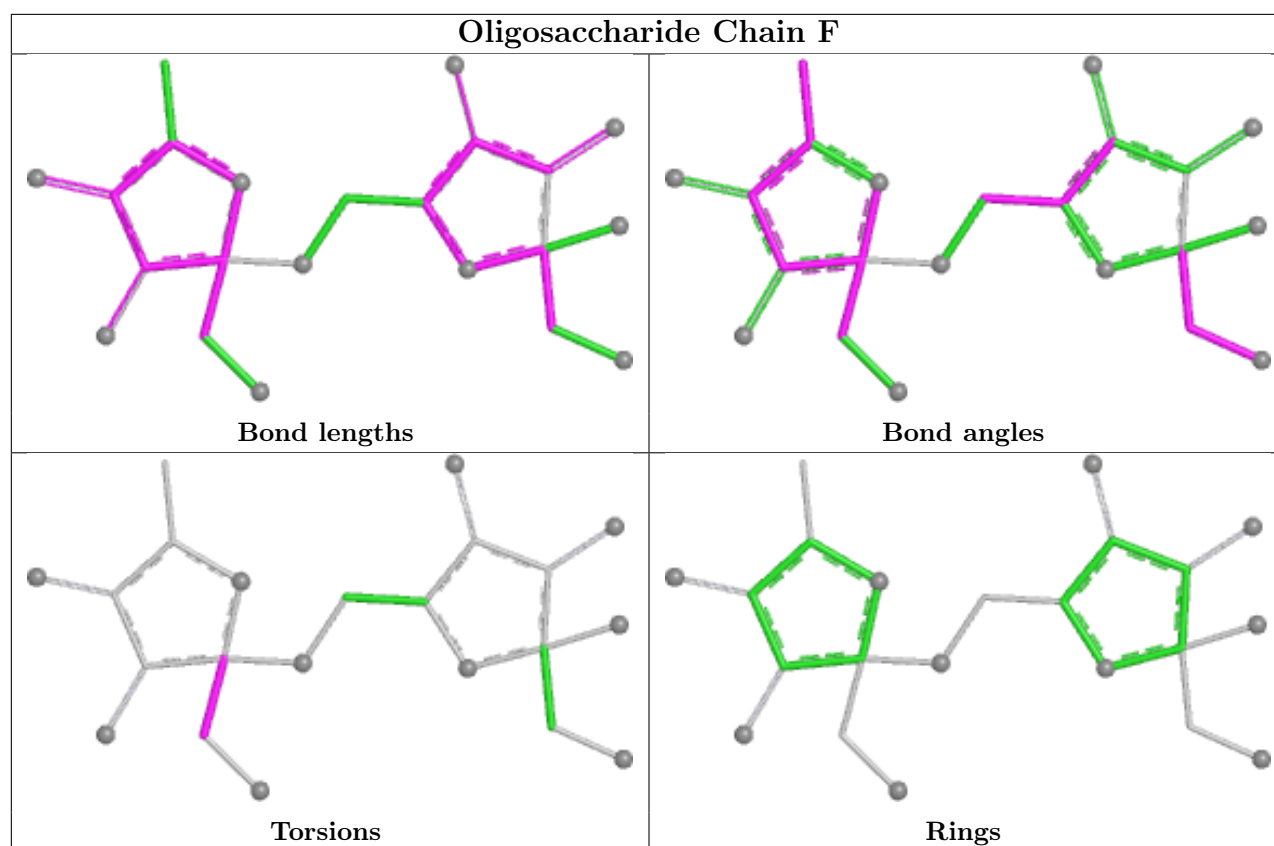
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	5	FRU	1	0
2	D	4	FRU	1	0
3	F	1	FRU	1	0

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









5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 11 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 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	440/466 (94%)	-0.37	0 100 100	7, 24, 45, 78	6 (1%)
1	B	440/466 (94%)	-0.24	13 (2%) 52 53	8, 25, 55, 86	10 (2%)
All	All	880/932 (94%)	-0.31	13 (1%) 72 73	7, 24, 49, 86	16 (1%)

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	436	TYR	3.1
1	B	129	GLY	3.0
1	B	205	SER	2.9
1	B	85[A]	TRP	2.8
1	B	433	ARG	2.8

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.

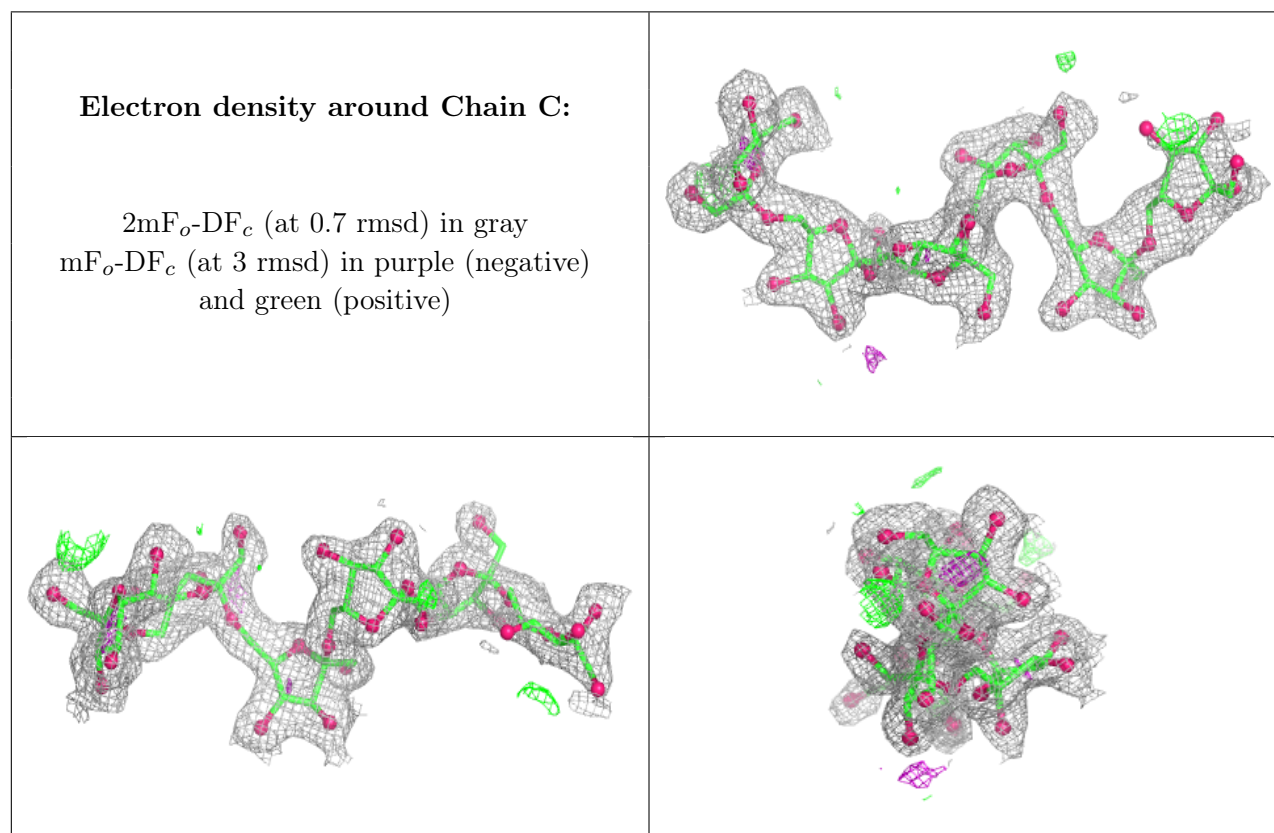
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	FRU	F	1	12/12	0.50	0.21	40,51,59,68	12
2	FRU	E	2	11/12	0.56	0.16	65,70,81,87	1
3	FRU	F	2	10/12	0.56	0.20	29,51,55,58	10
2	FRU	E	1	12/12	0.62	0.13	68,78,93,93	0
2	FRU	E	3	11/12	0.69	0.15	52,55,65,70	11

Continued on next page...

