Full wwPDB X-ray Structure Validation Report (i)

Dec 12, 2023 – 02:54 pm GMT

PDB ID	:	4D5D
Title	:	Crystal structure of CymA from Klebsiella oxytoca
Authors	:	van den Berg, B.; Bhamidimarri, S.P.; Kleinekathoefer, U.; Winterhalter, M.
Deposited on	:	2014-11-03
Resolution	:	1.90 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 (i) 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 (i)) were used in the production of this report:

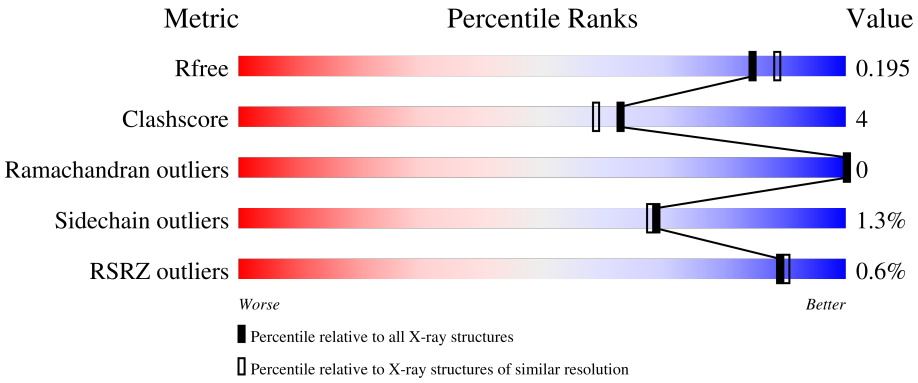
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.90 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}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

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 <=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	339	84% 8% 8%
1	B	339	% 84% 7% 9%
2	C	6	17% 67% 17%
2	E	6	50% 50%
3	D	7	14% 86%

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Mol	Chain	Length	Quality of chain
3	F	7	100%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 6314 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 CYMA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	311	Total	C	N	O	S	22	5	0
			2616	1669	437	508	2			
1	B	309	Total	C	N	O	S	27	3	0
			2589	1648	436	503	2			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-14	ALA	-	expression tag	UNP Q48391
A	-13	ASN	-	expression tag	UNP Q48391
A	-12	VAL	-	expression tag	UNP Q48391
A	-11	ARG	-	expression tag	UNP Q48391
A	-10	LEU	-	expression tag	UNP Q48391
A	-9	GLN	-	expression tag	UNP Q48391
A	-8	HIS	-	expression tag	UNP Q48391
A	-7	HIS	-	expression tag	UNP Q48391
A	-6	HIS	-	expression tag	UNP Q48391
A	-5	HIS	-	expression tag	UNP Q48391
A	-4	HIS	-	expression tag	UNP Q48391
A	-3	HIS	-	expression tag	UNP Q48391
A	-2	HIS	-	expression tag	UNP Q48391
A	-1	LEU	-	expression tag	UNP Q48391
A	0	GLU	-	expression tag	UNP Q48391
B	-14	ALA	-	expression tag	UNP Q48391
B	-13	ASN	-	expression tag	UNP Q48391
B	-12	VAL	-	expression tag	UNP Q48391
B	-11	ARG	-	expression tag	UNP Q48391
B	-10	LEU	-	expression tag	UNP Q48391
B	-9	GLN	-	expression tag	UNP Q48391
B	-8	HIS	-	expression tag	UNP Q48391
B	-7	HIS	-	expression tag	UNP Q48391
B	-6	HIS	-	expression tag	UNP Q48391
B	-5	HIS	-	expression tag	UNP Q48391

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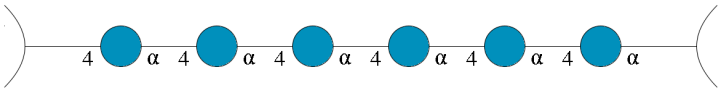
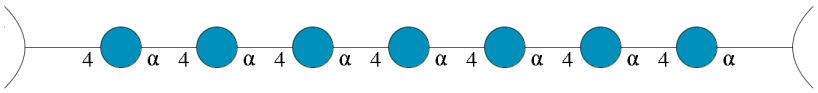
Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	HIS	-	expression tag	UNP Q48391
B	-3	HIS	-	expression tag	UNP Q48391
B	-2	HIS	-	expression tag	UNP Q48391
B	-1	LEU	-	expression tag	UNP Q48391
B	0	GLU	-	expression tag	UNP Q48391

- Molecule 2 is an oligosaccharide called Cyclohexakis-(1-4)-(alpha-D-glucopyranose).



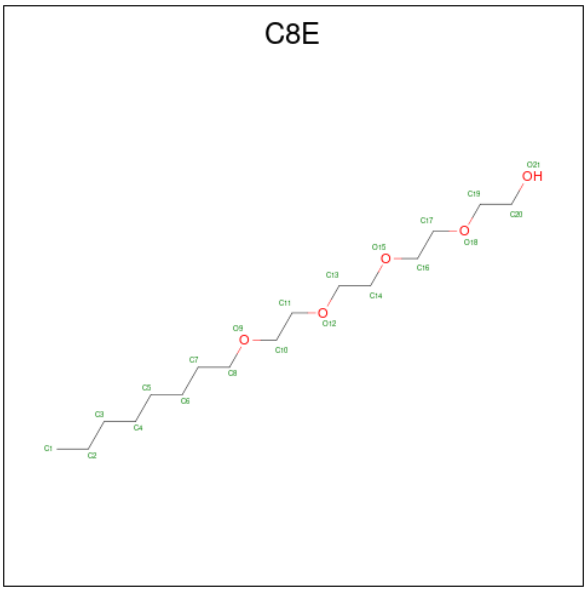
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
2	C	6	Total C O 66 36 30	0	0	0
2	E	6	Total C O 66 36 30	0	0	0

- Molecule 3 is an oligosaccharide called Cycloheptakis-(1-4)-(alpha-D-glucopyranose).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
3	D	7	Total C O 77 42 35	0	0	0
3	F	7	Total C O 77 42 35	0	0	0

- Molecule 4 is (HYDROXYETHYLOXY)TRI(ETHYLOXY)OCTANE (three-letter code: C8E) (formula: $C_{16}H_{34}O_5$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 14 10 4	0	0
4	A	1	Total C O 13 11 2	0	0
4	A	1	Total C O 13 11 2	0	0
4	A	1	Total C O 11 7 4	0	0
4	A	1	Total C O 13 9 4	0	0
4	A	1	Total C O 14 11 3	0	0
4	A	1	Total C O 18 14 4	3	0
4	A	1	Total C O 16 12 4	2	0
4	A	1	Total C O 8 5 3	0	0
4	A	1	Total C O 7 4 3	0	0
4	B	1	Total C O 13 11 2	0	0
4	B	1	Total C O 13 11 2	0	0
4	B	1	Total C O 14 12 2	2	0
4	B	1	Total C O 21 16 5	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 8 5 3	0	0
4	B	1	Total C O 13 11 2	3	0
4	B	1	Total C 8 8	0	0
4	B	1	Total C O 12 9 3	0	0
4	B	1	Total C 6 6	0	0
4	B	1	Total C O 12 10 2	4	0

- Molecule 5 is water.

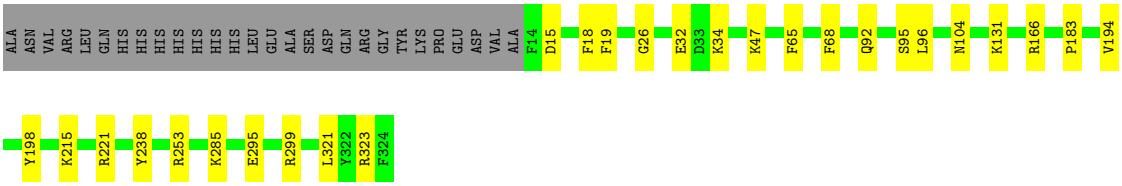
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	299	Total O 299 299	0	0
5	B	277	Total O 277 277	0	0

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.

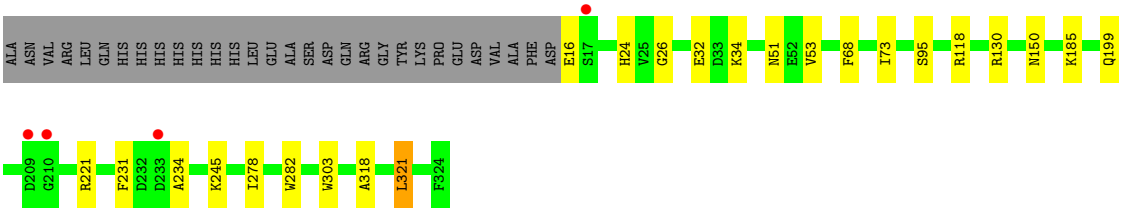
• Molecule 1: CYMA

Chain A:



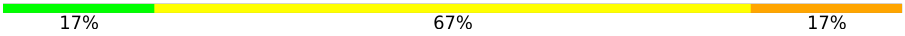
• Molecule 1: CYMA

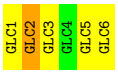
Chain B:



• Molecule 2: Cyclohexakis-(1-4)-(alpha-D-glucopyranose)

Chain C:



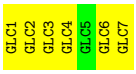


• Molecule 2: Cyclohexakis-(1-4)-(alpha-D-glucopyranose)

Chain E:

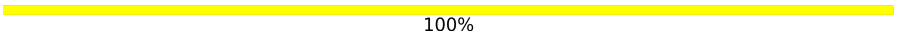
• Molecule 3: Cycloheptakis-(1-4)-(alpha-D-glucopyranose)

Chain D:



• Molecule 3: Cycloheptakis-(1-4)-(alpha-D-glucopyranose)

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α, β, γ	140.16Å 77.99Å 109.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.10 – 1.90 44.93 – 1.90	Depositor EDS
% Data completeness (in resolution range)	95.6 (47.10-1.90) 92.2 (44.93-1.90)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.35 (at 1.89Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.179 , 0.198 0.181 , 0.195	Depositor DCC
R_{free} test set	1975 reflections (2.14%)	wwPDB-VP
Wilson B-factor (Å^2)	27.1	Xtriage
Anisotropy	0.236	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 53.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	6314	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.74% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: C8E, GLC

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	0.41	0/2702	0.53	0/3644
1	B	0.41	0/2668	0.53	0/3597
All	All	0.41	0/5370	0.53	0/7241

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

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	2616	0	2451	20	0
1	B	2589	0	2422	23	0
2	C	66	0	54	1	0
2	E	66	0	54	3	0
3	D	77	0	63	1	0
3	F	77	0	63	0	0
4	A	127	0	164	6	0
4	B	120	0	191	11	0
5	A	299	0	0	10	1
5	B	277	0	0	3	1
All	All	6314	0	5462	48	1

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 4.

All (48) 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)
5:A:2294:HOH:O	3:D:3:GLC:O6	2.06	0.73
1:A:221:ARG:NH1	5:A:2207:HOH:O	2.22	0.71
1:B:221:ARG:NH1	5:B:2210:HOH:O	2.23	0.70
5:A:2288:HOH:O	1:B:118:ARG:NH2	2.26	0.68
1:A:295:GLU:OE2	5:A:2269:HOH:O	2.13	0.65
1:B:185:LYS:HZ2	4:B:1334:C8E:H102	1.60	0.64
1:A:253:ARG:HH11	1:A:253:ARG:HG2	1.67	0.59
1:A:323:ARG:NH2	5:A:2257:HOH:O	2.25	0.58
1:A:32:GLU:OE2	1:A:34:LYS:HD2	2.05	0.57
1:B:278:ILE:HG22	4:B:1328:C8E:H71	1.88	0.56
1:A:65:PHE:HZ	1:A:96:LEU:HD22	1.71	0.56
1:B:26:GLY:HA3	1:B:321:LEU:HD12	1.88	0.56
1:B:32:GLU:OE2	1:B:34:LYS:HD2	2.07	0.54
1:A:221:ARG:HD3	4:A:1329:C8E:H101	1.89	0.53
1:A:92:GLN:OE1	4:A:1327:C8E:H71	2.11	0.51
1:B:282:TRP:CD1	4:B:1328:C8E:H72	2.47	0.50
1:A:221:ARG:NH1	5:A:2203:HOH:O	2.45	0.49
1:B:68:PHE:HB3	1:B:95:SER:HB3	1.94	0.49
1:B:118:ARG:HG3	1:B:130:ARG:NH1	2.29	0.48
5:B:2275:HOH:O	2:E:2:GLC:H62	2.14	0.47
1:A:131:LYS:HG2	1:A:166:ARG:HG2	1.97	0.47
1:A:26:GLY:HA3	1:A:321:LEU:HD23	1.96	0.46
1:B:185:LYS:HZ3	4:B:1334:C8E:C4	2.29	0.46
1:B:278:ILE:CG2	4:B:1328:C8E:H71	2.46	0.46
1:B:199:GLN:NE2	5:B:2156:HOH:O	2.50	0.45
1:A:238:TYR:OH	1:A:285:LYS:HE3	2.18	0.44
1:A:47:LYS:NZ	5:A:2044:HOH:O	2.45	0.44
1:B:318:ALA:HB1	4:B:1330:C8E:H172	1.99	0.44
1:A:19:PHE:O	1:B:118:ARG:HD3	2.17	0.44
5:A:2160:HOH:O	2:C:2:GLC:H62	2.18	0.44
1:A:198:TYR:HB2	4:A:1329:C8E:H61	2.00	0.43
1:B:185:LYS:NZ	4:B:1334:C8E:H71	2.32	0.43
1:A:183:PRO:HD2	1:A:194:VAL:O	2.18	0.43
1:B:185:LYS:NZ	4:B:1334:C8E:H102	2.33	0.43
1:A:299:ARG:NH1	5:A:2271:HOH:O	2.33	0.43
1:B:73:ILE:CD1	4:B:1331:C8E:H101	2.49	0.42
1:A:215:LYS:HB2	1:A:215:LYS:HE2	1.83	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:1336:C8E:H111	1:B:53:VAL:HB	2.02	0.41
1:B:231:PHE:HB2	1:B:234:ALA:HB3	2.01	0.41
1:B:245:LYS:HE2	1:B:245:LYS:HB3	1.82	0.41
1:A:253:ARG:HG2	1:A:253:ARG:NH1	2.33	0.41
1:A:68:PHE:HB3	1:A:95:SER:HB3	2.01	0.41
1:B:303:TRP:CZ2	2:E:2:GLC:H4	2.56	0.41
1:B:24:HIS:CD2	1:B:321:LEU:HG	2.55	0.41
1:B:185:LYS:HD2	4:B:1334:C8E:H102	2.02	0.41
4:A:1331:C8E:H141	5:A:2146:HOH:O	2.21	0.40
4:A:1336:C8E:H101	4:B:1331:C8E:H131	2.03	0.40
2:E:1:GLC:O5	2:E:6:GLC:H61	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:A:2216:HOH:O	5:B:2052:HOH:O[2_555]	2.16	0.04

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	314/339 (93%)	310 (99%)	4 (1%)	0	100	100
1	B	310/339 (91%)	307 (99%)	3 (1%)	0	100	100
All	All	624/678 (92%)	617 (99%)	7 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	276/295 (94%)	273 (99%)	3 (1%)	73	73
1	B	272/295 (92%)	267 (98%)	5 (2%)	59	55
All	All	548/590 (93%)	540 (98%)	8 (2%)	69	62

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ASP
1	A	18	PHE
1	A	104	ASN
1	B	16	GLU
1	B	51[A]	ASN
1	B	51[B]	ASN
1	B	150	ASN
1	B	321	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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

26 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	GLC	C	1	2	11,11,12	0.39	0	15,15,17	2.34	6 (40%)
2	GLC	C	2	2	11,11,12	0.57	0	15,15,17	2.01	6 (40%)
2	GLC	C	3	2	11,11,12	0.57	0	15,15,17	2.39	5 (33%)
2	GLC	C	4	2	11,11,12	0.41	0	15,15,17	0.79	0
2	GLC	C	5	2	11,11,12	0.62	0	15,15,17	1.69	3 (20%)
2	GLC	C	6	2	11,11,12	0.64	0	15,15,17	1.43	3 (20%)
3	GLC	D	1	3	11,11,12	0.45	0	15,15,17	2.46	6 (40%)
3	GLC	D	2	3	11,11,12	0.41	0	15,15,17	0.97	1 (6%)
3	GLC	D	3	3	11,11,12	0.50	0	15,15,17	1.00	0
3	GLC	D	4	3	11,11,12	0.52	0	15,15,17	1.88	3 (20%)
3	GLC	D	5	3	11,11,12	0.43	0	15,15,17	0.59	0
3	GLC	D	6	3	11,11,12	0.53	0	15,15,17	1.69	5 (33%)
3	GLC	D	7	3	11,11,12	0.43	0	15,15,17	1.77	4 (26%)
2	GLC	E	1	2	11,11,12	0.44	0	15,15,17	1.84	4 (26%)
2	GLC	E	2	2	11,11,12	0.69	0	15,15,17	1.61	4 (26%)
2	GLC	E	3	2	11,11,12	0.59	0	15,15,17	2.27	4 (26%)
2	GLC	E	4	2	11,11,12	0.44	0	15,15,17	0.89	1 (6%)
2	GLC	E	5	2	11,11,12	0.50	0	15,15,17	1.66	2 (13%)
2	GLC	E	6	2	11,11,12	0.63	0	15,15,17	1.19	1 (6%)
3	GLC	F	1	3	11,11,12	0.49	0	15,15,17	2.22	5 (33%)
3	GLC	F	2	3	11,11,12	0.53	0	15,15,17	1.25	3 (20%)
3	GLC	F	3	3	11,11,12	0.39	0	15,15,17	1.04	1 (6%)
3	GLC	F	4	3	11,11,12	0.53	0	15,15,17	1.77	2 (13%)
3	GLC	F	5	3	11,11,12	0.47	0	15,15,17	1.03	1 (6%)
3	GLC	F	6	3	11,11,12	0.50	0	15,15,17	1.34	2 (13%)
3	GLC	F	7	3	11,11,12	0.55	0	15,15,17	2.05	4 (26%)