Continued from previous page...

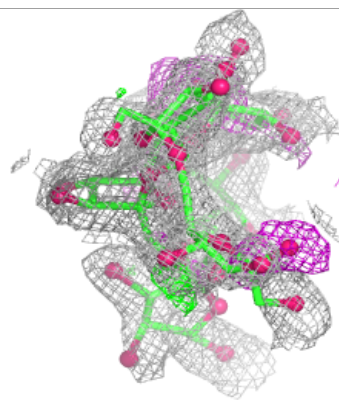
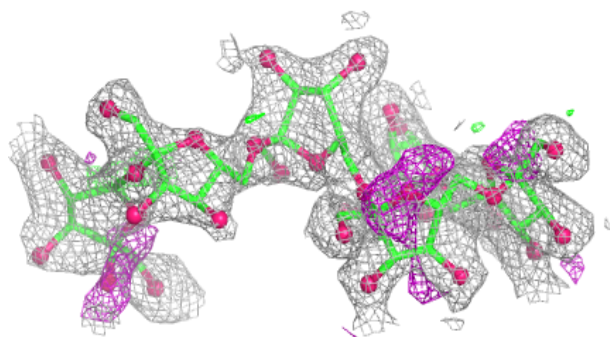
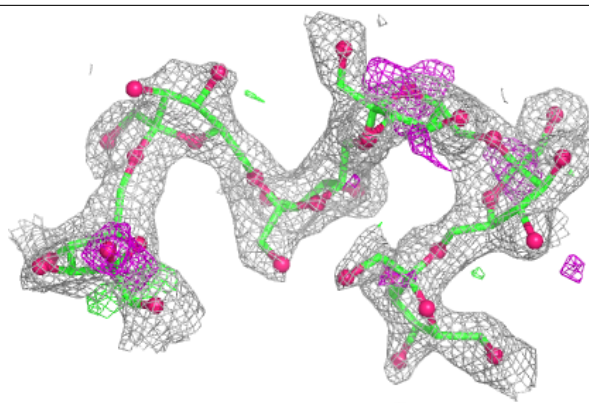
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FRU	D	6	11/12	0.72	0.12	35,50,57,61	0
2	FRU	C	1	12/12	0.74	0.12	54,62,66,68	0
2	FRU	E	4	11/12	0.77	0.14	28,44,48,48	11
2	FRU	D	2	11/12	0.78	0.11	44,46,53,58	0
2	FRU	D	1	12/12	0.78	0.13	38,48,59,61	0
2	FRU	E	5	11/12	0.79	0.13	30,39,53,58	11
2	FRU	E	6	11/12	0.81	0.15	16,27,40,41	11
2	FRU	D	5	11/12	0.83	0.11	27,33,46,49	0
2	FRU	D	3	11/12	0.84	0.10	35,42,52,59	0
2	FRU	D	4	11/12	0.84	0.12	25,31,36,39	0
2	FRU	C	3	11/12	0.85	0.10	28,37,44,44	0
2	FRU	C	2	11/12	0.87	0.09	38,50,65,65	0
2	FRU	C	5	11/12	0.91	0.08	25,30,42,45	0
2	FRU	C	6	11/12	0.91	0.09	15,23,32,37	0
2	FRU	C	4	11/12	0.93	0.07	23,28,33,34	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

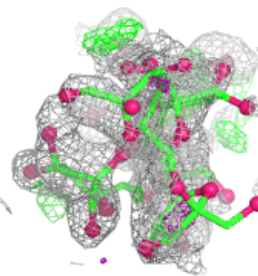
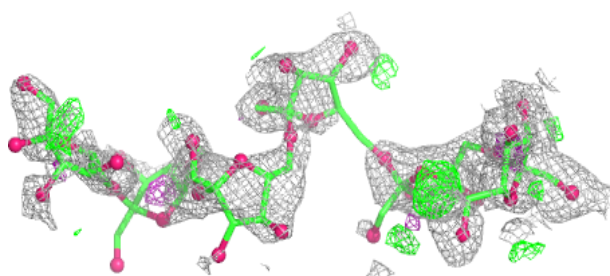
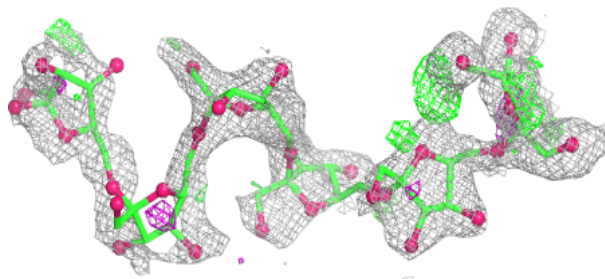


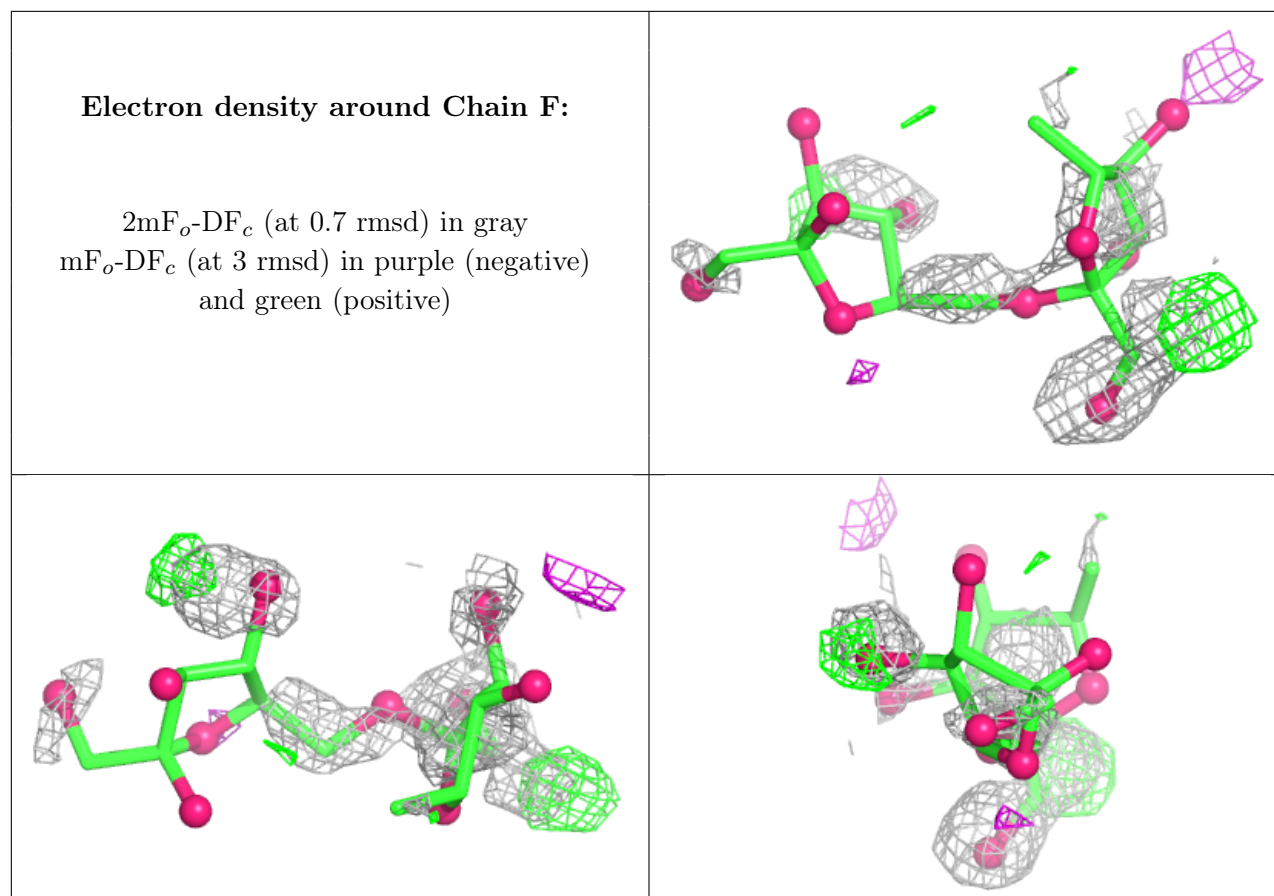
Electron density around Chain D:

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

**Electron density around Chain E:**

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





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.

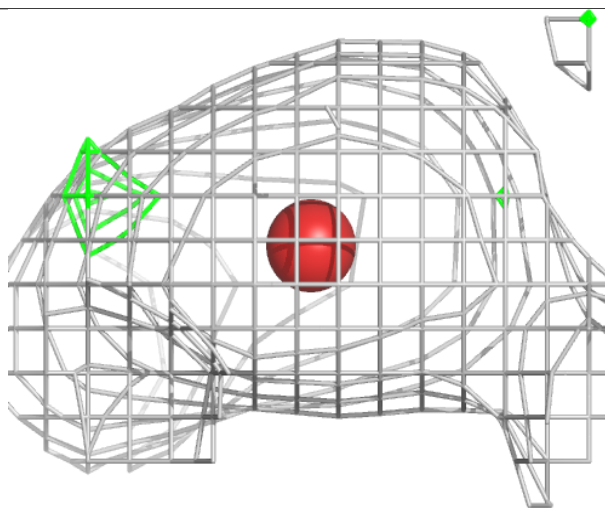
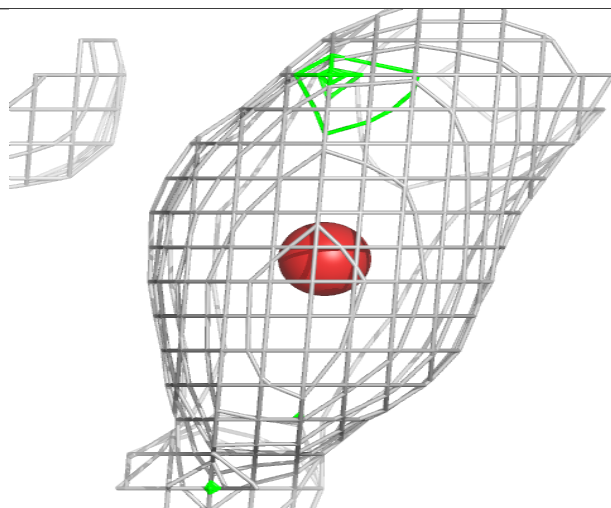
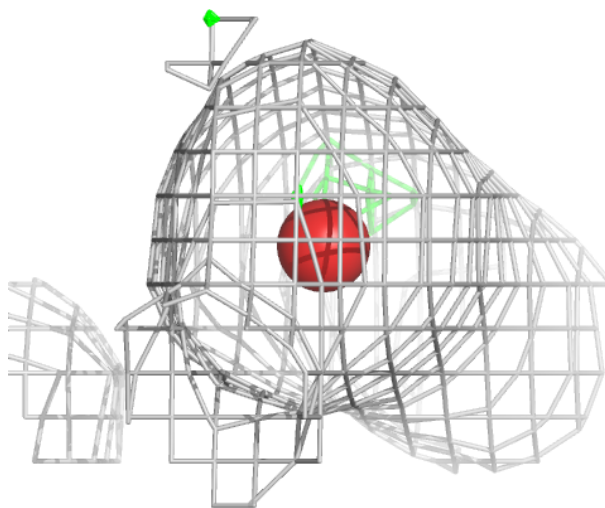
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	BR	B	502	1/1	0.93	0.07	45,45,45,45	1
5	BR	A	507	1/1	0.95	0.23	79,79,79,79	0
5	BR	B	503	1/1	0.96	0.05	38,38,38,38	1
5	BR	A	506	1/1	0.98	0.04	38,38,38,38	1
5	BR	A	505	1/1	0.99	0.03	25,25,25,25	0
4	CA	A	501	1/1	0.99	0.03	14,14,14,14	0
4	CA	B	501	1/1	0.99	0.01	15,15,15,15	0
5	BR	A	502	1/1	0.99	0.02	25,25,25,25	1
5	BR	A	503	1/1	0.99	0.03	24,24,24,24	1
5	BR	B	504	1/1	0.99	0.03	34,34,34,34	1
5	BR	A	504	1/1	1.00	0.01	22,22,22,22	1

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.