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	GLC	C	1	2	-	0/2/19/22	0/1/1/1
2	GLC	C	2	2	-	2/2/19/22	0/1/1/1
2	GLC	C	3	2	-	2/2/19/22	0/1/1/1
2	GLC	C	4	2	-	1/2/19/22	0/1/1/1
2	GLC	C	5	2	-	1/2/19/22	0/1/1/1
2	GLC	C	6	2	-	2/2/19/22	0/1/1/1
3	GLC	D	1	3	-	1/2/19/22	0/1/1/1
3	GLC	D	2	3	-	0/2/19/22	0/1/1/1
3	GLC	D	3	3	-	0/2/19/22	0/1/1/1
3	GLC	D	4	3	-	2/2/19/22	0/1/1/1
3	GLC	D	5	3	-	0/2/19/22	0/1/1/1
3	GLC	D	6	3	-	1/2/19/22	0/1/1/1
3	GLC	D	7	3	-	1/2/19/22	0/1/1/1
2	GLC	E	1	2	-	1/2/19/22	0/1/1/1
2	GLC	E	2	2	-	2/2/19/22	0/1/1/1
2	GLC	E	3	2	-	0/2/19/22	0/1/1/1
2	GLC	E	4	2	-	2/2/19/22	0/1/1/1
2	GLC	E	5	2	-	0/2/19/22	0/1/1/1
2	GLC	E	6	2	-	2/2/19/22	0/1/1/1
3	GLC	F	1	3	-	0/2/19/22	0/1/1/1
3	GLC	F	2	3	-	2/2/19/22	0/1/1/1
3	GLC	F	3	3	-	0/2/19/22	0/1/1/1
3	GLC	F	4	3	-	2/2/19/22	0/1/1/1
3	GLC	F	5	3	-	0/2/19/22	0/1/1/1
3	GLC	F	6	3	-	0/2/19/22	0/1/1/1
3	GLC	F	7	3	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (76) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	E	3	GLC	C1-O5-C5	6.73	121.31	112.19
2	C	3	GLC	C1-O5-C5	5.98	120.29	112.19
3	D	1	GLC	C1-O5-C5	5.70	119.91	112.19
3	D	4	GLC	C1-O5-C5	5.68	119.89	112.19
2	C	2	GLC	C1-O5-C5	5.42	119.53	112.19
2	C	1	GLC	C1-O5-C5	5.03	119.01	112.19
3	F	4	GLC	C1-O5-C5	4.91	118.84	112.19
3	F	1	GLC	C1-O5-C5	4.78	118.67	112.19
3	F	7	GLC	C1-O5-C5	4.74	118.61	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	1	GLC	C3-C4-C5	4.54	118.33	110.24
2	E	5	GLC	C1-O5-C5	4.47	118.25	112.19
3	D	7	GLC	C1-O5-C5	4.19	117.87	112.19
3	D	1	GLC	C3-C4-C5	4.09	117.53	110.24
3	D	1	GLC	O5-C1-C2	3.92	116.82	110.77
3	F	1	GLC	O5-C5-C6	3.79	113.14	107.20
3	F	7	GLC	C3-C4-C5	3.78	116.98	110.24
2	C	5	GLC	C1-O5-C5	3.77	117.30	112.19
2	E	2	GLC	C1-O5-C5	3.73	117.24	112.19
3	D	6	GLC	C1-O5-C5	3.67	117.17	112.19
2	E	1	GLC	C1-O5-C5	3.67	117.16	112.19
2	E	1	GLC	C3-C4-C5	3.64	116.73	110.24
3	F	6	GLC	C1-O5-C5	3.59	117.06	112.19
2	C	3	GLC	C3-C4-C5	3.54	116.56	110.24
3	F	7	GLC	O5-C5-C4	3.29	118.82	110.83
2	C	5	GLC	O5-C5-C4	3.27	118.78	110.83
3	F	1	GLC	O5-C1-C2	3.23	115.76	110.77
2	E	1	GLC	O5-C5-C6	3.15	112.15	107.20
2	E	5	GLC	O5-C5-C4	3.04	118.23	110.83
3	F	3	GLC	O4-C4-C3	-3.02	103.36	110.35
2	C	3	GLC	C1-C2-C3	3.02	113.38	109.67
3	F	1	GLC	C3-C4-C5	2.95	115.50	110.24
2	E	3	GLC	C3-C4-C5	2.94	115.49	110.24
2	E	2	GLC	O4-C4-C3	-2.94	103.56	110.35
3	F	7	GLC	C1-C2-C3	2.93	113.27	109.67
2	E	6	GLC	O5-C5-C6	2.92	111.78	107.20
2	C	1	GLC	O5-C5-C6	2.90	111.74	107.20
2	C	3	GLC	C2-C3-C4	2.89	115.90	110.89
3	F	1	GLC	O4-C4-C3	-2.85	103.75	110.35
3	F	2	GLC	O4-C4-C3	-2.74	104.02	110.35
2	C	1	GLC	O4-C4-C3	-2.71	104.08	110.35
2	C	2	GLC	O5-C5-C4	2.69	117.38	110.83
3	D	2	GLC	O4-C4-C3	-2.67	104.17	110.35
3	D	6	GLC	O2-C2-C3	-2.66	104.82	110.14
2	C	6	GLC	O5-C5-C6	2.61	111.29	107.20
3	D	7	GLC	C1-C2-C3	2.60	112.87	109.67
2	C	5	GLC	C3-C4-C5	2.56	114.81	110.24
2	E	2	GLC	O5-C5-C4	2.55	117.03	110.83
3	D	1	GLC	O5-C5-C6	2.55	111.20	107.20
2	C	3	GLC	O5-C1-C2	2.51	114.64	110.77
3	F	2	GLC	C6-C5-C4	2.50	118.87	113.00
2	C	1	GLC	O5-C1-C2	2.47	114.58	110.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	D	4	GLC	O5-C5-C4	2.46	116.80	110.83
2	C	6	GLC	O5-C1-C2	2.41	114.49	110.77
2	E	2	GLC	C6-C5-C4	2.36	118.53	113.00
2	C	2	GLC	O4-C4-C3	-2.34	104.93	110.35
2	C	6	GLC	O4-C4-C5	-2.31	103.57	109.30
3	D	7	GLC	O5-C5-C4	2.30	116.42	110.83
3	D	4	GLC	C3-C4-C5	2.27	114.29	110.24
3	F	5	GLC	C3-C4-C5	2.26	114.28	110.24
2	E	4	GLC	O4-C4-C3	-2.24	105.17	110.35
3	D	1	GLC	C1-C2-C3	2.23	112.41	109.67
3	D	7	GLC	C3-C4-C5	2.21	114.18	110.24
2	C	2	GLC	C3-C4-C5	2.20	114.17	110.24
2	E	1	GLC	O5-C1-C2	2.17	114.12	110.77
3	F	2	GLC	O4-C4-C5	2.17	114.68	109.30
2	C	2	GLC	O5-C1-C2	2.14	114.08	110.77
2	C	1	GLC	O2-C2-C3	-2.14	105.85	110.14
2	C	2	GLC	C1-C2-C3	-2.13	107.05	109.67
3	F	4	GLC	O5-C1-C2	2.12	114.05	110.77
3	D	1	GLC	O5-C5-C4	2.12	115.98	110.83
3	F	6	GLC	O5-C1-C2	2.10	114.01	110.77
2	E	3	GLC	O5-C5-C4	2.07	115.86	110.83
2	E	3	GLC	O5-C1-C2	2.05	113.94	110.77
3	D	6	GLC	C1-C2-C3	2.03	112.17	109.67
3	D	6	GLC	O5-C1-C2	2.03	113.90	110.77
3	D	6	GLC	O5-C5-C4	2.03	115.76	110.83

There are no chirality outliers.

All (26) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	4	GLC	O5-C5-C6-O6
3	D	4	GLC	C4-C5-C6-O6
2	E	2	GLC	O5-C5-C6-O6
3	F	2	GLC	C4-C5-C6-O6
2	C	2	GLC	C4-C5-C6-O6
2	E	2	GLC	C4-C5-C6-O6
2	C	2	GLC	O5-C5-C6-O6
3	F	7	GLC	C4-C5-C6-O6
2	C	6	GLC	O5-C5-C6-O6
2	C	6	GLC	C4-C5-C6-O6
3	F	2	GLC	O5-C5-C6-O6
2	C	3	GLC	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	D	1	GLC	C4-C5-C6-O6
3	F	4	GLC	O5-C5-C6-O6
2	E	6	GLC	C4-C5-C6-O6
3	F	4	GLC	C4-C5-C6-O6
3	F	7	GLC	O5-C5-C6-O6
2	C	3	GLC	C4-C5-C6-O6
2	E	6	GLC	O5-C5-C6-O6
2	E	4	GLC	C4-C5-C6-O6
2	E	4	GLC	O5-C5-C6-O6
2	C	4	GLC	C4-C5-C6-O6
2	E	1	GLC	C4-C5-C6-O6
3	D	7	GLC	C4-C5-C6-O6
3	D	6	GLC	C4-C5-C6-O6
2	C	5	GLC	O5-C5-C6-O6

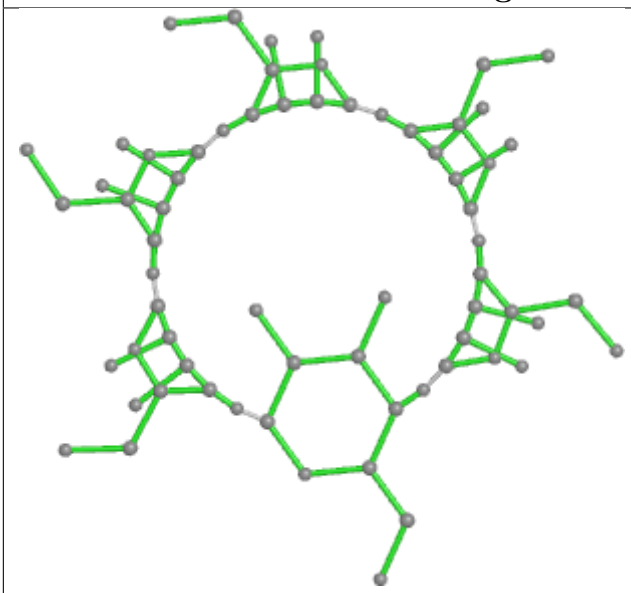
There are no ring outliers.

5 monomers are involved in 5 short contacts:

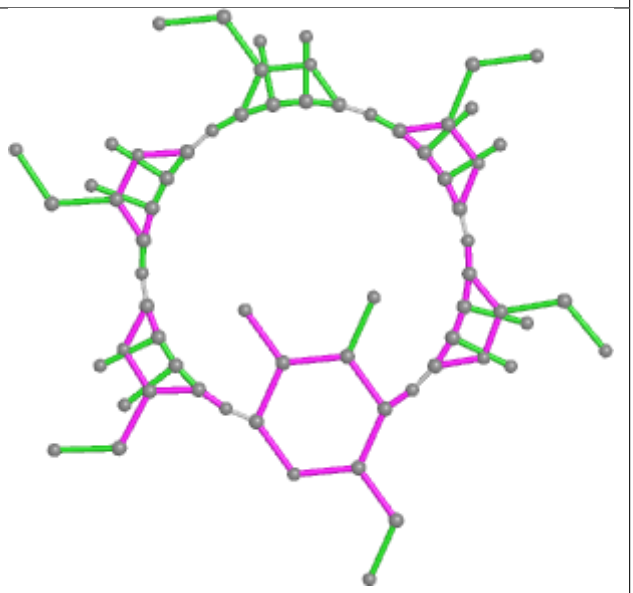
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	2	GLC	2	0
2	E	6	GLC	1	0
2	E	1	GLC	1	0
3	D	3	GLC	1	0
2	C	2	GLC	1	0

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

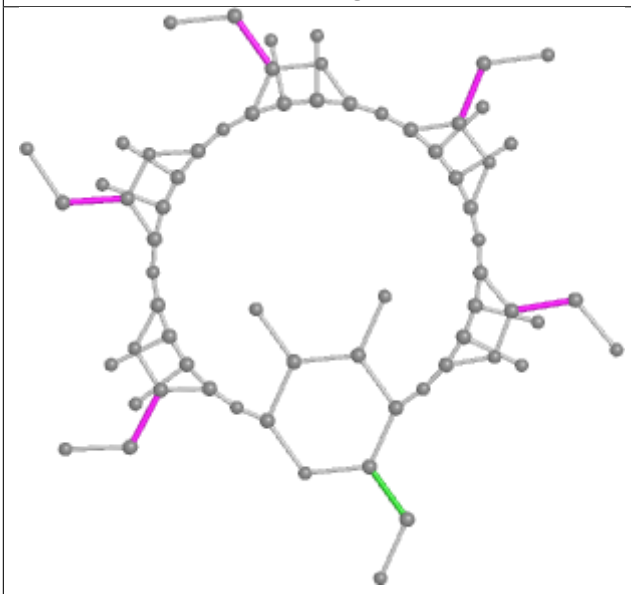
Oligosaccharide Chain C



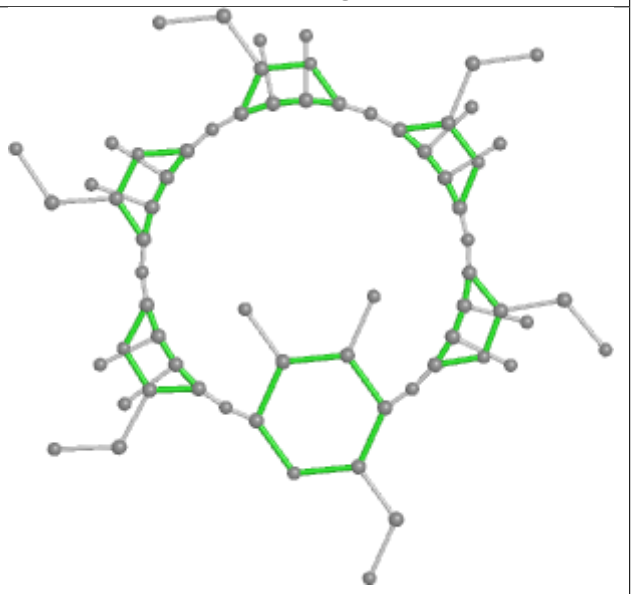
Bond lengths



Bond angles

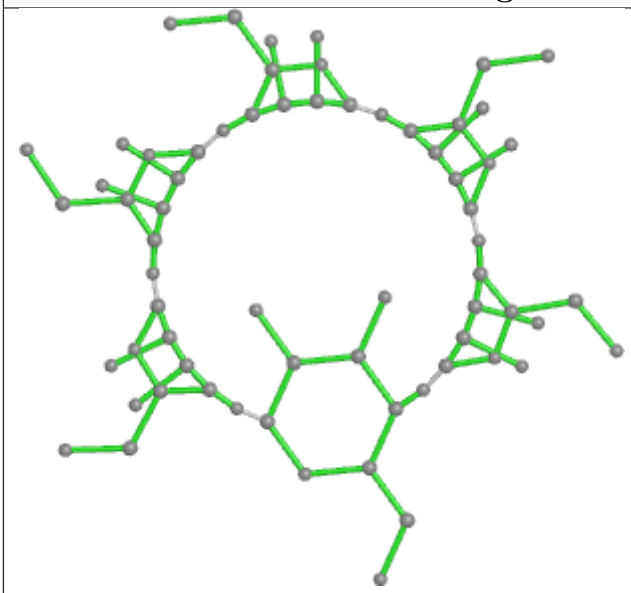


Torsions

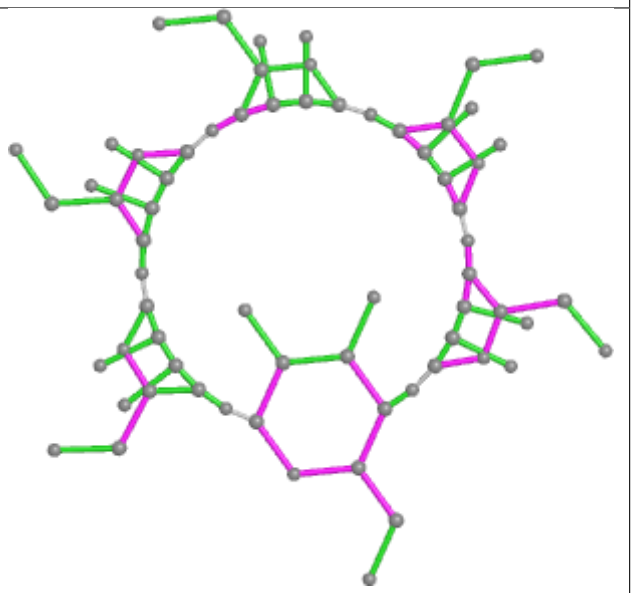


Rings

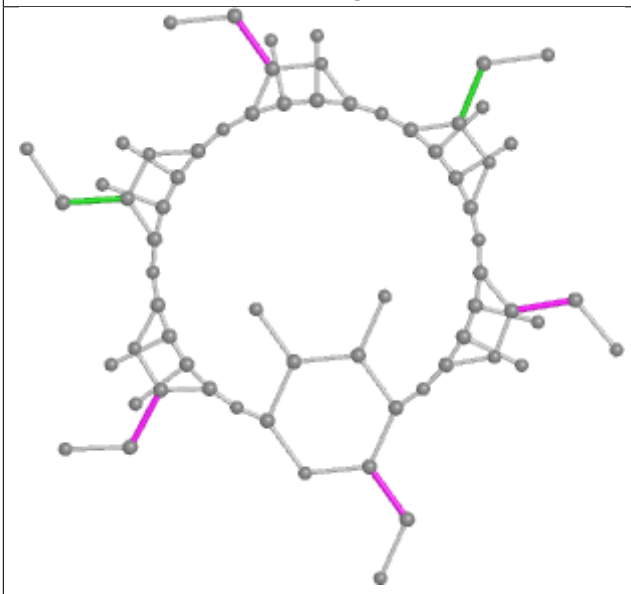
Oligosaccharide Chain E



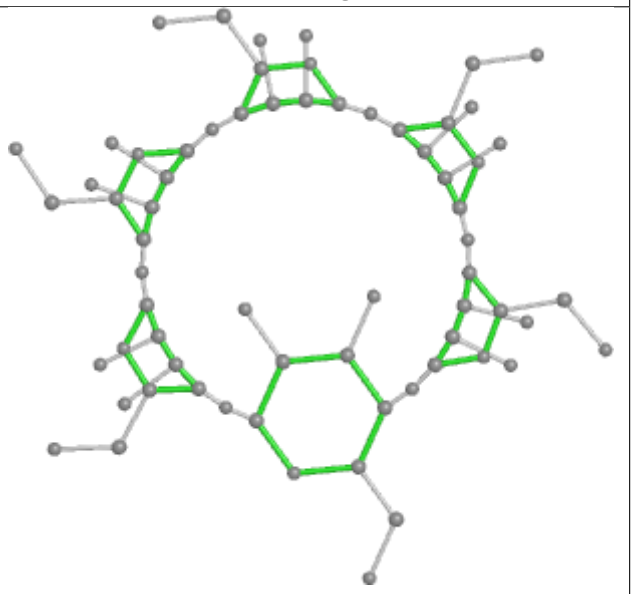
Bond lengths



Bond angles

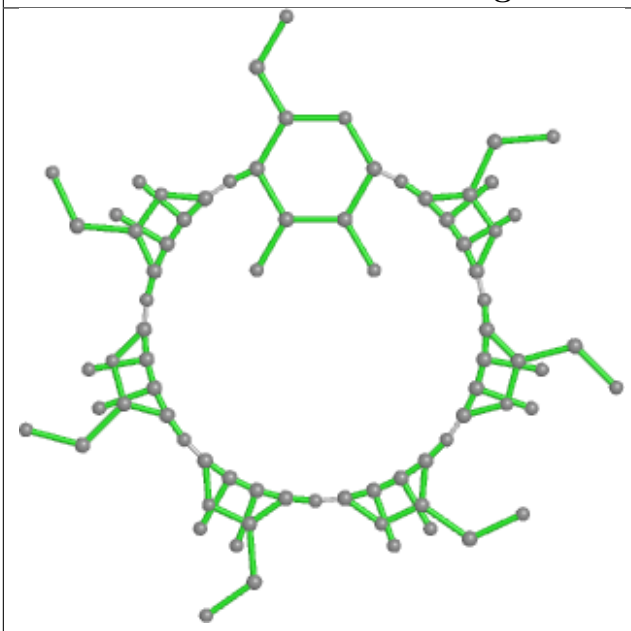


Torsions

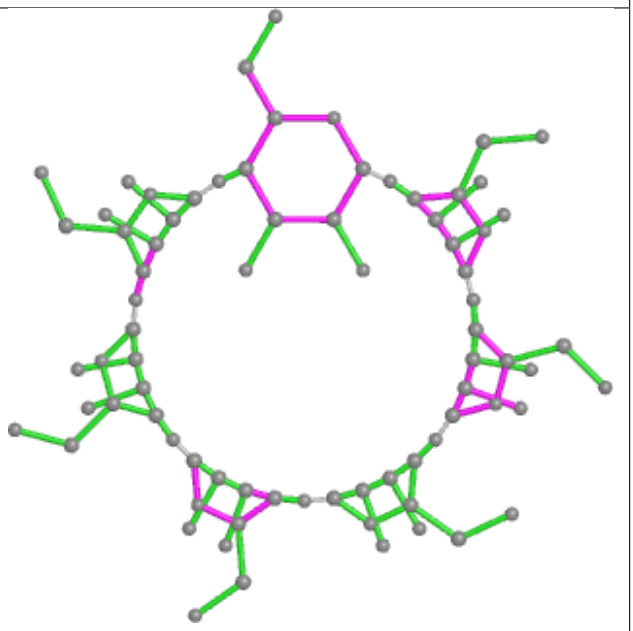


Rings

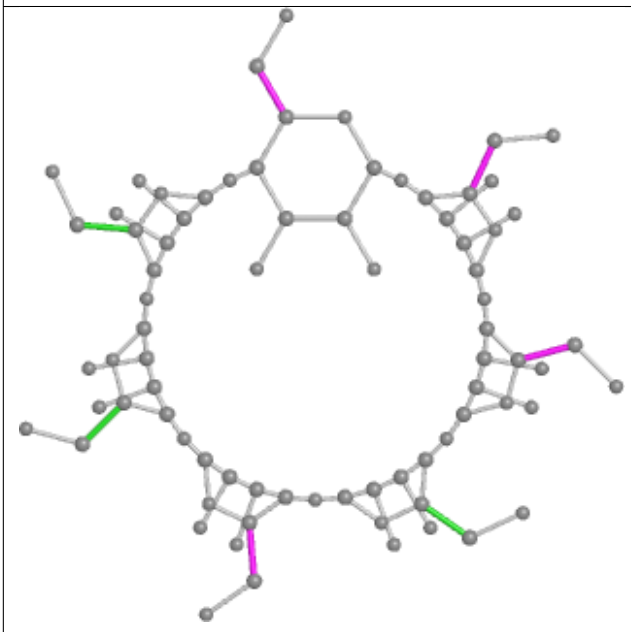
Oligosaccharide Chain D



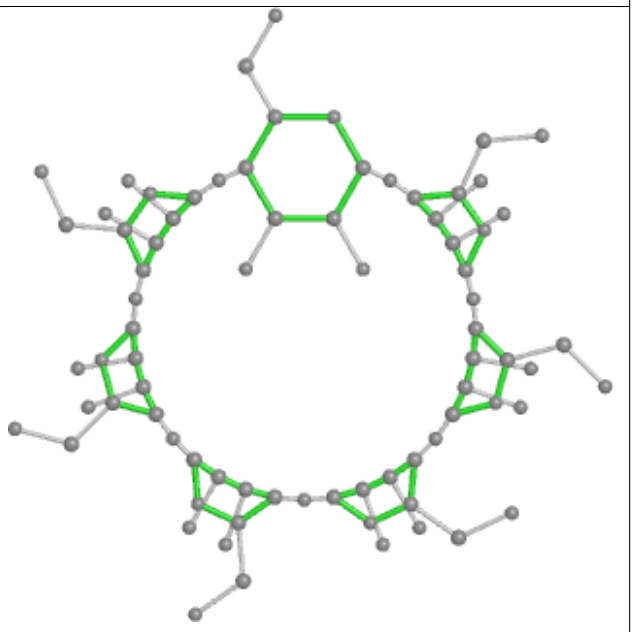
Bond lengths



Bond angles

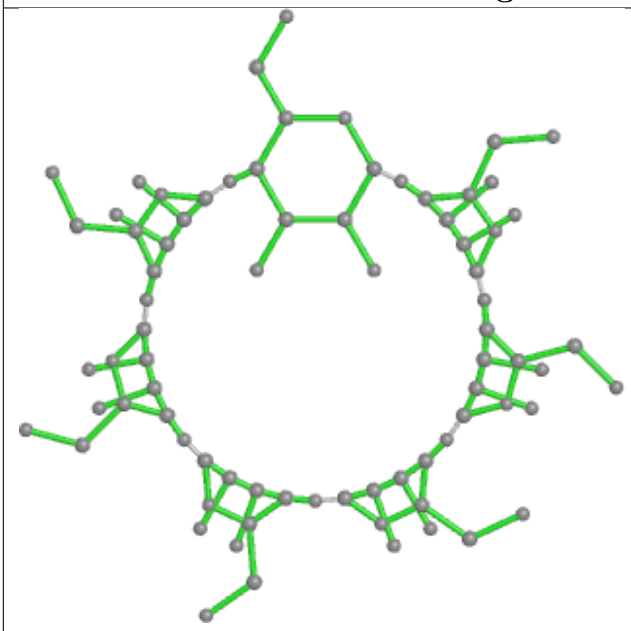


Torsions

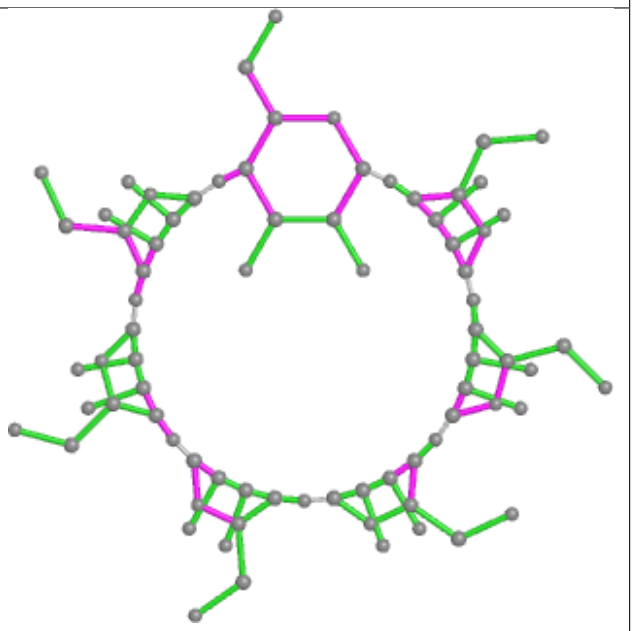


Rings

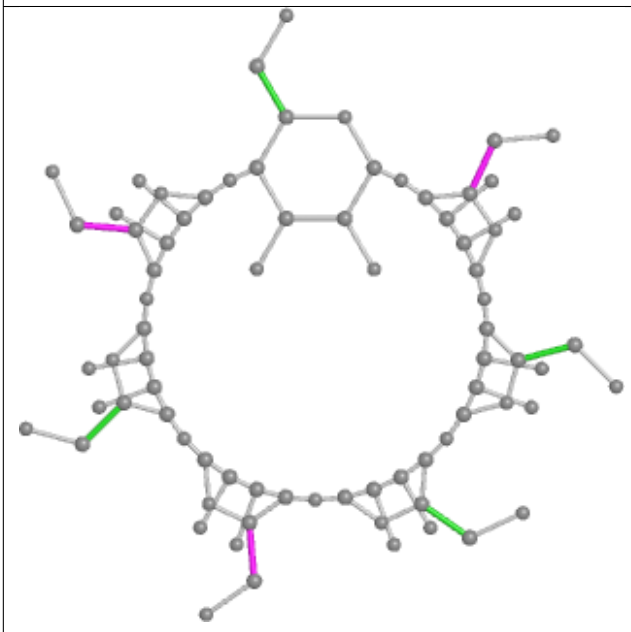
Oligosaccharide Chain F



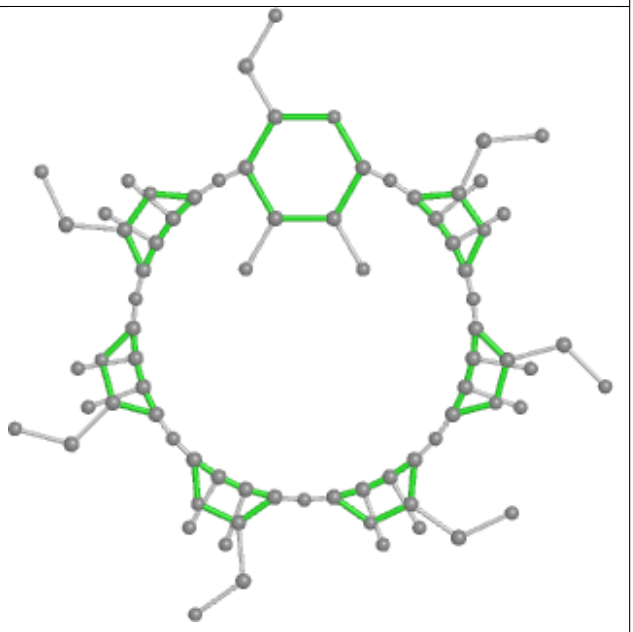
Bond lengths



Bond angles



Torsions



Rings

5.6 Ligand geometry

20 ligands 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	$\#\|Z\|>2$
4	C8E	A	1331	-	12,12,20	0.54	0	11,11,19	0.29	0
4	C8E	B	1329	-	13,13,20	0.39	0	12,12,19	0.46	0
4	C8E	B	1334	-	11,11,20	0.38	0	10,10,19	0.27	0
4	C8E	B	1328	-	12,12,20	0.32	0	11,11,19	0.48	0
4	C8E	A	1334	-	15,15,20	0.48	0	14,14,19	0.21	0
4	C8E	A	1336	-	6,6,20	0.40	0	5,5,19	0.30	0
4	C8E	B	1336	-	11,11,20	0.33	0	10,10,19	0.53	0
4	C8E	B	1327	-	12,12,20	0.29	0	11,11,19	0.63	0
4	C8E	B	1333	-	7,7,20	0.29	0	6,6,19	0.38	0
4	C8E	B	1335	-	5,5,20	0.32	0	4,4,19	0.33	0
4	C8E	B	1330	-	20,20,20	0.46	0	19,19,19	0.55	0
4	C8E	A	1332	-	13,13,20	0.39	0	12,12,19	0.41	0
4	C8E	A	1329	-	12,12,20	0.39	0	11,11,19	0.58	0
4	C8E	A	1328	-	11,11,20	0.44	0	9,9,19	0.31	0
4	C8E	B	1332	-	12,12,20	0.37	0	11,11,19	0.29	0
4	C8E	A	1327	-	13,13,20	0.42	0	12,12,19	0.33	0
4	C8E	B	1331	-	7,7,20	0.42	0	6,6,19	0.45	0
4	C8E	A	1333	-	17,17,20	0.35	0	16,16,19	0.52	0
4	C8E	A	1330	-	10,10,20	0.42	0	9,9,19	0.33	0
4	C8E	A	1335	-	7,7,20	0.37	0	6,6,19	0.32	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
4	C8E	A	1331	-	-	1/10/10/18	-
4	C8E	B	1329	-	-	6/11/11/18	-
4	C8E	B	1334	-	-	4/9/9/18	-
4	C8E	B	1328	-	-	4/10/10/18	-
4	C8E	A	1334	-	-	7/13/13/18	-
4	C8E	A	1336	-	-	2/4/4/18	-
4	C8E	B	1336	-	-	5/9/9/18	-
4	C8E	B	1327	-	-	7/10/10/18	-
4	C8E	B	1333	-	-	1/5/5/18	-
4	C8E	B	1335	-	-	1/3/3/18	-
4	C8E	B	1330	-	-	7/18/18/18	-
4	C8E	A	1332	-	-	6/11/11/18	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	C8E	A	1329	-	-	6/10/10/18	-
4	C8E	A	1328	-	-	2/7/7/18	-
4	C8E	B	1332	-	-	7/10/10/18	-
4	C8E	A	1327	-	-	4/11/11/18	-
4	C8E	B	1331	-	-	2/5/5/18	-
4	C8E	A	1333	-	-	6/15/15/18	-
4	C8E	A	1330	-	-	5/8/8/18	-
4	C8E	A	1335	-	-	4/5/5/18	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (87) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1334	C8E	O12-C13-C14-O15
4	A	1334	C8E	O15-C16-C17-O18
4	A	1333	C8E	O12-C13-C14-O15
4	A	1334	C8E	O9-C10-C11-O12
4	A	1334	C8E	O12-C13-C14-O15
4	A	1330	C8E	O12-C13-C14-O15
4	B	1330	C8E	O15-C16-C17-O18
4	B	1327	C8E	C6-C7-C8-O9
4	A	1330	C8E	O15-C16-C17-O18
4	A	1329	C8E	O9-C10-C11-O12
4	A	1335	C8E	O9-C10-C11-O12
4	A	1336	C8E	O12-C13-C14-O15
4	B	1332	C8E	C2-C3-C4-C5
4	B	1327	C8E	C4-C5-C6-C7
4	B	1328	C8E	C2-C3-C4-C5
4	B	1335	C8E	C2-C3-C4-C5
4	B	1327	C8E	C2-C3-C4-C5
4	B	1332	C8E	C4-C5-C6-C7
4	B	1333	C8E	C4-C5-C6-C7
4	B	1336	C8E	C3-C4-C5-C6
4	A	1332	C8E	C6-C7-C8-O9
4	A	1332	C8E	O9-C10-C11-O12
4	A	1332	C8E	C5-C6-C7-C8
4	B	1327	C8E	C3-C4-C5-C6

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Mol	Chain	Res	Type	Atoms
4	A	1329	C8E	C5-C6-C7-C8
4	A	1329	C8E	C4-C5-C6-C7
4	B	1330	C8E	C3-C4-C5-C6
4	B	1329	C8E	C2-C3-C4-C5
4	B	1330	C8E	C6-C7-C8-O9
4	B	1334	C8E	C4-C5-C6-C7
4	B	1327	C8E	O9-C10-C11-O12
4	B	1330	C8E	C2-C3-C4-C5
4	B	1331	C8E	O9-C10-C11-O12
4	B	1332	C8E	C5-C6-C7-C8
4	B	1332	C8E	C3-C4-C5-C6
4	A	1329	C8E	C3-C4-C5-C6
4	A	1332	C8E	O12-C13-C14-O15
4	A	1329	C8E	C2-C3-C4-C5
4	B	1328	C8E	O9-C10-C11-O12
4	A	1332	C8E	C2-C3-C4-C5
4	B	1336	C8E	O9-C10-C11-O12
4	A	1327	C8E	C6-C7-C8-O9
4	A	1336	C8E	C10-C11-O12-C13
4	A	1333	C8E	C1-C2-C3-C4
4	B	1329	C8E	C6-C7-C8-O9
4	A	1330	C8E	C16-C17-O18-C19
4	B	1336	C8E	C1-C2-C3-C4
4	B	1334	C8E	C5-C6-C7-C8
4	A	1334	C8E	C20-C19-O18-C17
4	B	1330	C8E	C4-C5-C6-C7
4	A	1335	C8E	C13-C14-O15-C16
4	A	1330	C8E	C10-C11-O12-C13
4	A	1327	C8E	C14-C13-O12-C11
4	A	1332	C8E	C10-C11-O12-C13
4	B	1328	C8E	C6-C7-C8-O9
4	A	1331	C8E	C17-C16-O15-C14
4	A	1327	C8E	C13-C14-O15-C16
4	A	1327	C8E	C7-C8-O9-C10
4	A	1334	C8E	C7-C8-O9-C10
4	A	1333	C8E	C7-C8-O9-C10
4	B	1336	C8E	C7-C8-O9-C10
4	B	1327	C8E	C1-C2-C3-C4
4	A	1333	C8E	C10-C11-O12-C13
4	B	1330	C8E	O18-C19-C20-O21
4	B	1332	C8E	C1-C2-C3-C4
4	B	1327	C8E	C10-C11-O12-C13

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Mol	Chain	Res	Type	Atoms
4	A	1328	C8E	C1-C2-C3-C4
4	B	1329	C8E	C4-C5-C6-C7
4	B	1328	C8E	C10-C11-O12-C13
4	A	1333	C8E	C13-C14-O15-C16
4	A	1329	C8E	C10-C11-O12-C13
4	B	1331	C8E	C14-C13-O12-C11
4	A	1334	C8E	C17-C16-O15-C14
4	A	1335	C8E	C14-C13-O12-C11
4	A	1330	C8E	C13-C14-O15-C16
4	A	1335	C8E	O12-C13-C14-O15
4	B	1332	C8E	C7-C8-O9-C10
4	B	1334	C8E	C10-C11-O12-C13
4	A	1334	C8E	C6-C7-C8-O9
4	B	1332	C8E	O9-C10-C11-O12
4	B	1329	C8E	O9-C10-C11-O12
4	B	1330	C8E	O12-C13-C14-O15
4	B	1336	C8E	C2-C3-C4-C5
4	B	1329	C8E	C7-C8-O9-C10
4	A	1328	C8E	O9-C10-C11-O12
4	B	1329	C8E	C1-C2-C3-C4
4	A	1333	C8E	C17-C16-O15-C14

There are no ring outliers.