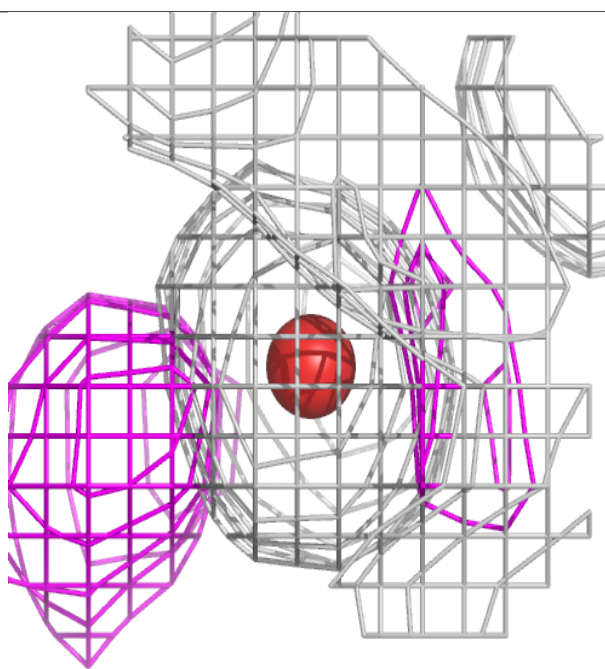
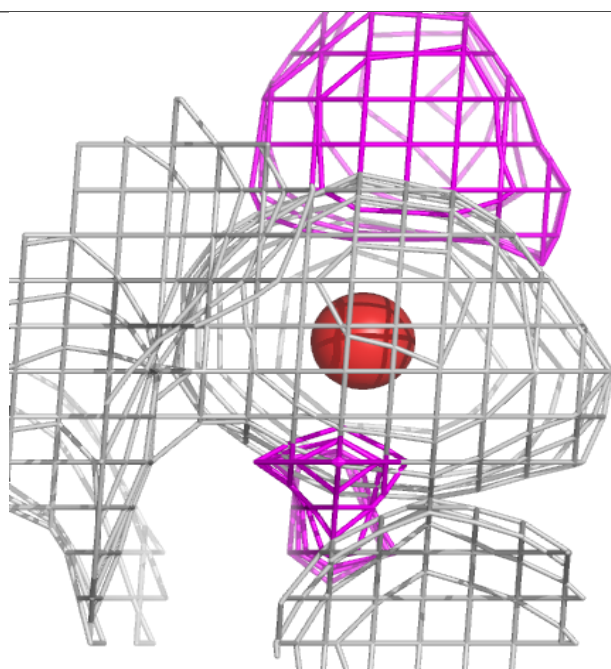
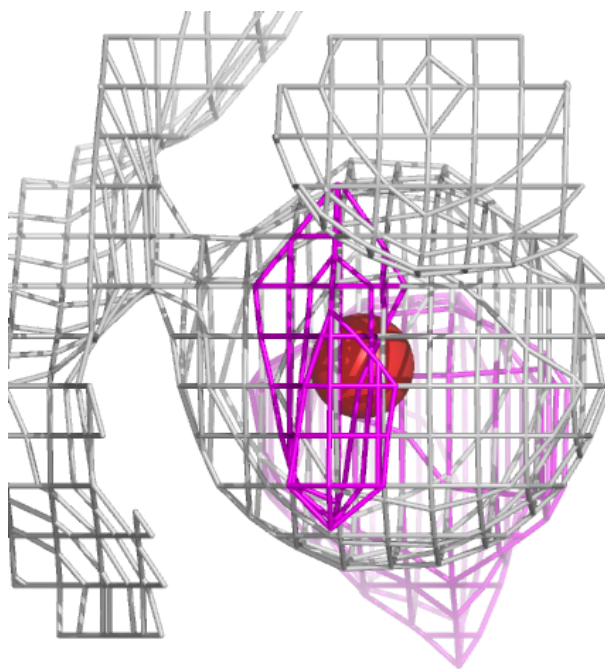
Electron density around BR B 502:

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



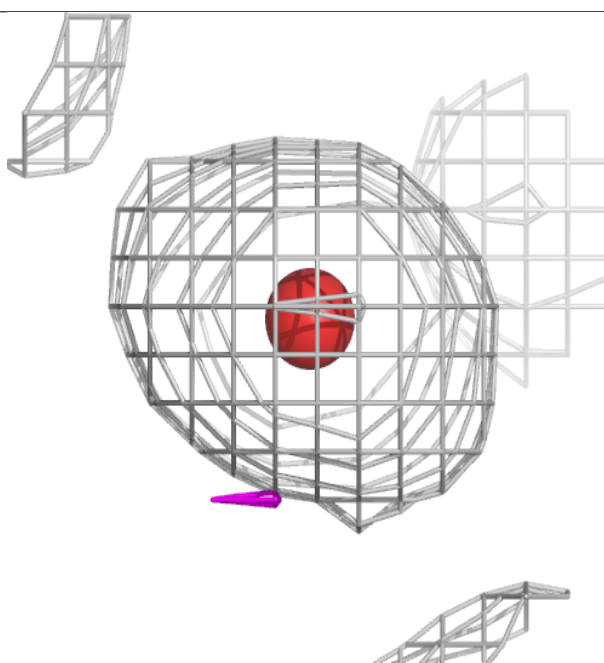
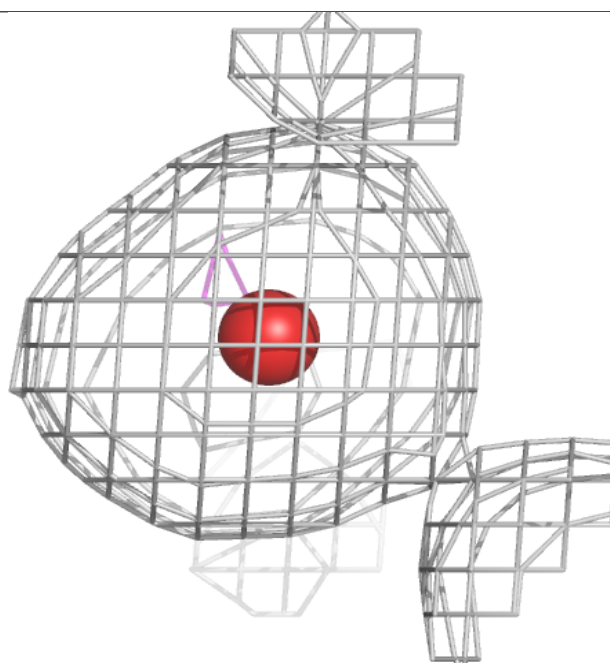
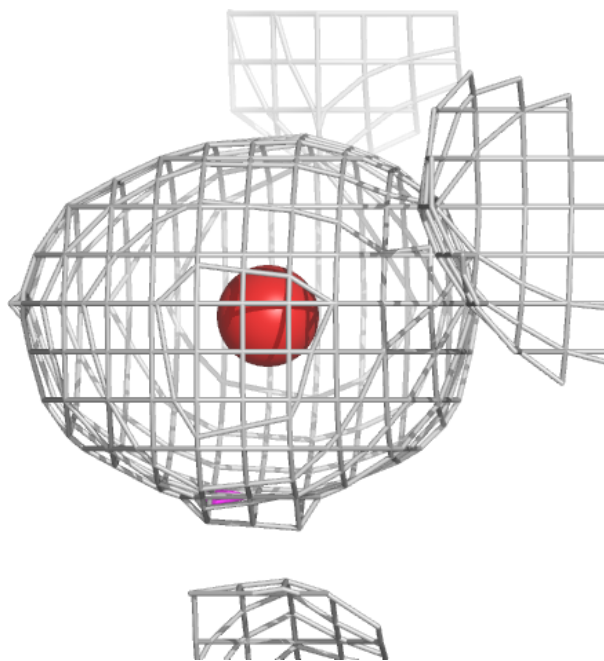
Electron density around BR A 507:

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



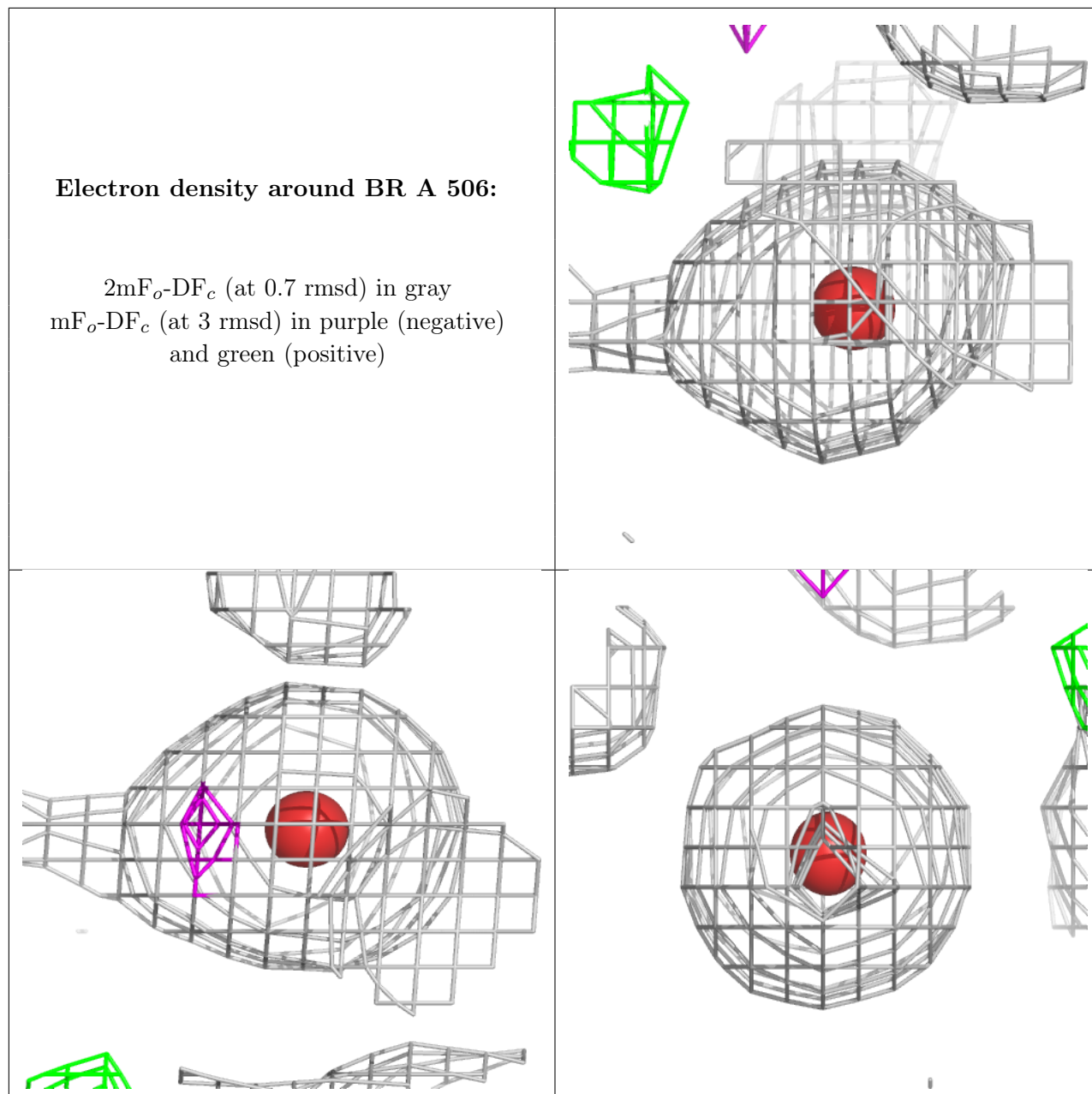
Electron density around BR B 503:

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



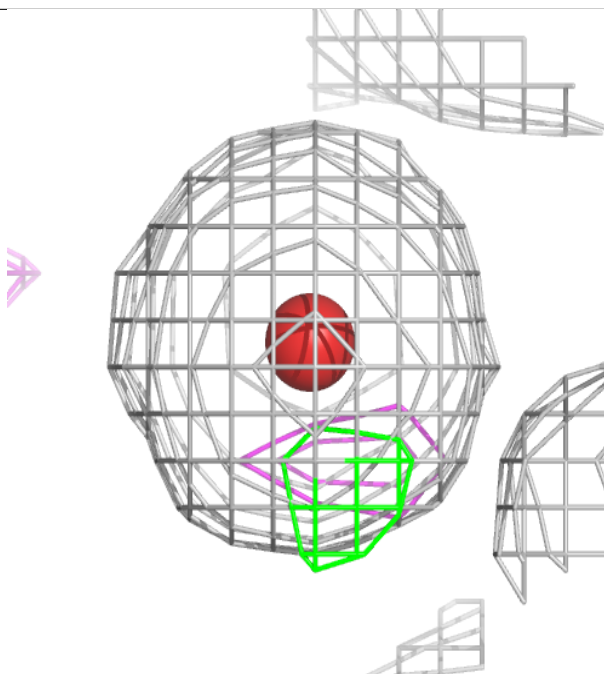
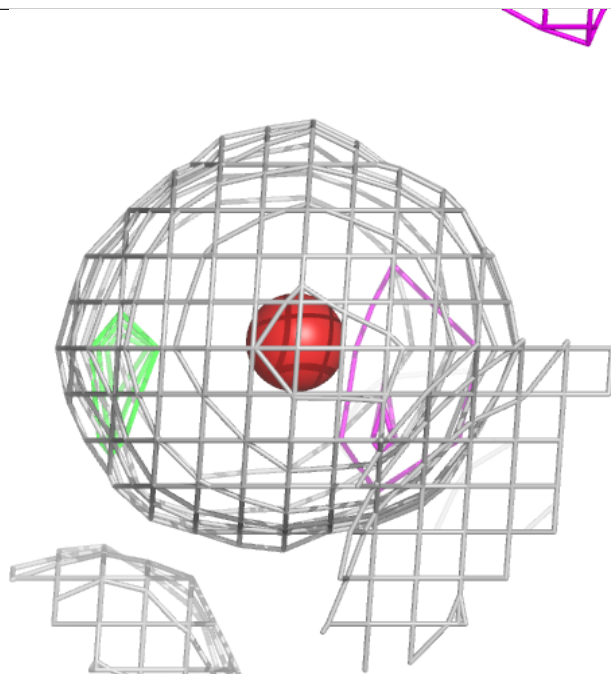
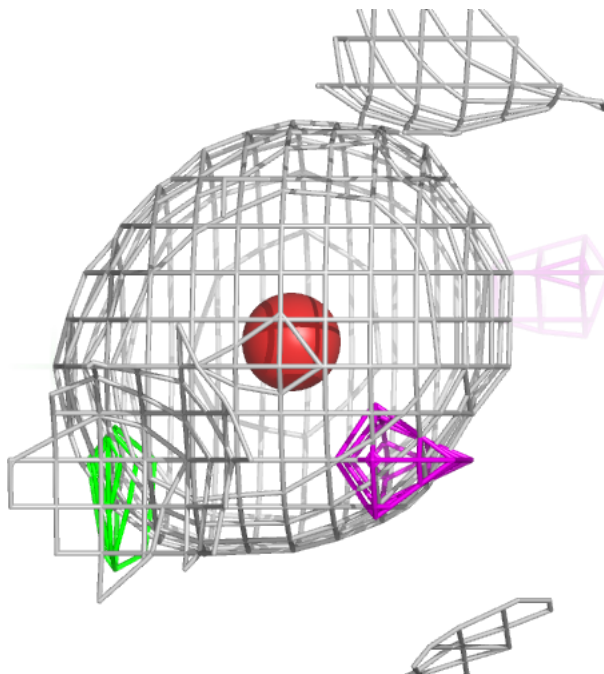
Electron density around BR A 506:

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



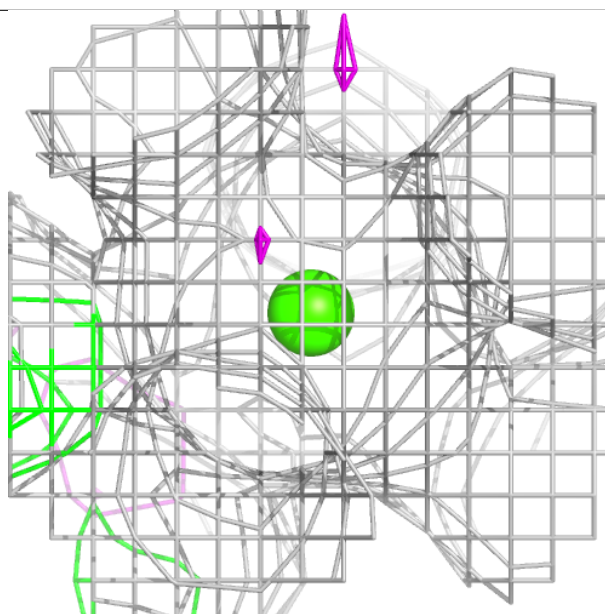
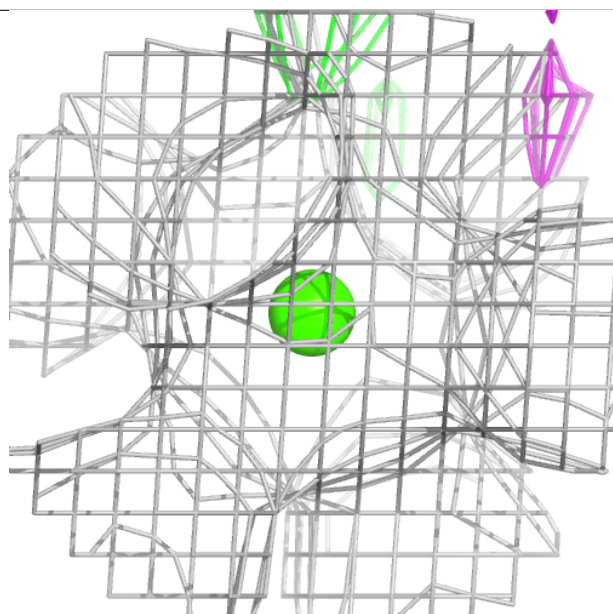
Electron density around BR A 505:

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



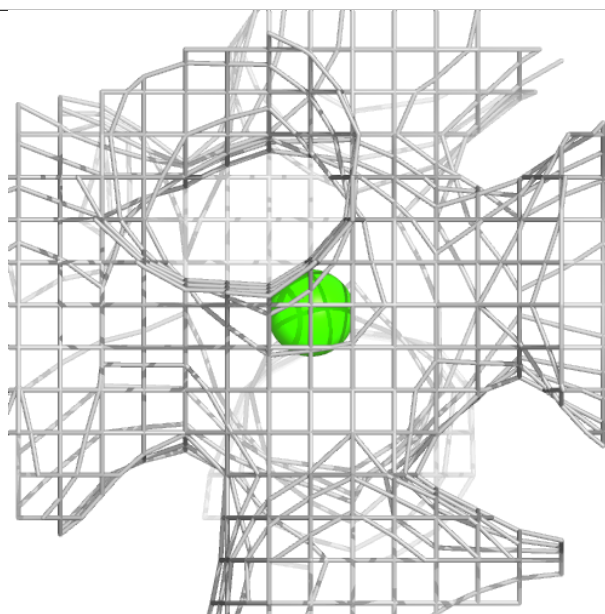
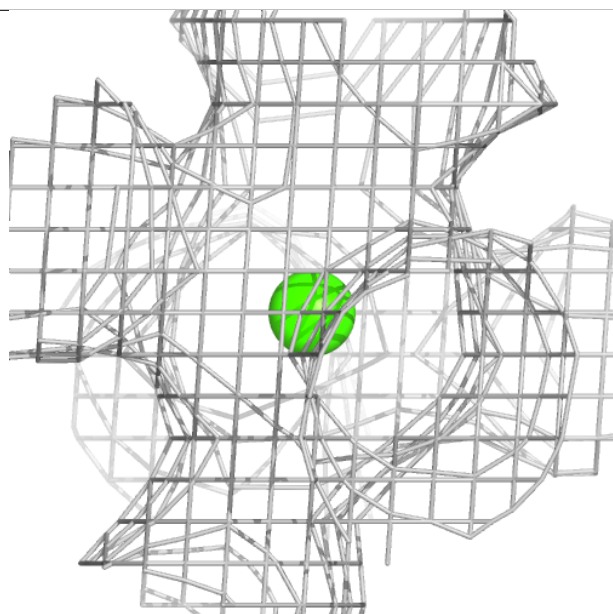
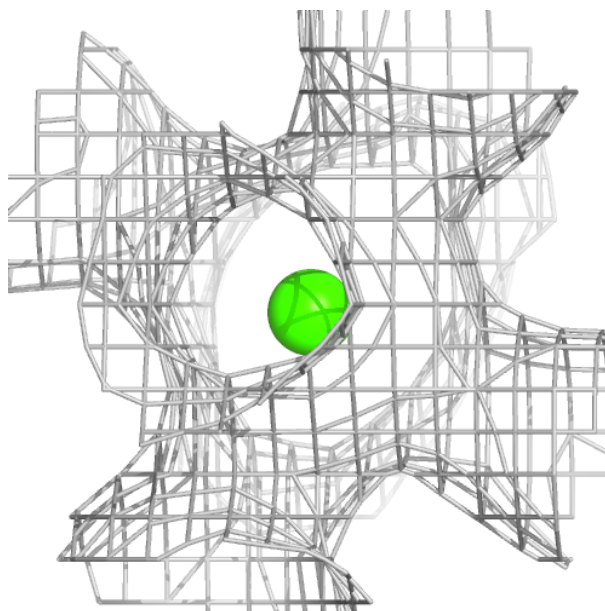
Electron density around CA A 501:

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



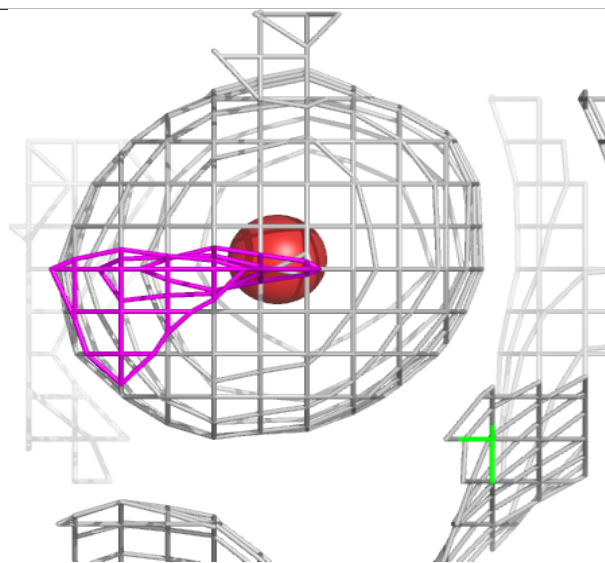
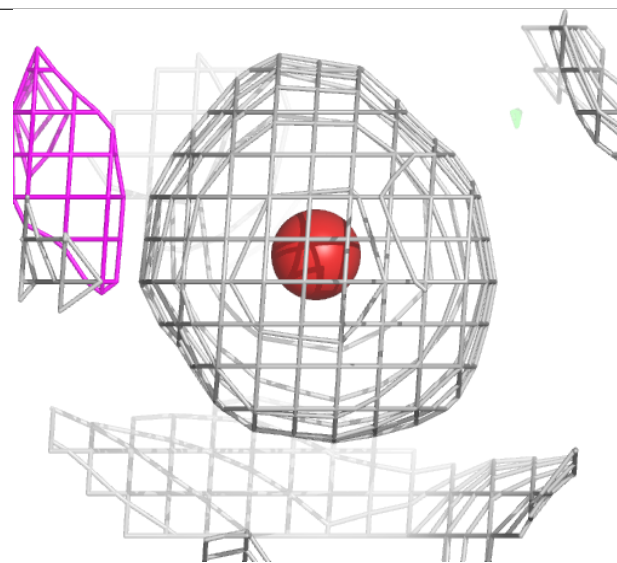
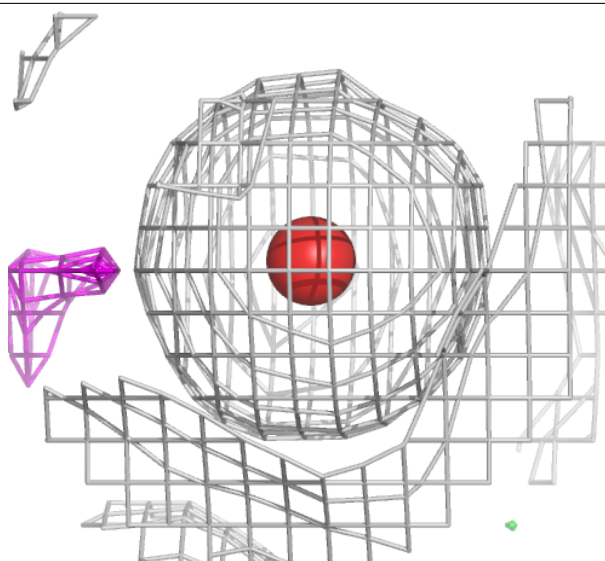
Electron density around CA B 501:

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



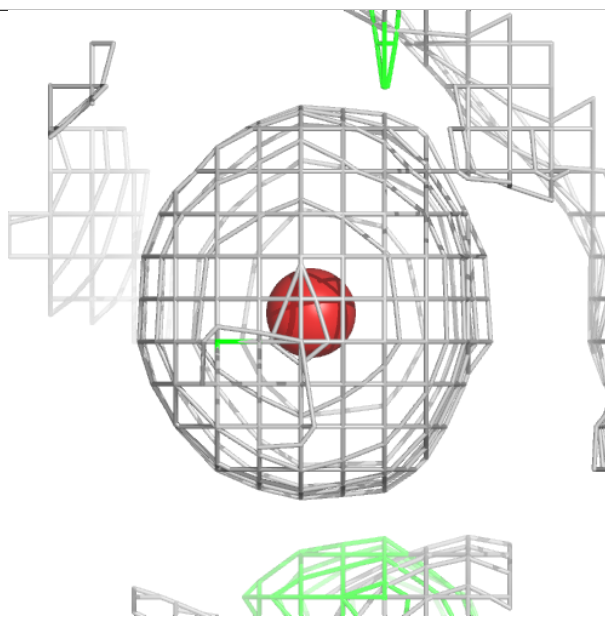
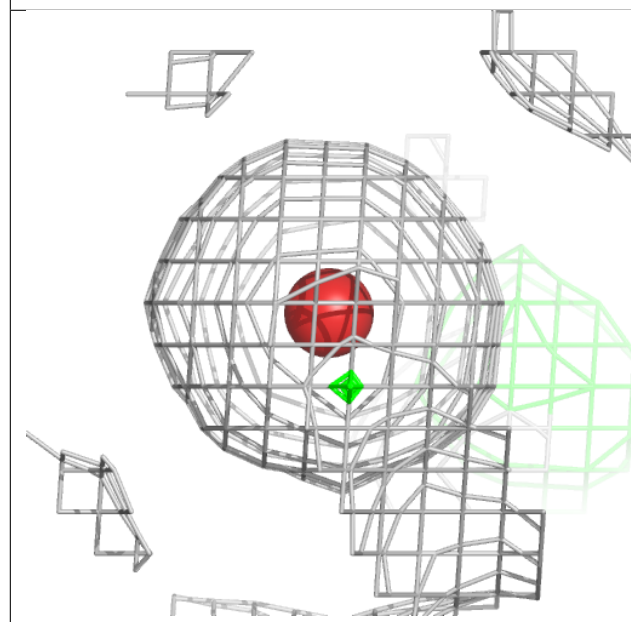
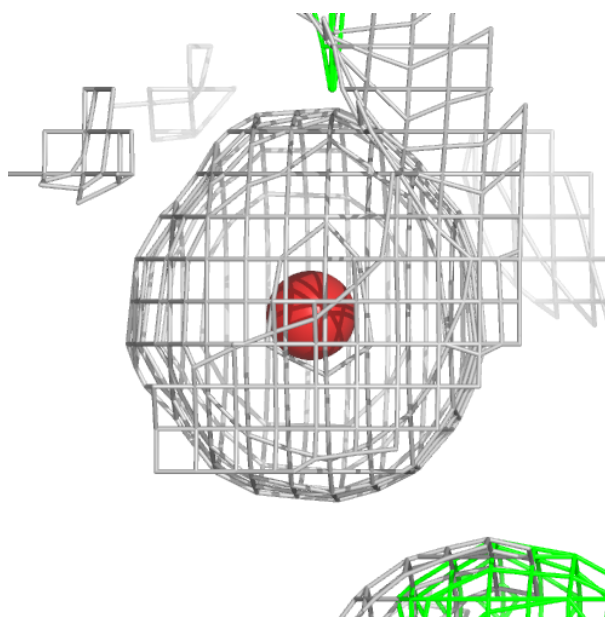
Electron density around BR A 502:

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



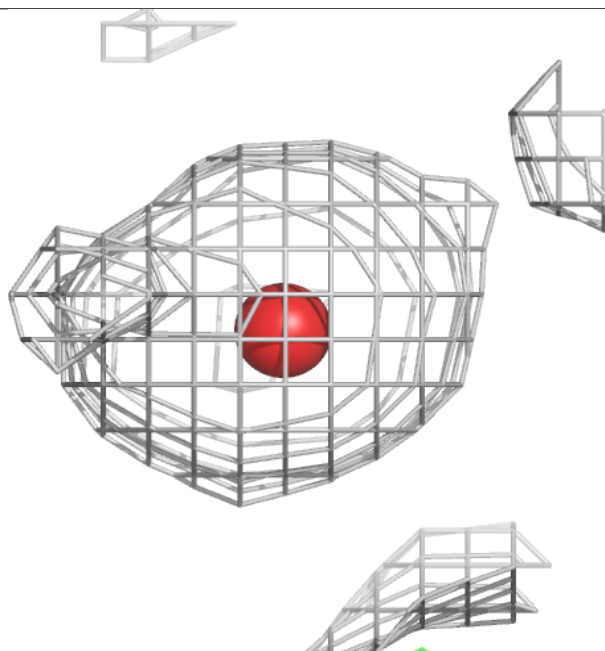
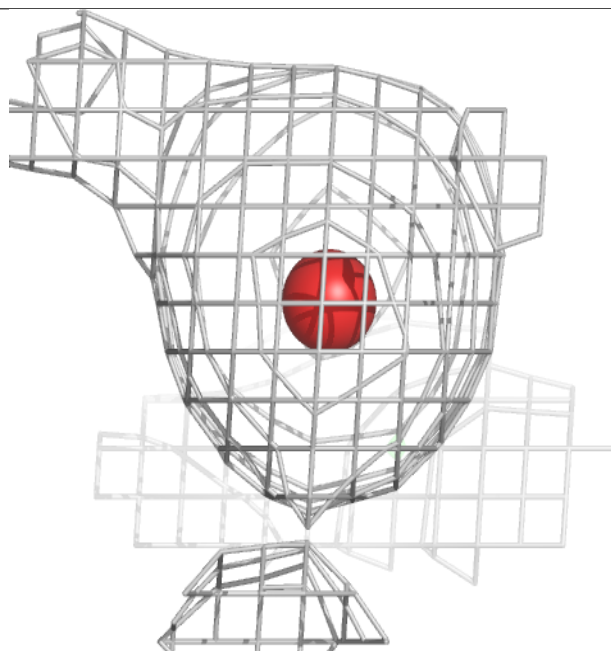
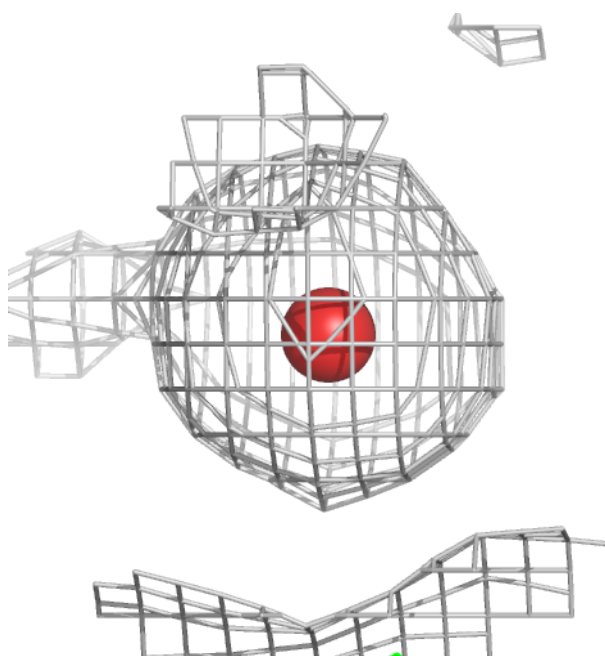
Electron density around BR A 503:

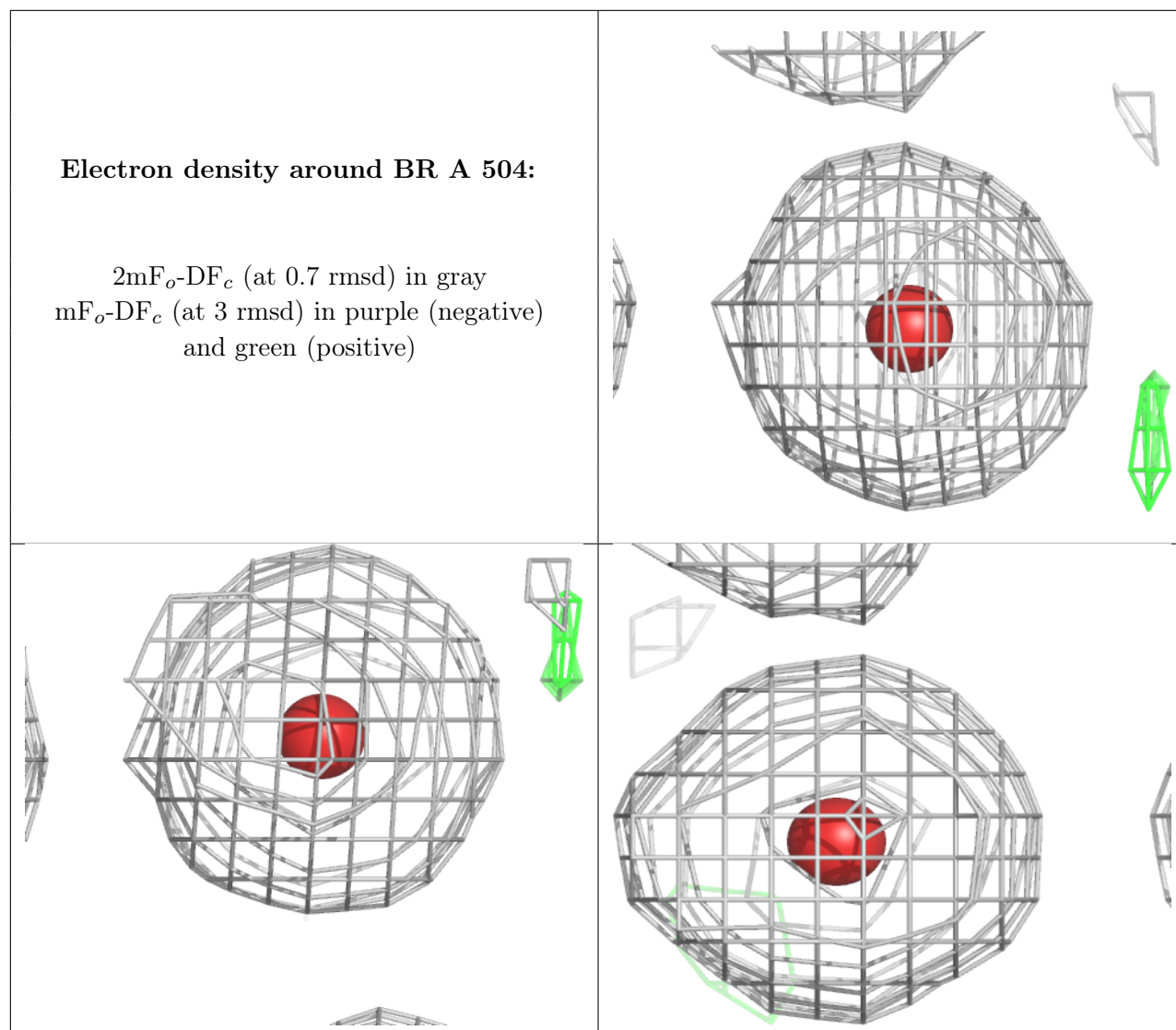
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around BR 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)





6.5 Other polymers ⓘ

There are no such residues in this entry.