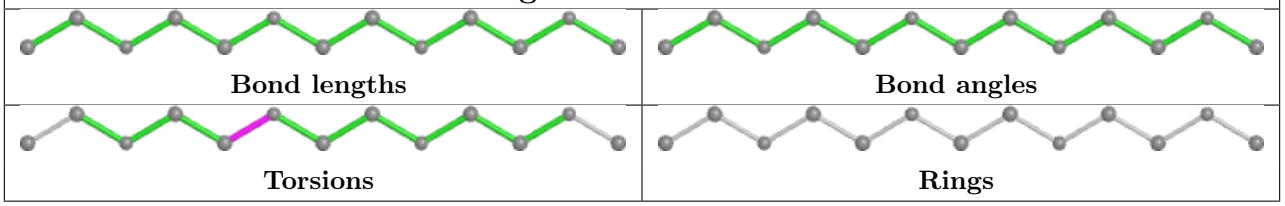
8 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1331	C8E	1	0
4	B	1334	C8E	5	0
4	B	1328	C8E	3	0
4	A	1336	C8E	2	0
4	B	1330	C8E	1	0
4	A	1329	C8E	2	0
4	A	1327	C8E	1	0
4	B	1331	C8E	2	0

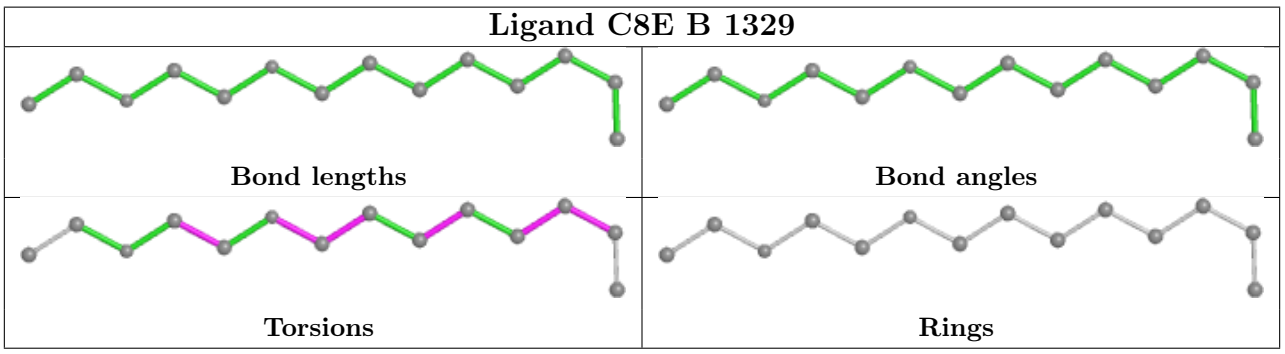
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

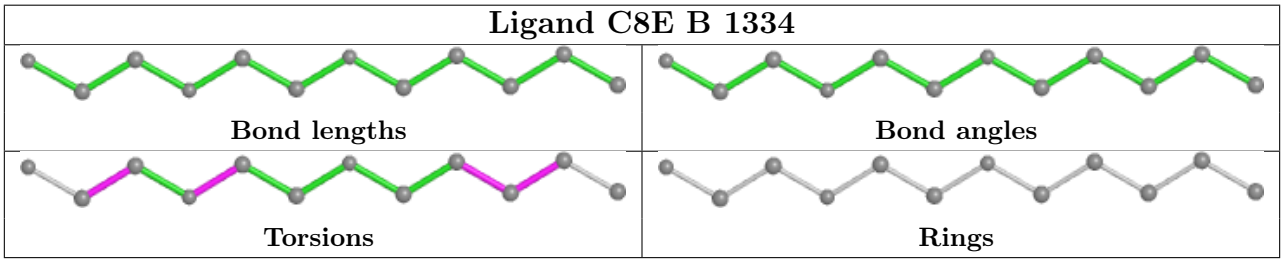
Ligand C8E A 1331



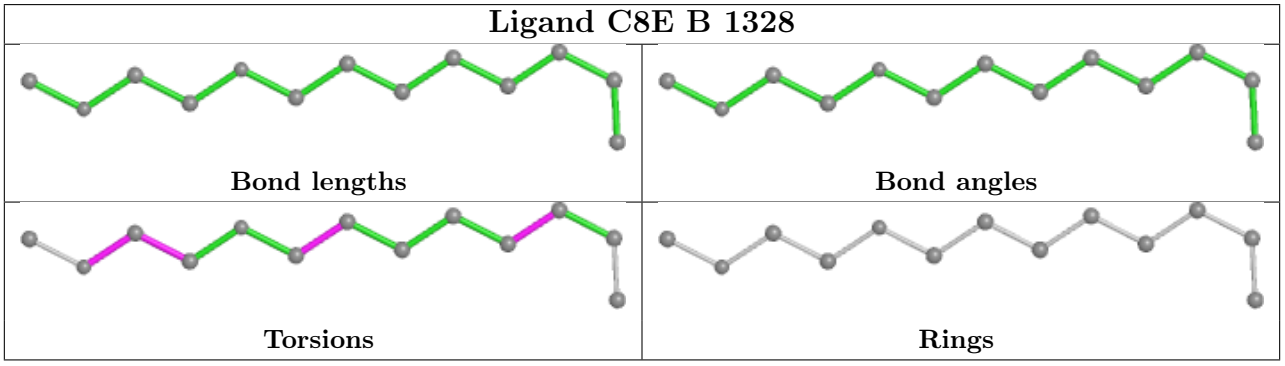
Ligand C8E B 1329



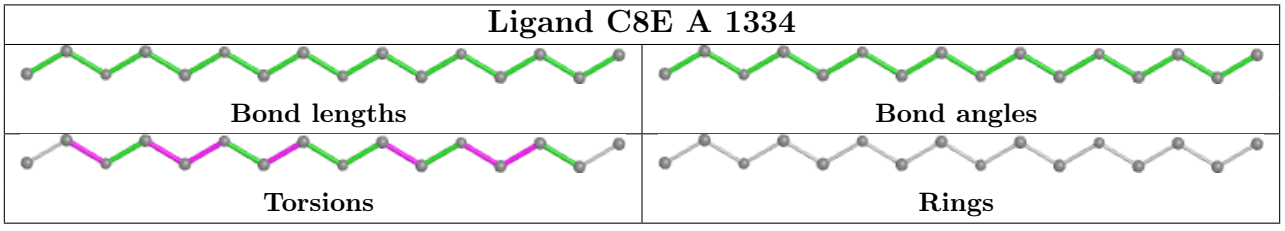
Ligand C8E B 1334



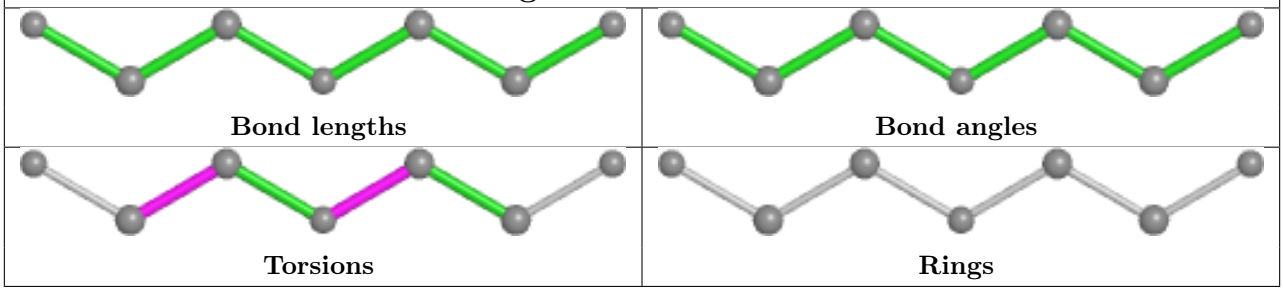
Ligand C8E B 1328



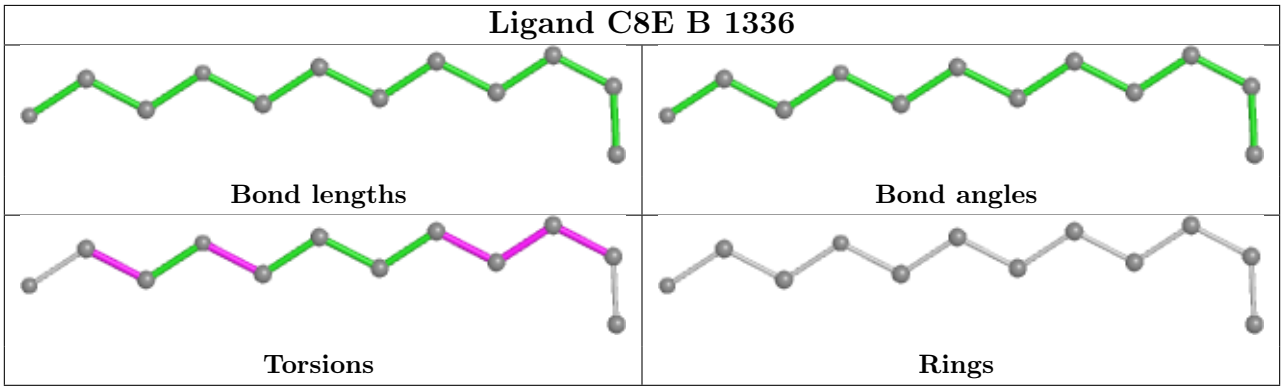
Ligand C8E A 1334



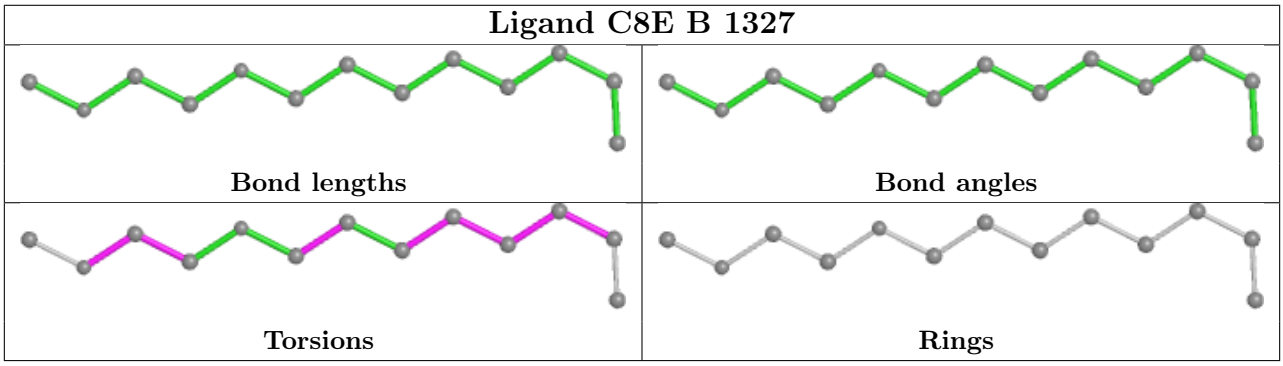
Ligand C8E A 1336



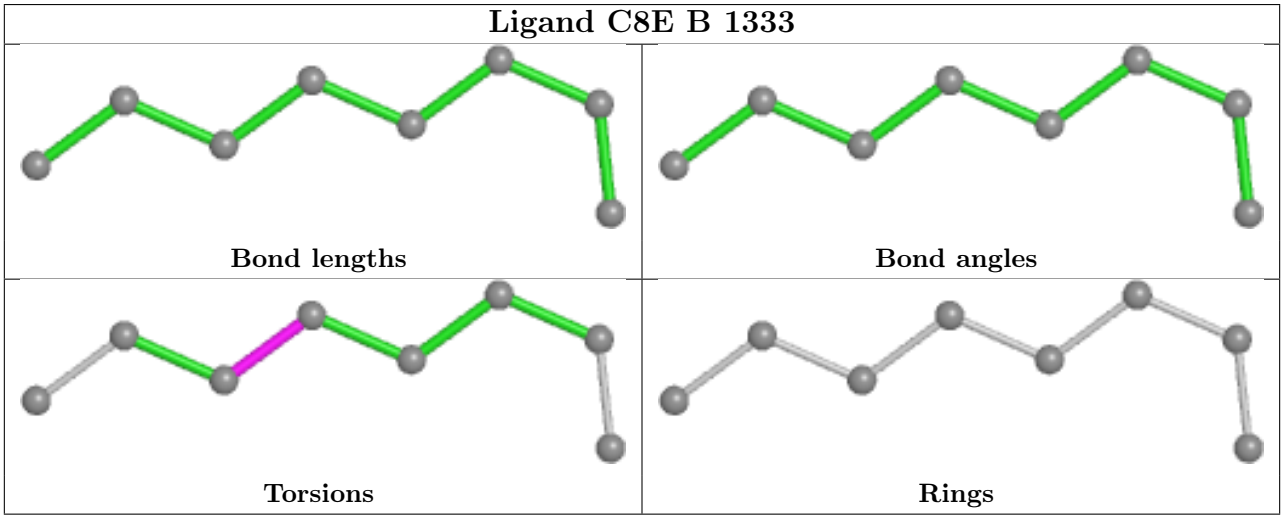
Ligand C8E B 1336



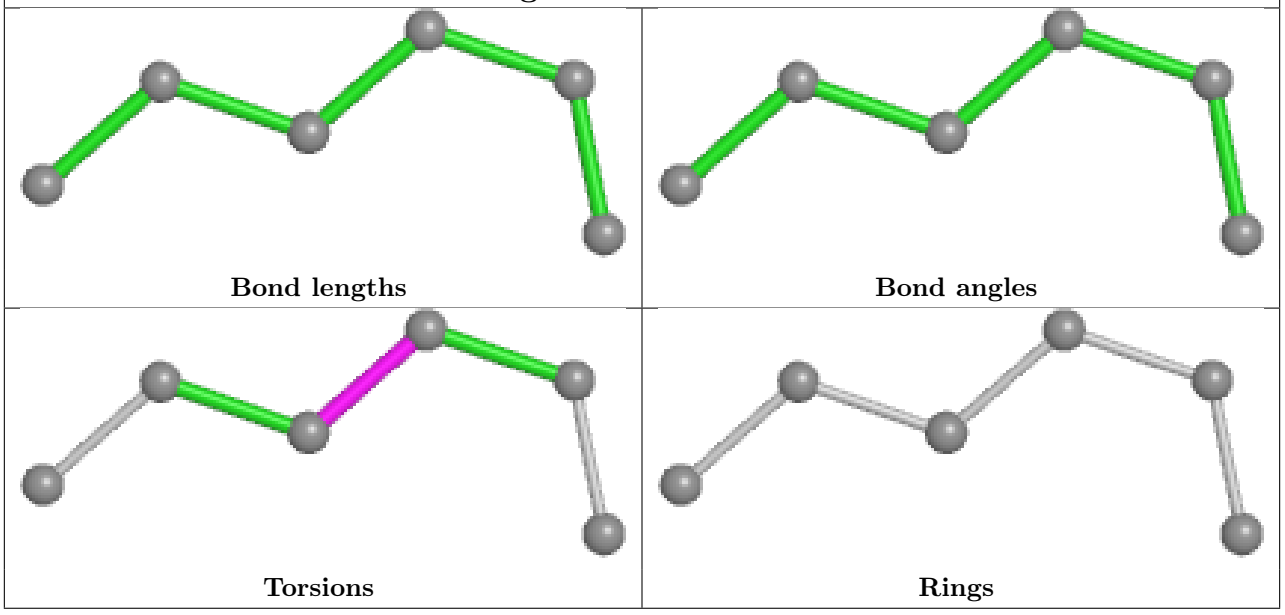
Ligand C8E B 1327



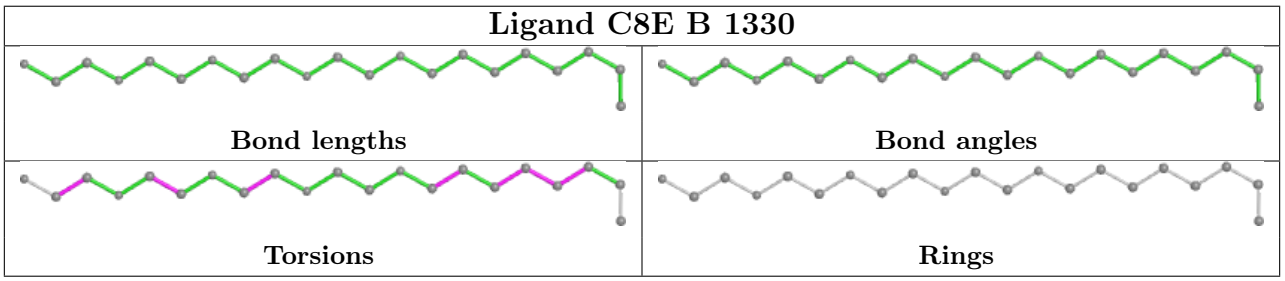
Ligand C8E B 1333



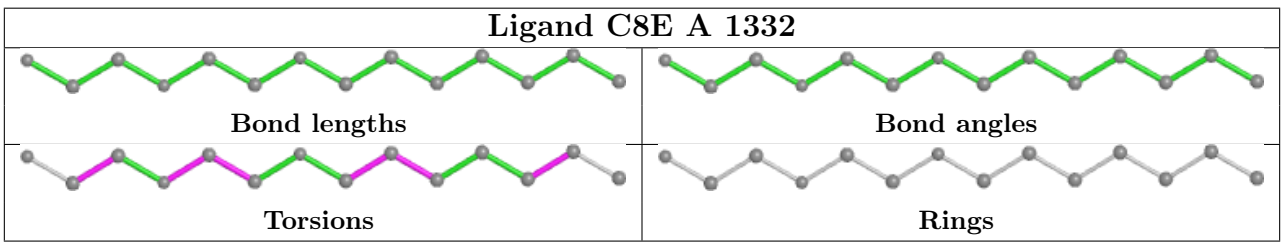
Ligand C8E B 1335



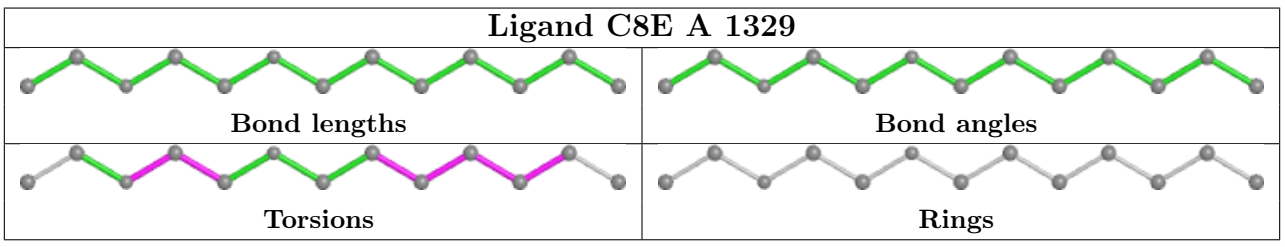
Ligand C8E B 1330



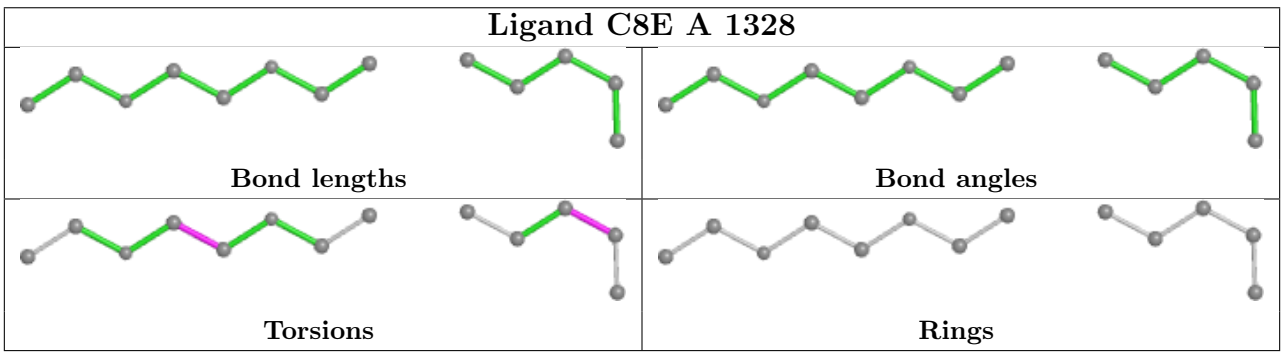
Ligand C8E A 1332



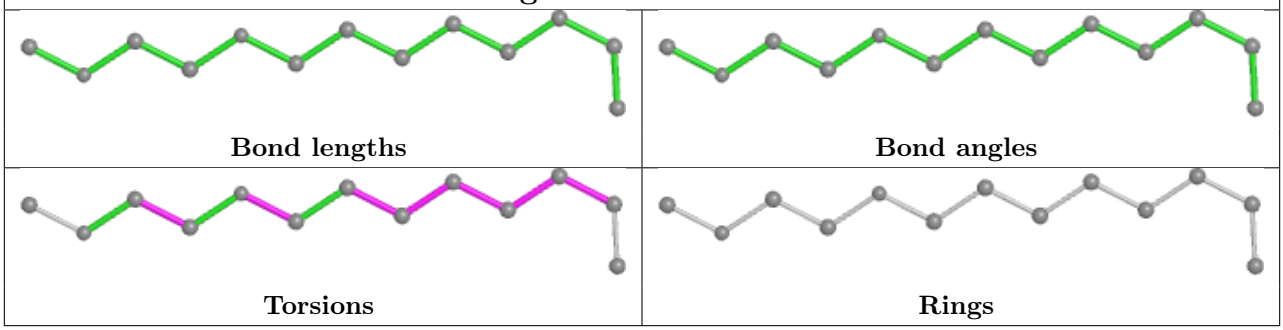
Ligand C8E A 1329



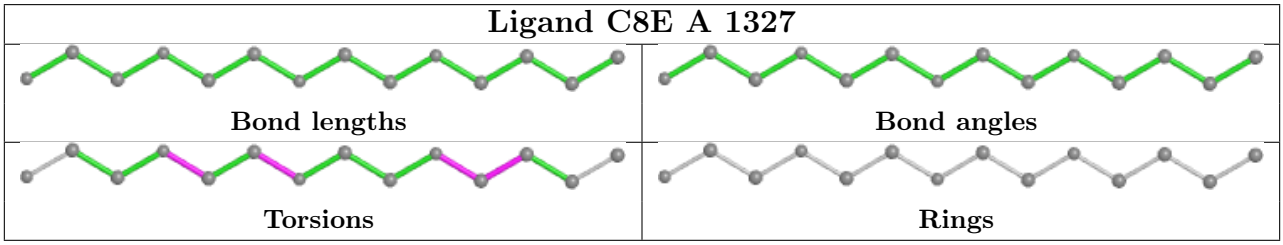
Ligand C8E A 1328



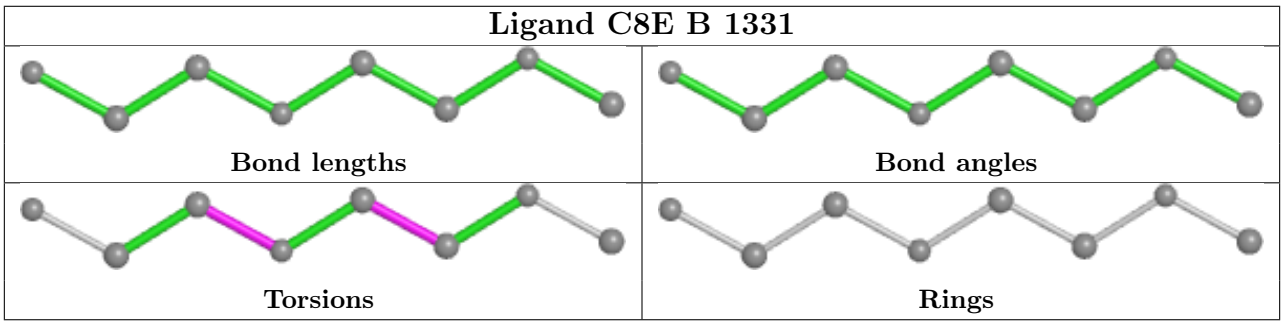
Ligand C8E B 1332



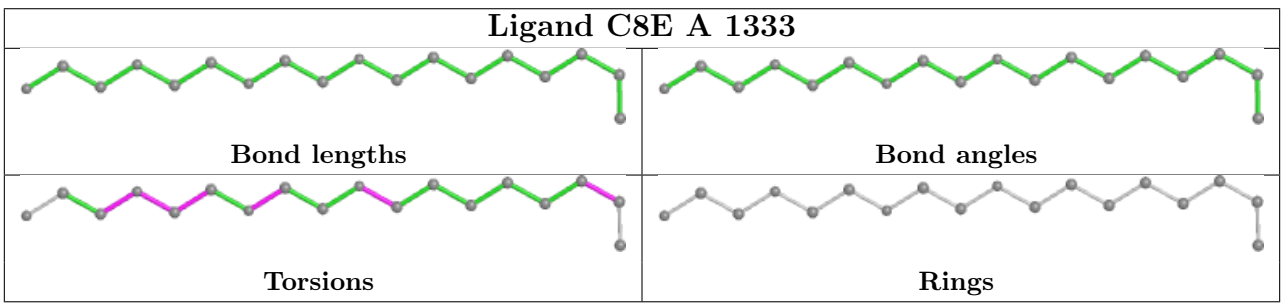
Ligand C8E A 1327



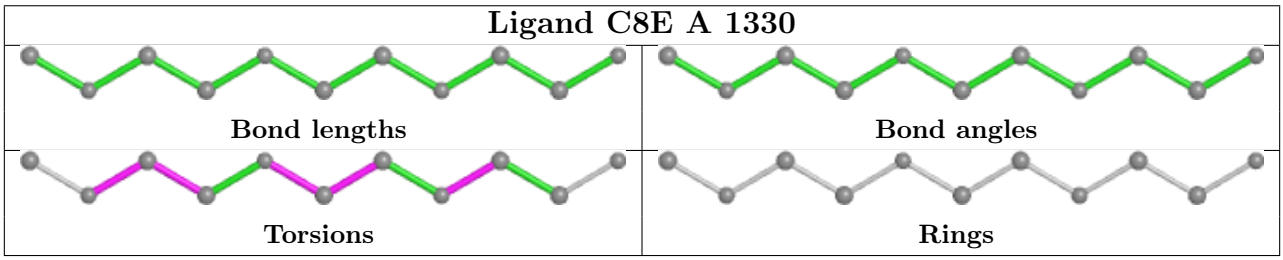
Ligand C8E B 1331

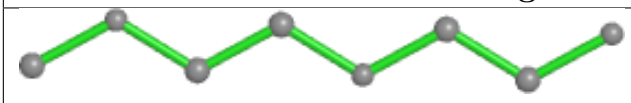
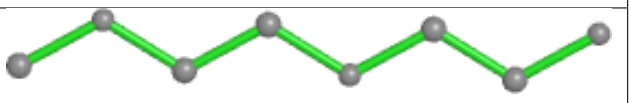
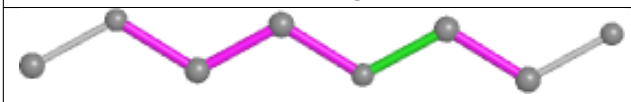
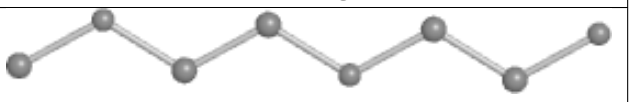


Ligand C8E A 1333



Ligand C8E A 1330



Ligand C8E A 1335	
 Bond lengths	 Bond angles
 Torsions	 Rings

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 95^{th} 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	311/339 (91%)	-0.14	0 100 100	22, 31, 47, 82	5 (1%)
1	B	309/339 (91%)	-0.19	4 (1%) 77 79	23, 32, 48, 72	9 (2%)
All	All	620/678 (91%)	-0.16	4 (0%) 89 90	22, 32, 48, 82	14 (2%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	17	SER	4.6
1	B	210	GLY	2.6
1	B	209	ASP	2.4
1	B	233	ASP	2.1

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

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

6.3 Carbohydrates

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, 95^{th} 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	GLC	F	5	11/12	0.76	0.15	45,55,62,63	0
3	GLC	D	2	11/12	0.82	0.19	46,57,72,84	0
3	GLC	D	4	11/12	0.84	0.25	50,55,63,66	0
2	GLC	C	2	11/12	0.84	0.21	52,62,70,77	0
3	GLC	F	1	11/12	0.85	0.25	55,59,65,69	0

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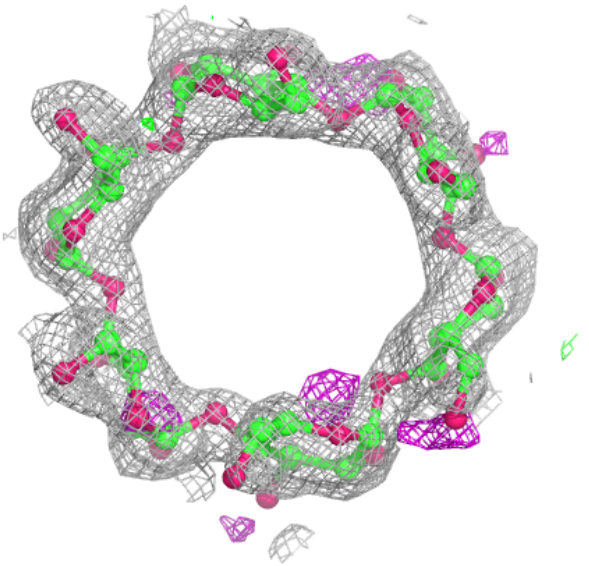
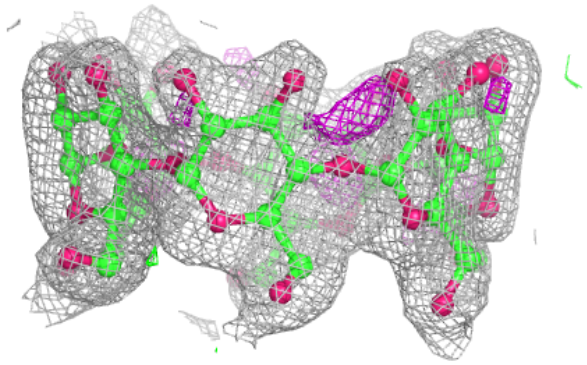
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	GLC	F	7	11/12	0.85	0.27	57,63,67,77	0
2	GLC	C	5	11/12	0.87	0.15	48,52,61,66	0
2	GLC	E	2	11/12	0.87	0.12	43,45,49,50	0
3	GLC	F	2	11/12	0.87	0.23	43,59,70,79	0
3	GLC	F	4	11/12	0.87	0.20	59,63,67,69	0
3	GLC	D	1	11/12	0.87	0.16	52,56,73,78	0
2	GLC	C	3	11/12	0.87	0.27	59,67,70,77	0
3	GLC	D	5	11/12	0.88	0.34	51,55,64,67	0
2	GLC	E	6	11/12	0.88	0.16	49,53,59,59	0
3	GLC	F	3	11/12	0.89	0.12	46,53,56,57	0
3	GLC	D	6	11/12	0.89	0.28	46,58,72,77	0
2	GLC	C	4	11/12	0.89	0.19	45,48,56,61	0
2	GLC	C	6	11/12	0.89	0.15	40,44,47,51	0
3	GLC	F	6	11/12	0.90	0.10	33,39,44,45	0
3	GLC	D	7	11/12	0.90	0.16	42,47,53,55	0
3	GLC	D	3	11/12	0.91	0.17	37,43,48,53	0
2	GLC	C	1	11/12	0.91	0.20	47,51,59,63	0
2	GLC	E	1	11/12	0.92	0.15	50,55,62,64	0
2	GLC	E	5	11/12	0.93	0.10	39,45,55,61	0
2	GLC	E	3	11/12	0.93	0.11	37,43,45,49	0
2	GLC	E	4	11/12	0.96	0.09	34,37,43,46	0

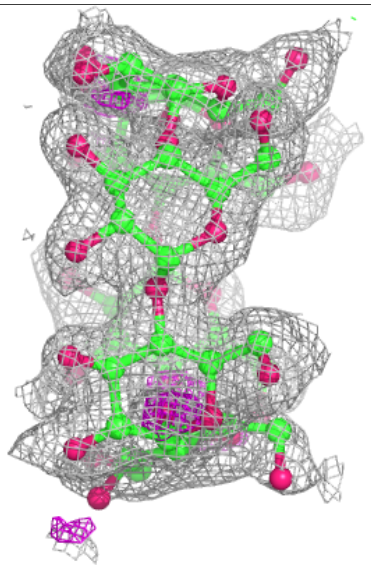
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain C:

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

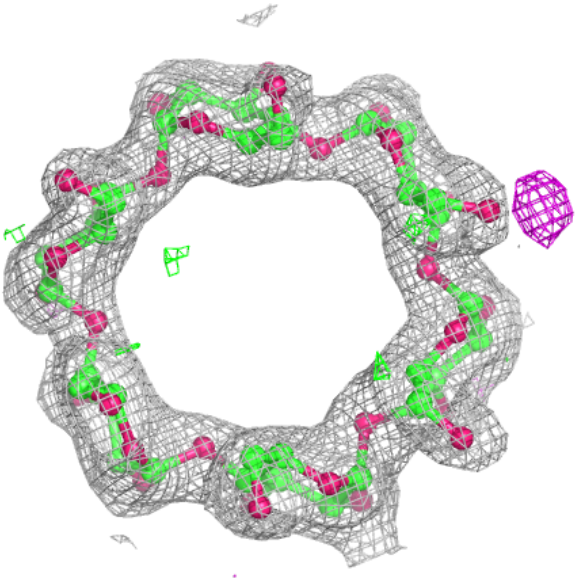


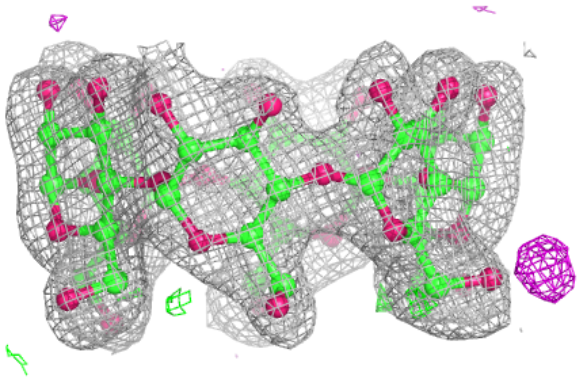


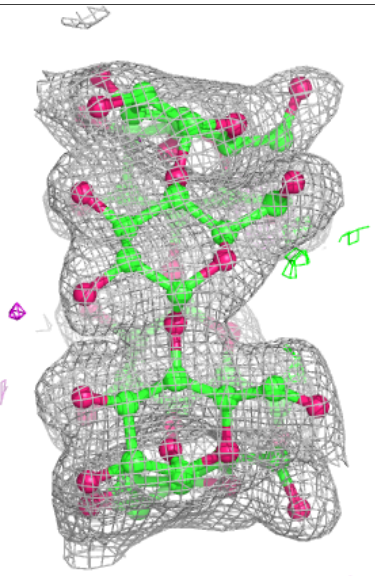


Electron density around Chain E:

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

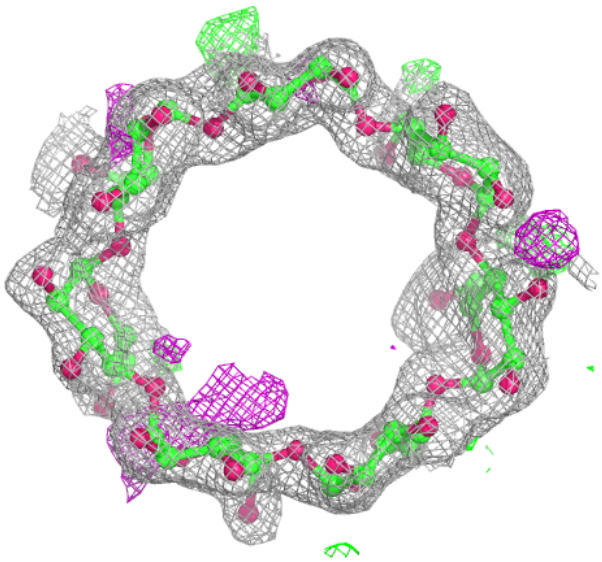


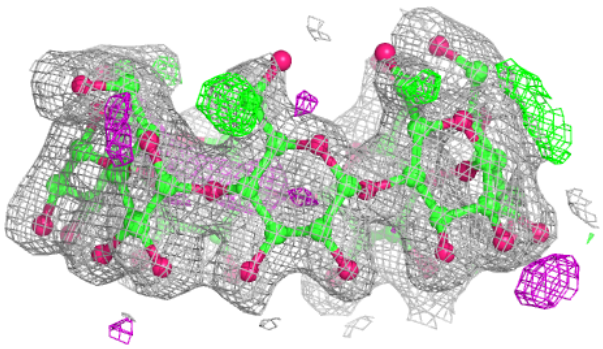


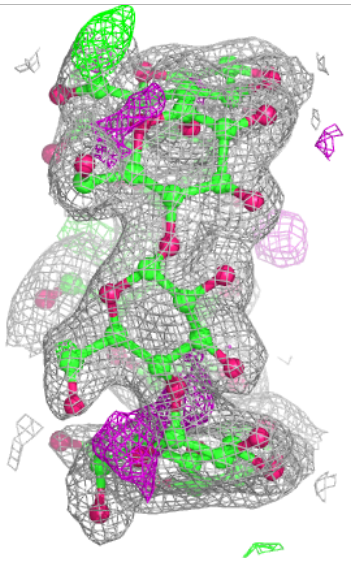


Electron density around Chain D:

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

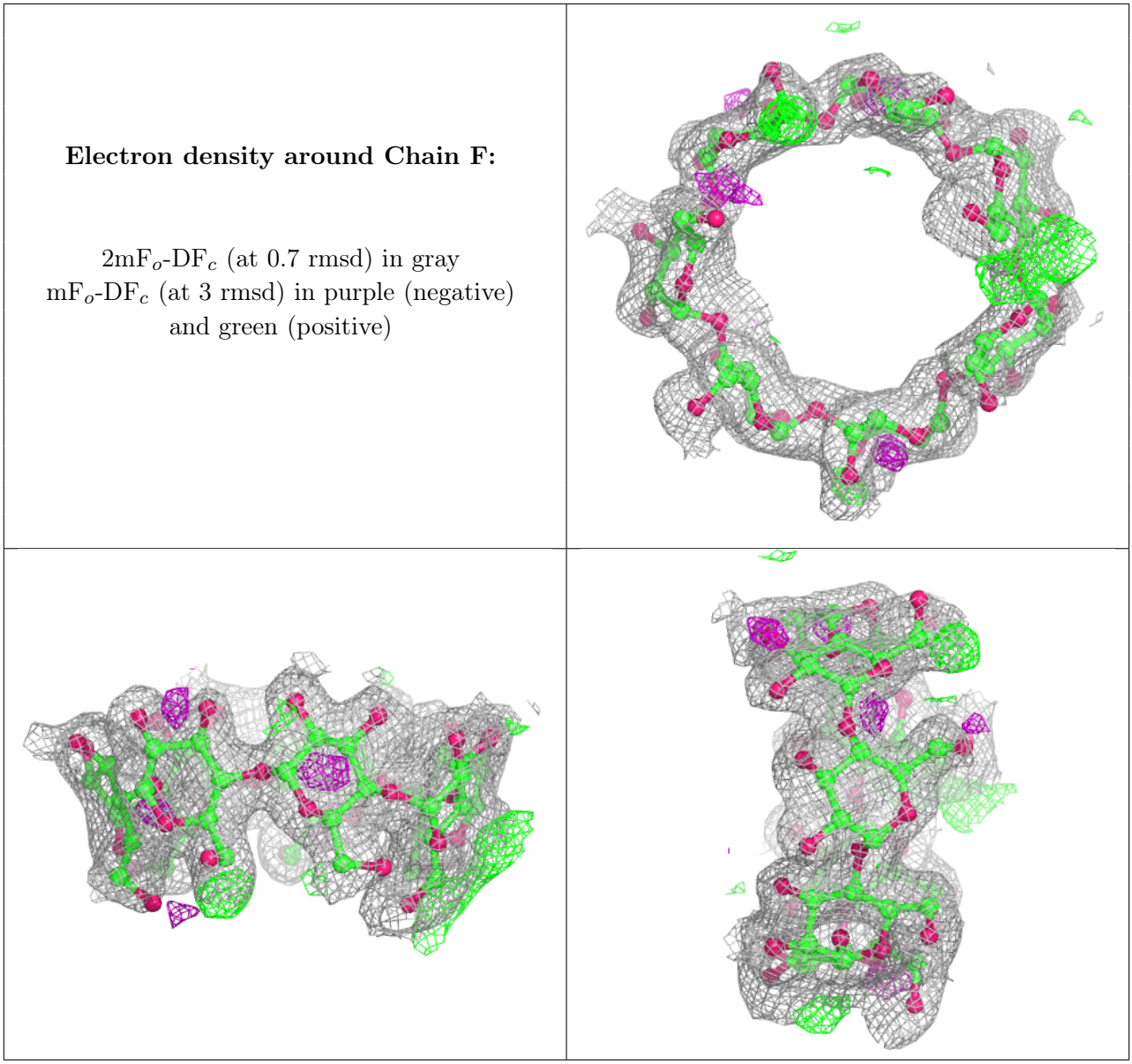






Electron density around Chain F:

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



6.4 Ligands

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, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	C8E	A	1328	13/21	0.71	0.20	47,52,58,59	0
4	C8E	B	1332	13/21	0.71	0.19	39,51,69,70	3
4	C8E	A	1331	13/21	0.74	0.18	57,72,84,88	0
4	C8E	B	1335	6/21	0.76	0.24	52,56,59,60	0
4	C8E	B	1329	14/21	0.78	0.21	47,51,55,57	2

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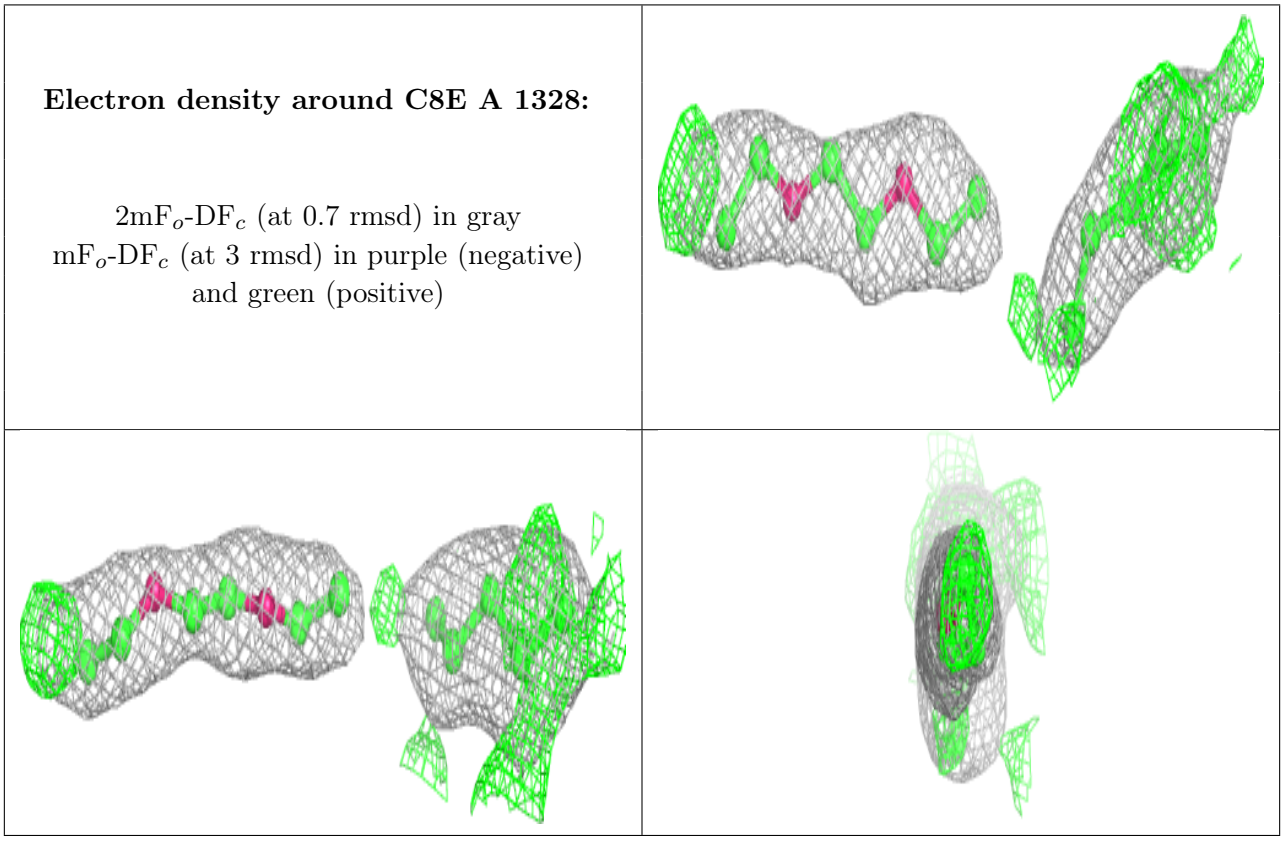
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	C8E	A	1329	13/21	0.79	0.23	50,65,80,82	0
4	C8E	B	1331	8/21	0.79	0.20	58,66,66,68	0
4	C8E	B	1328	13/21	0.80	0.23	43,55,69,73	0
4	C8E	A	1332	14/21	0.80	0.20	43,51,64,64	0
4	C8E	B	1330	21/21	0.80	0.20	50,59,67,72	0
4	C8E	B	1333	8/21	0.81	0.19	38,49,53,53	0
4	C8E	A	1334	16/21	0.81	0.18	46,56,66,69	2
4	C8E	B	1334	12/21	0.82	0.14	40,56,60,64	0
4	C8E	A	1336	7/21	0.82	0.19	59,65,68,70	0
4	C8E	A	1327	14/21	0.85	0.14	36,44,61,63	0
4	C8E	B	1336	12/21	0.86	0.13	41,48,57,60	4
4	C8E	A	1333	18/21	0.87	0.21	43,50,61,62	3
4	C8E	B	1327	13/21	0.90	0.14	43,51,56,57	0
4	C8E	A	1330	11/21	0.91	0.13	54,61,65,68	0
4	C8E	A	1335	8/21	0.94	0.17	42,46,55,55	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.

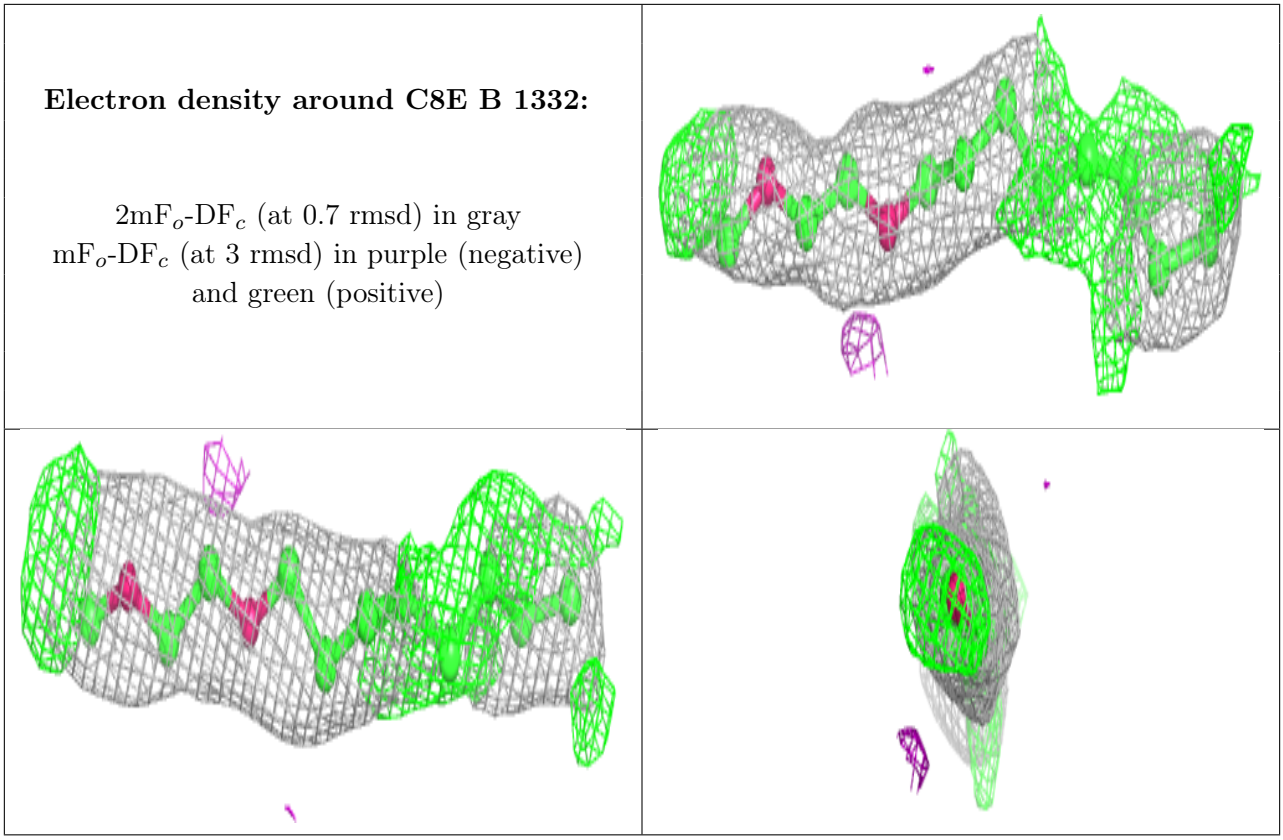
Electron density around C8E A 1328:

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



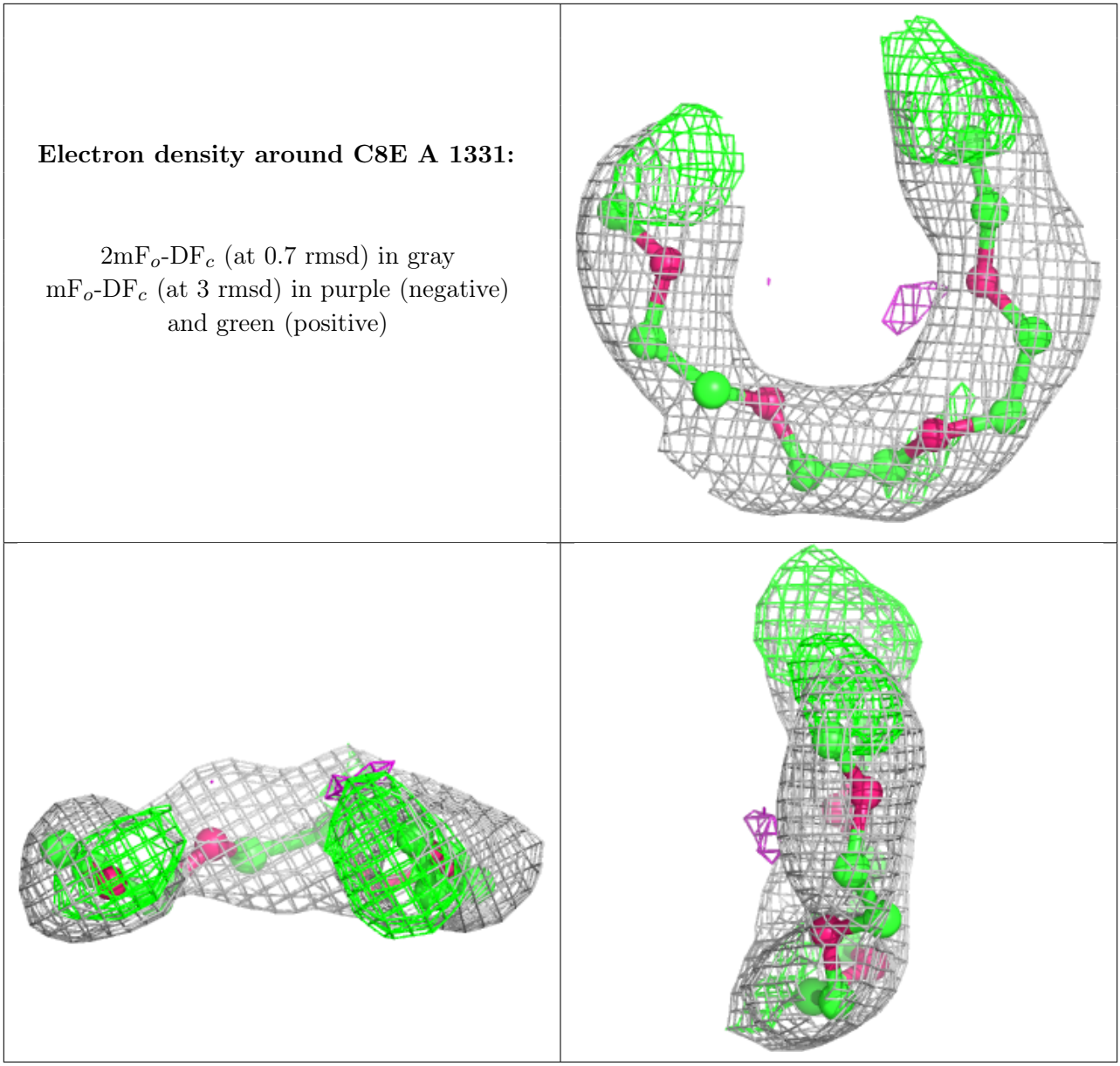
Electron density around C8E B 1332:

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



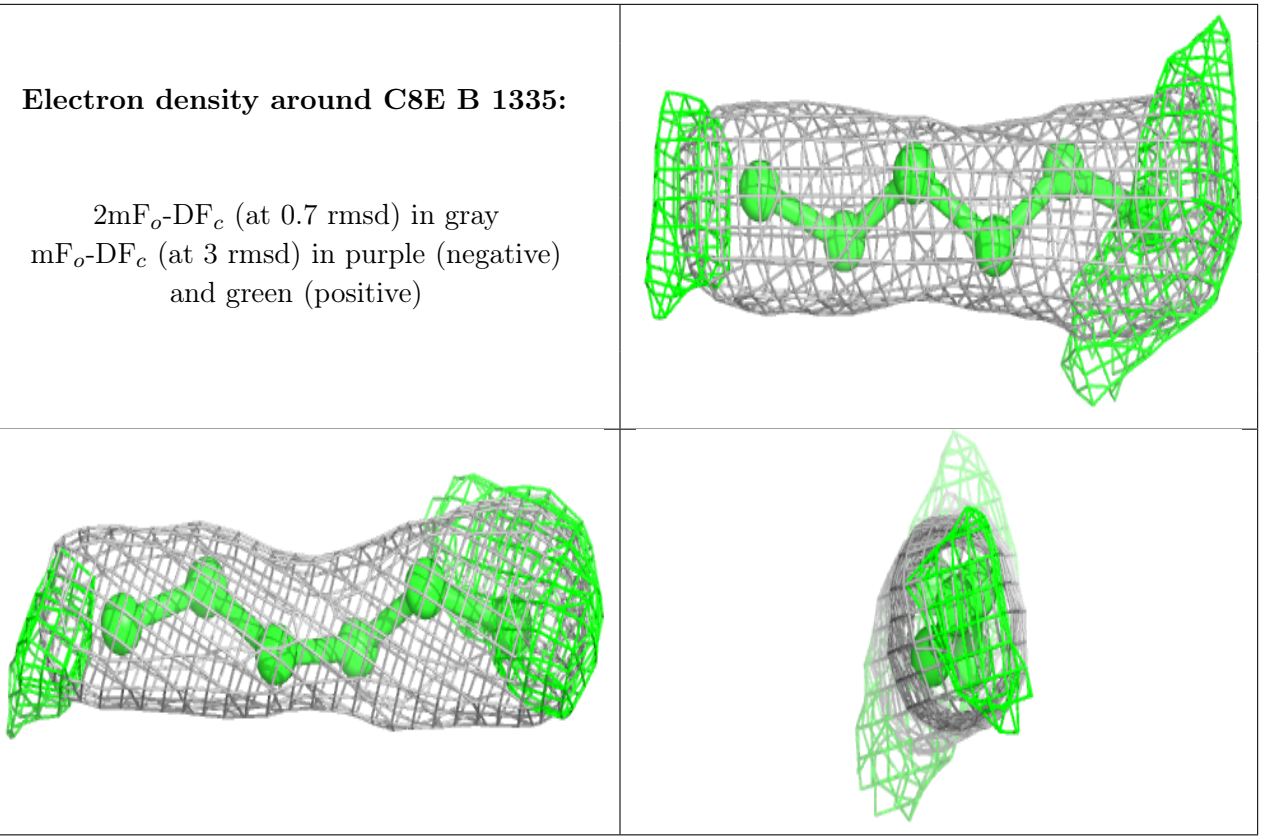
Electron density around C8E A 1331:

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



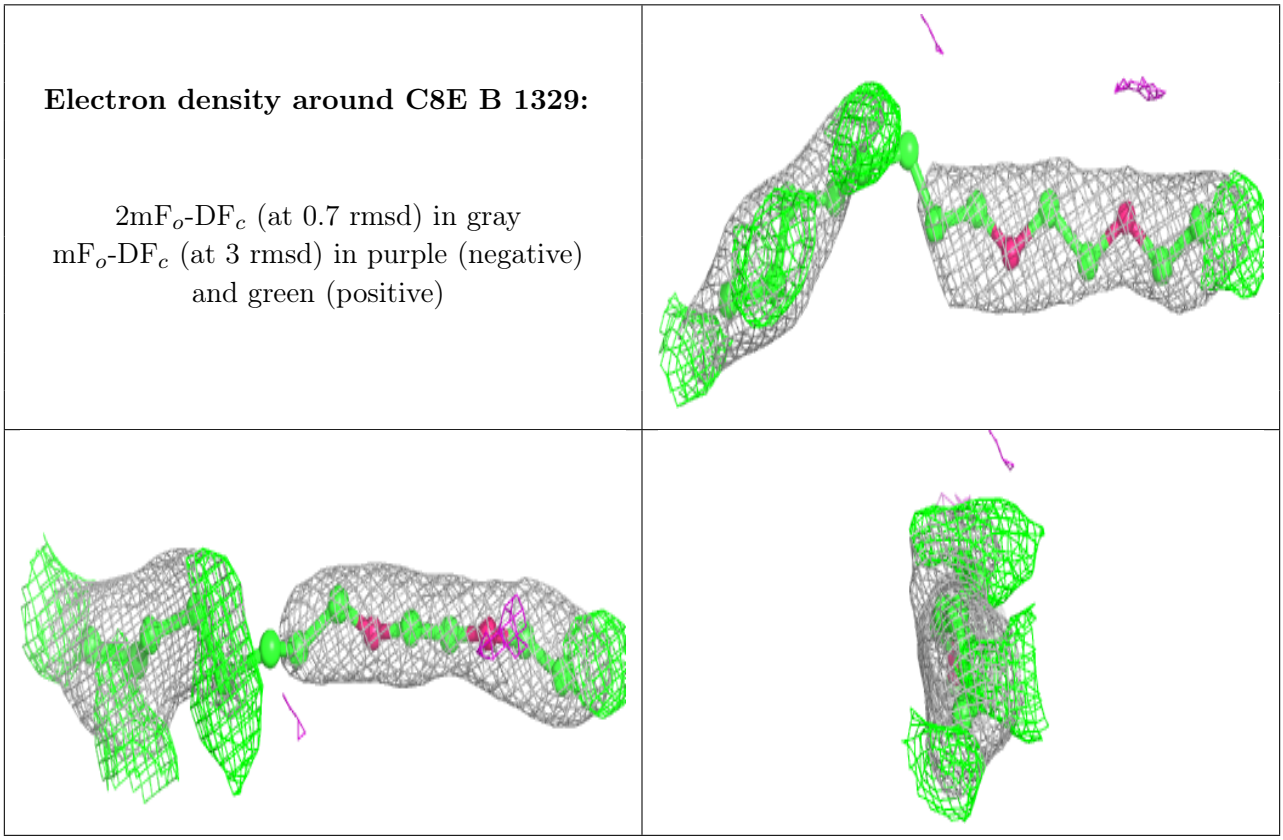
Electron density around C8E B 1335:

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



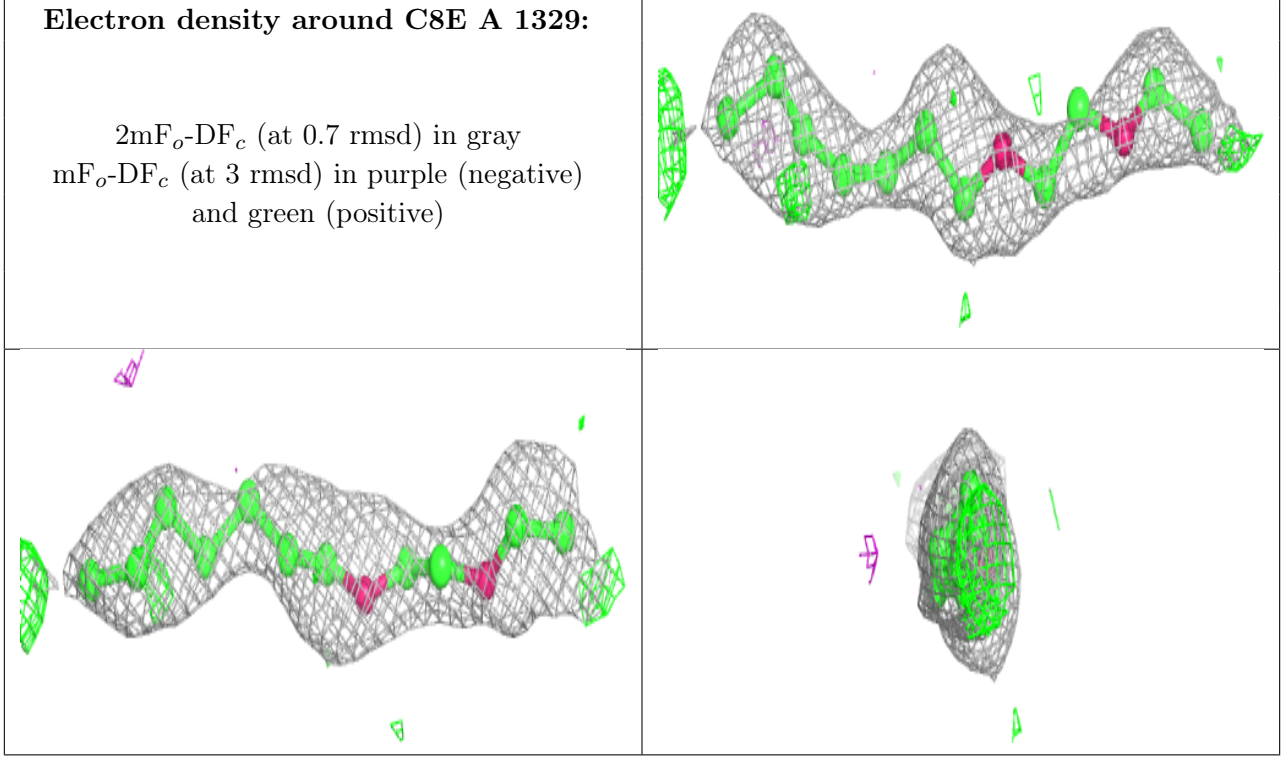
Electron density around C8E B 1329:

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



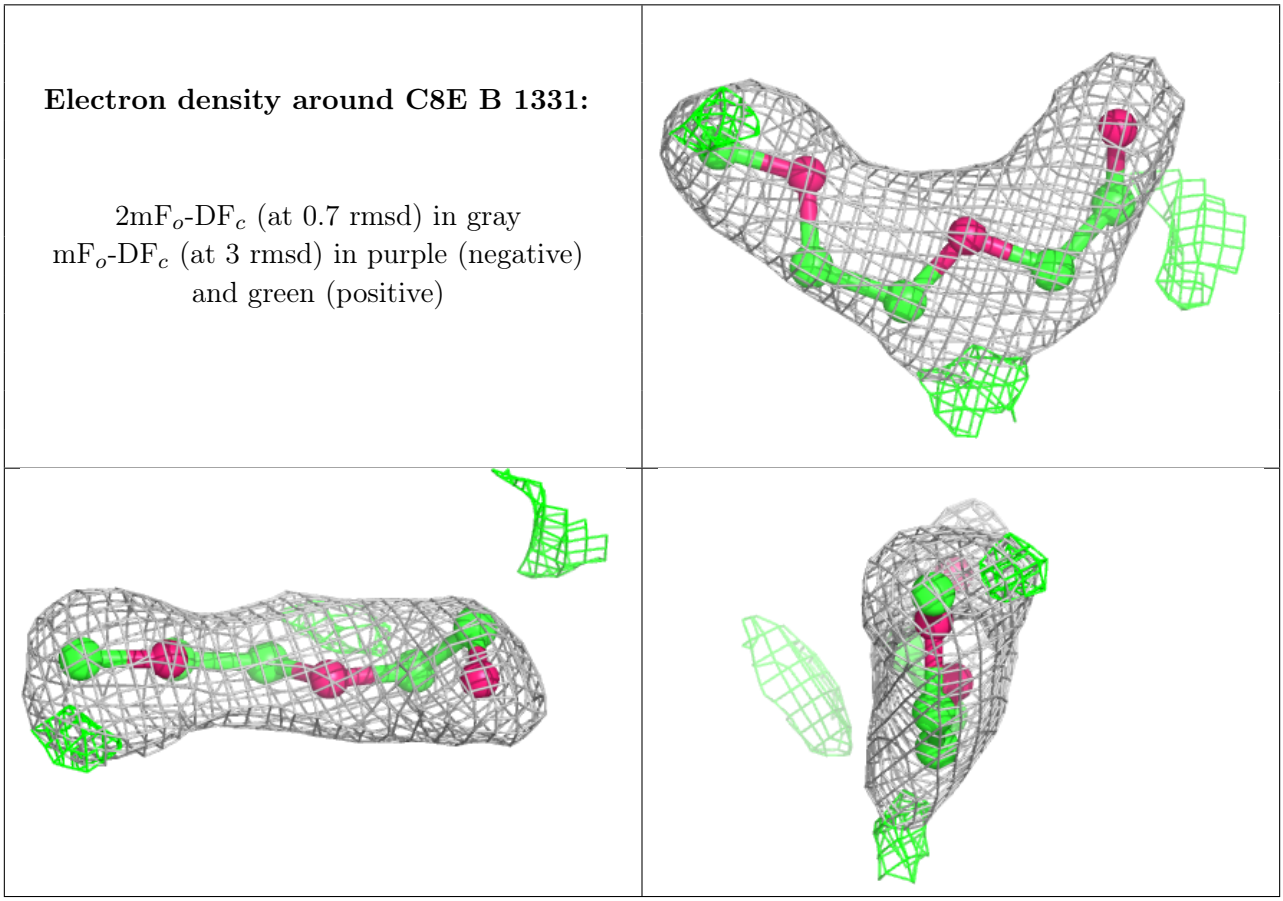
Electron density around C8E A 1329:

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



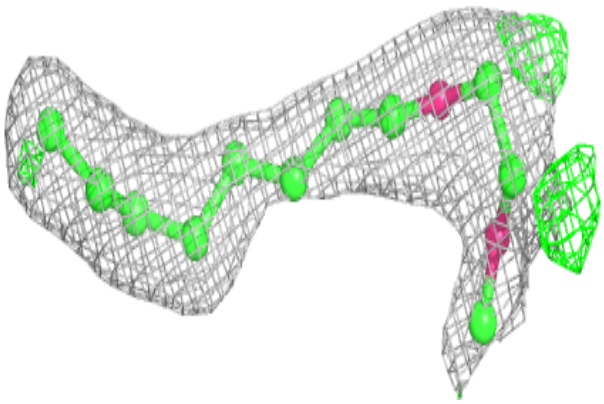
Electron density around C8E B 1331:

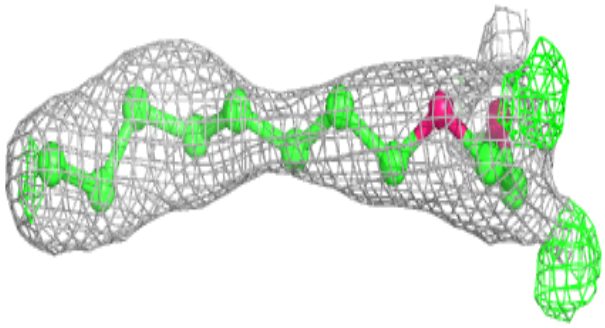
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

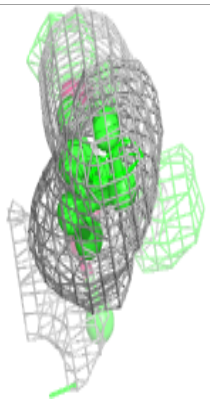


Electron density around C8E B 1328:

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

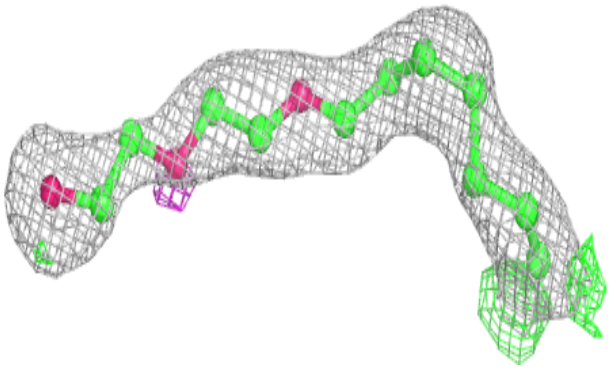


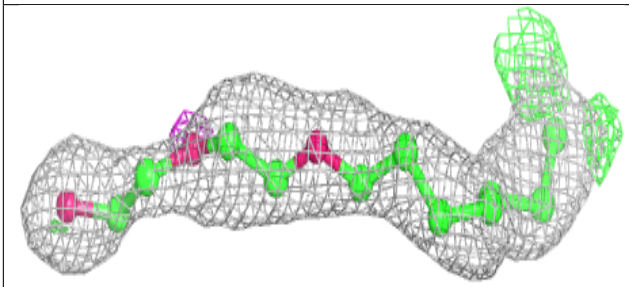


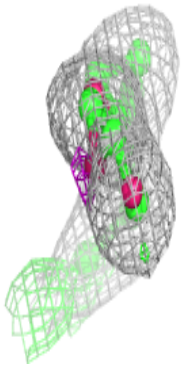


Electron density around C8E A 1332:

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

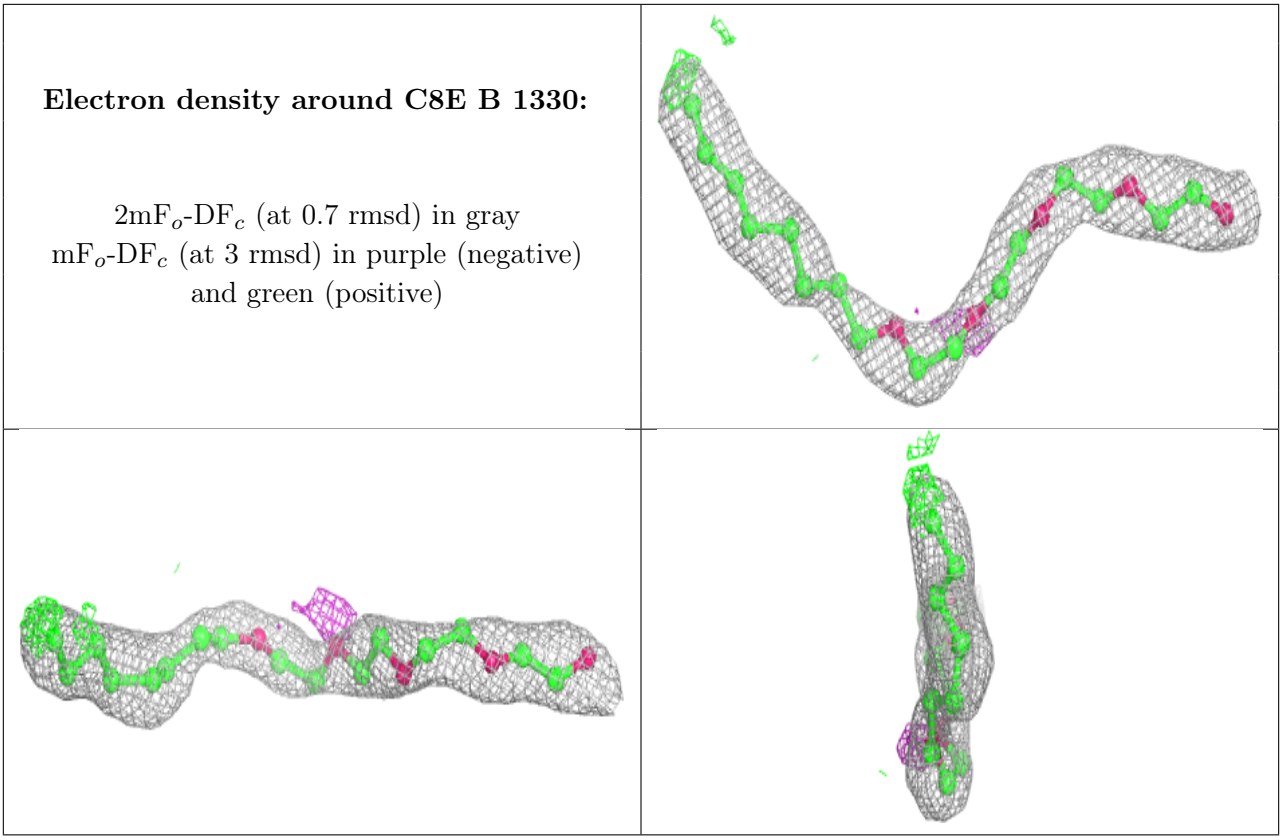






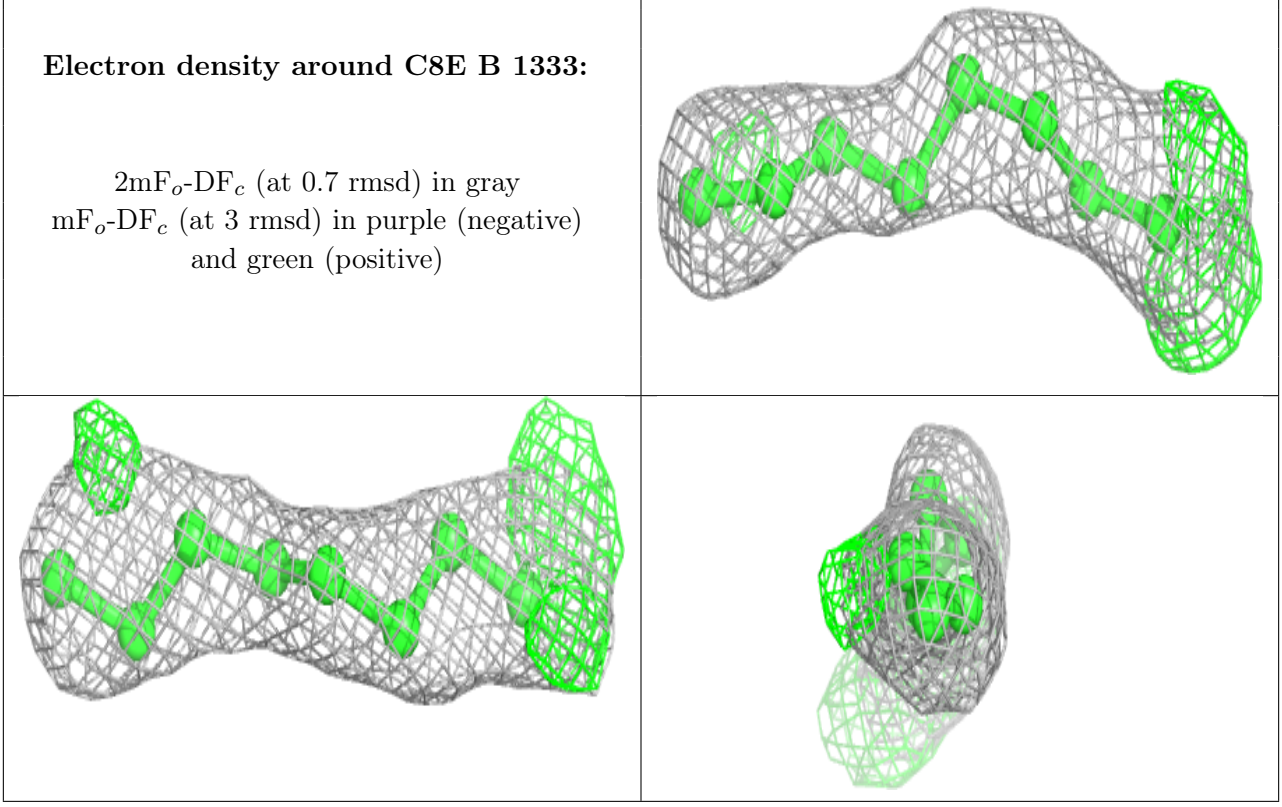
Electron density around C8E B 1330:

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



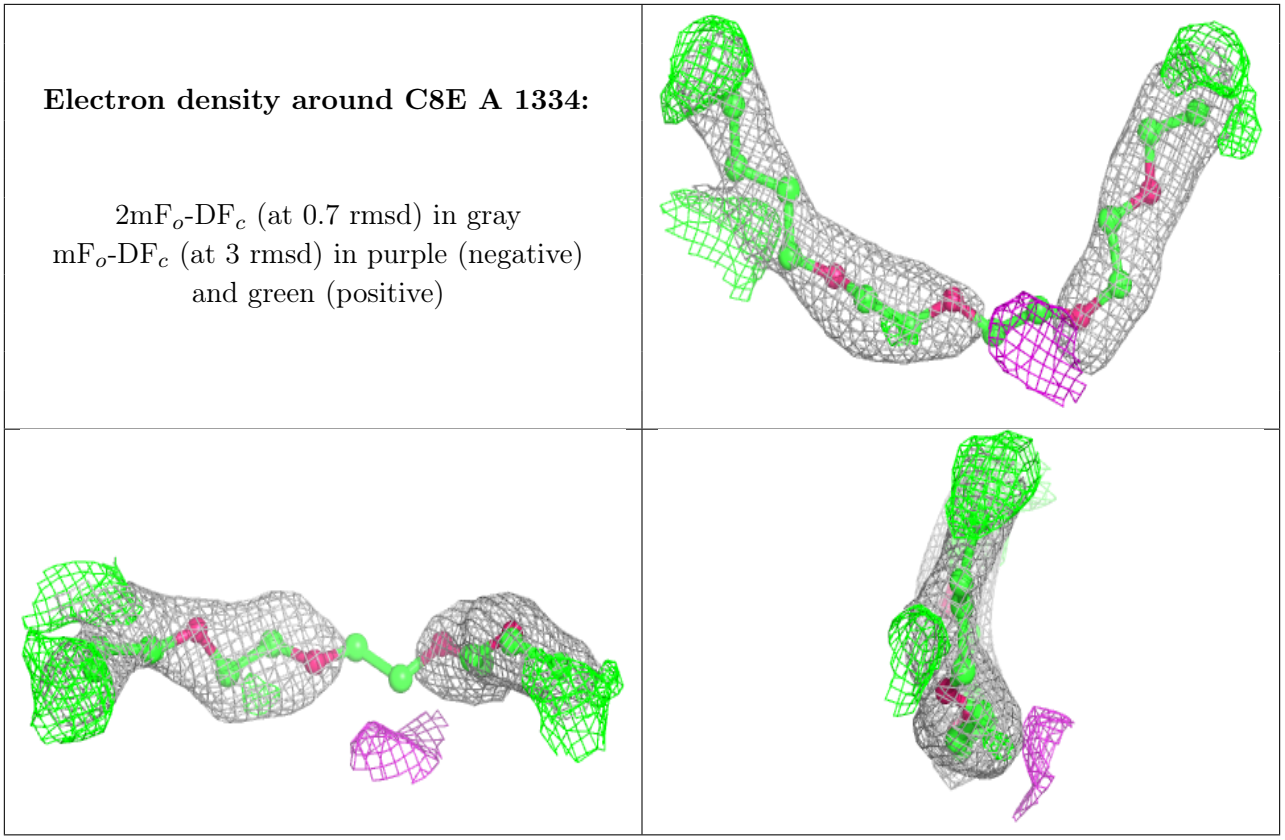
Electron density around C8E B 1333:

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



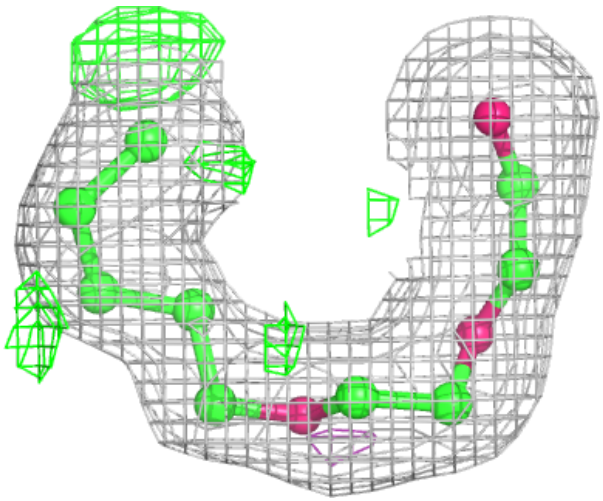
Electron density around C8E A 1334:

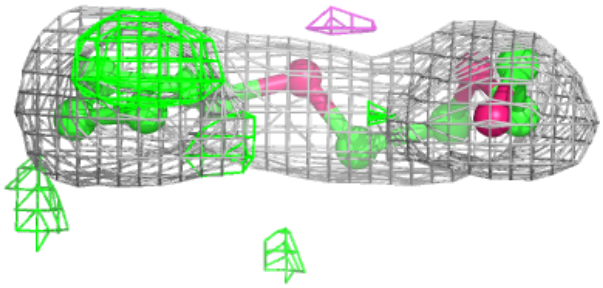
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

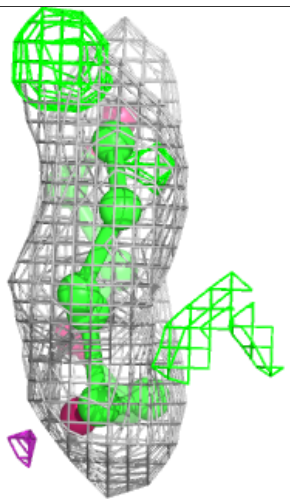


Electron density around C8E B 1334:

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

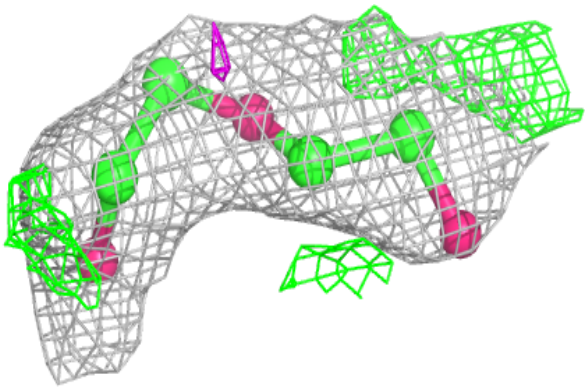


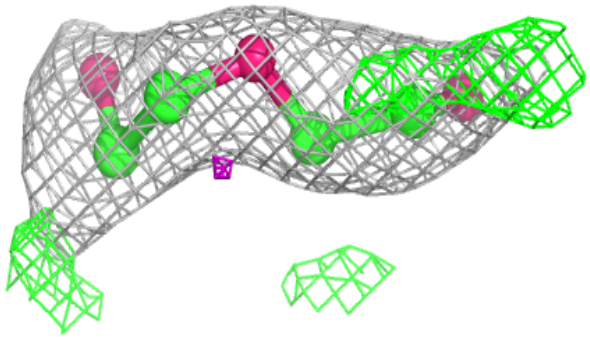


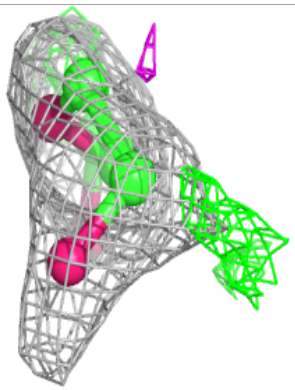


Electron density around C8E A 1336:

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

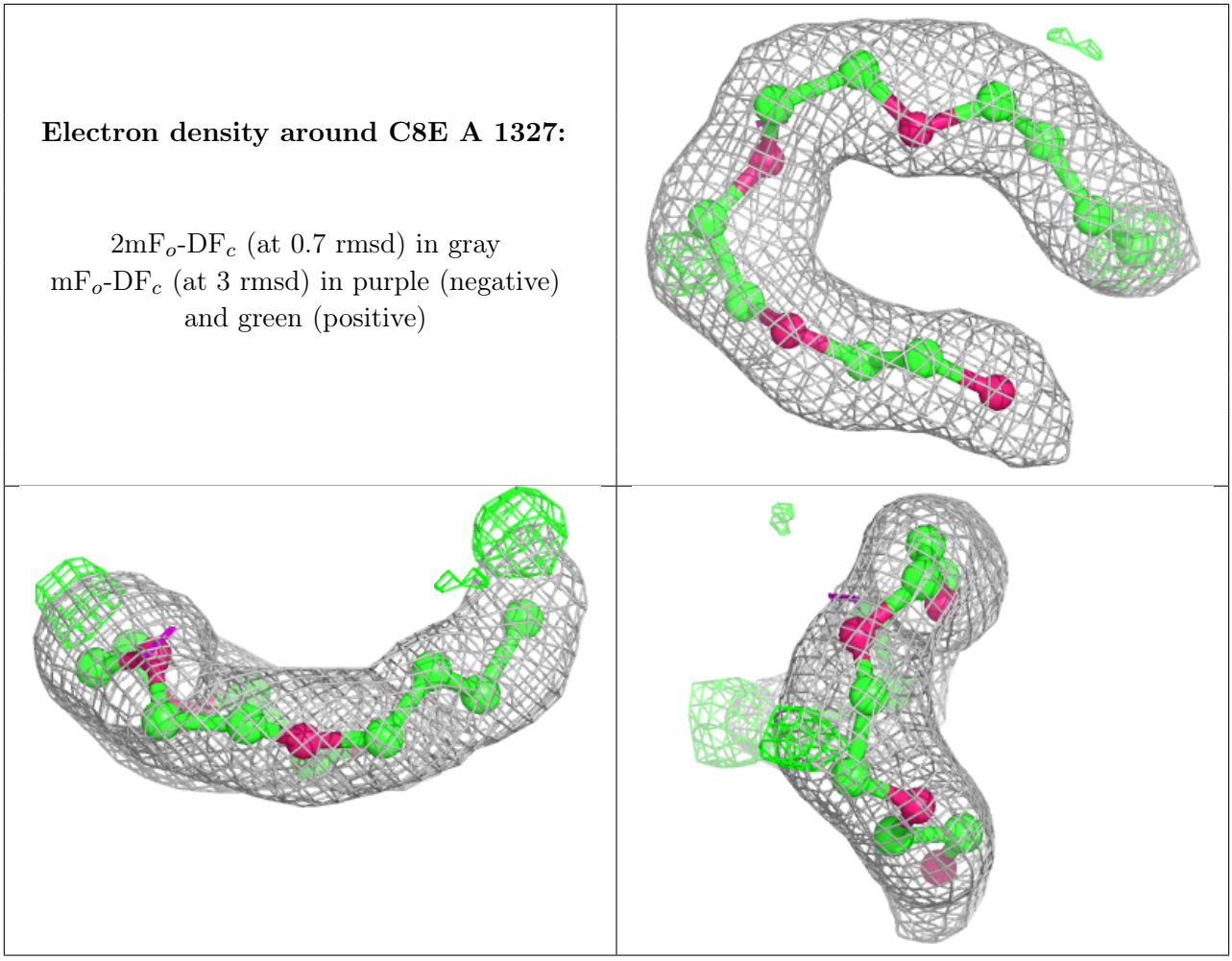






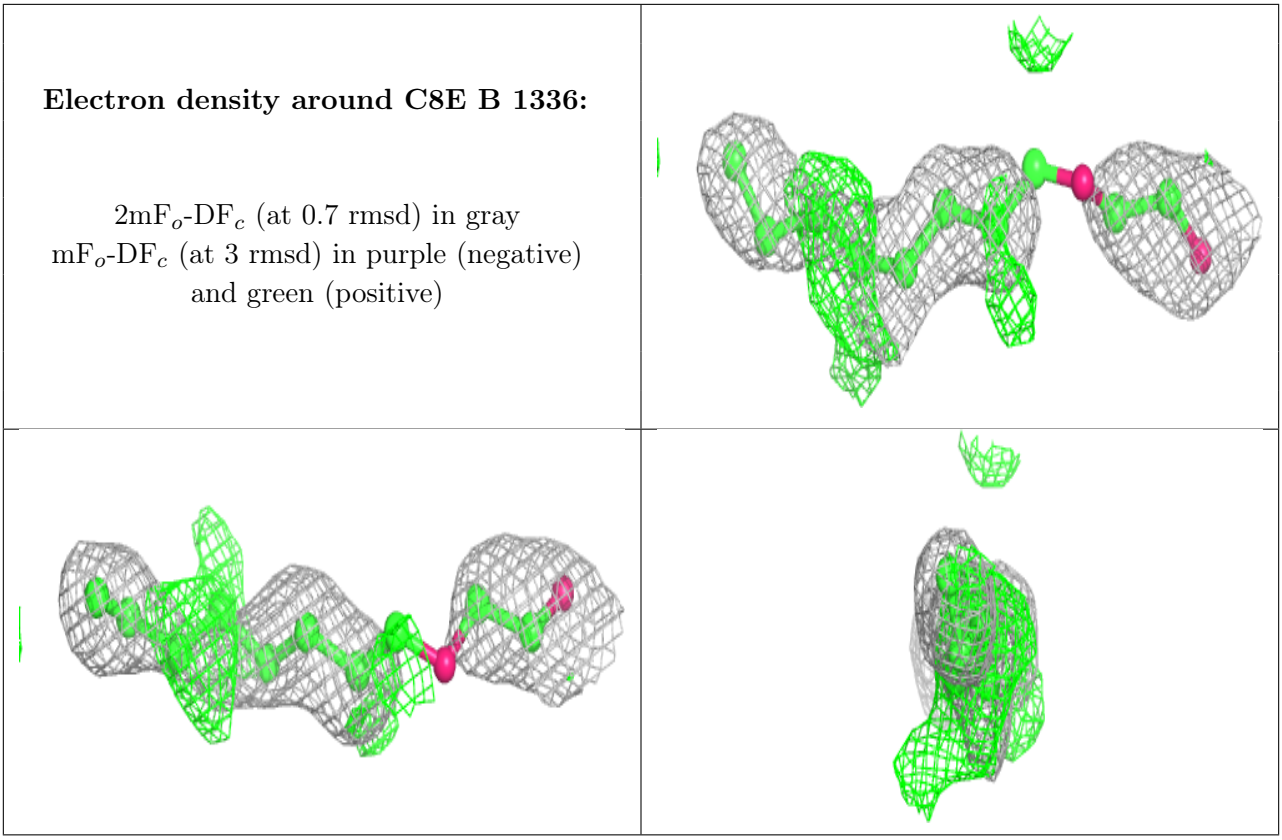
Electron density around C8E A 1327:

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



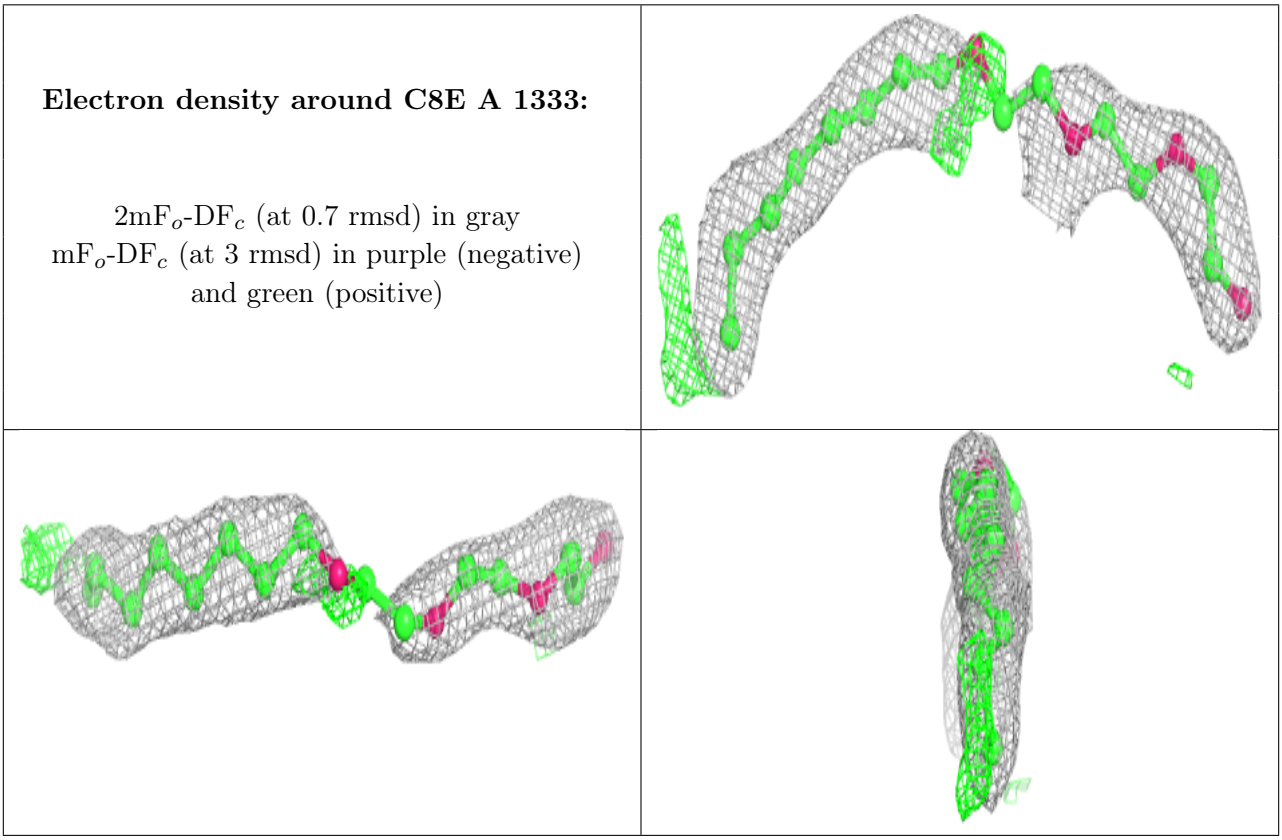
Electron density around C8E B 1336:

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



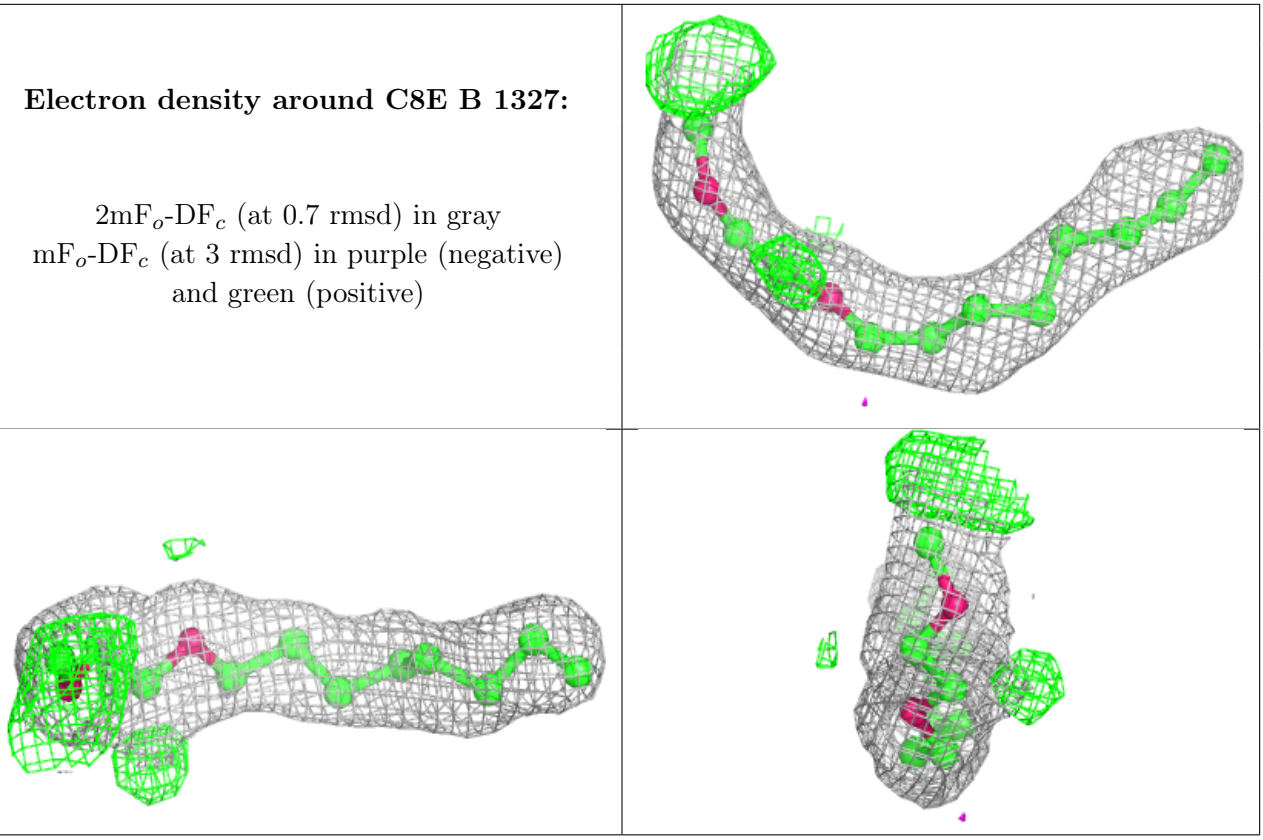
Electron density around C8E A 1333:

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



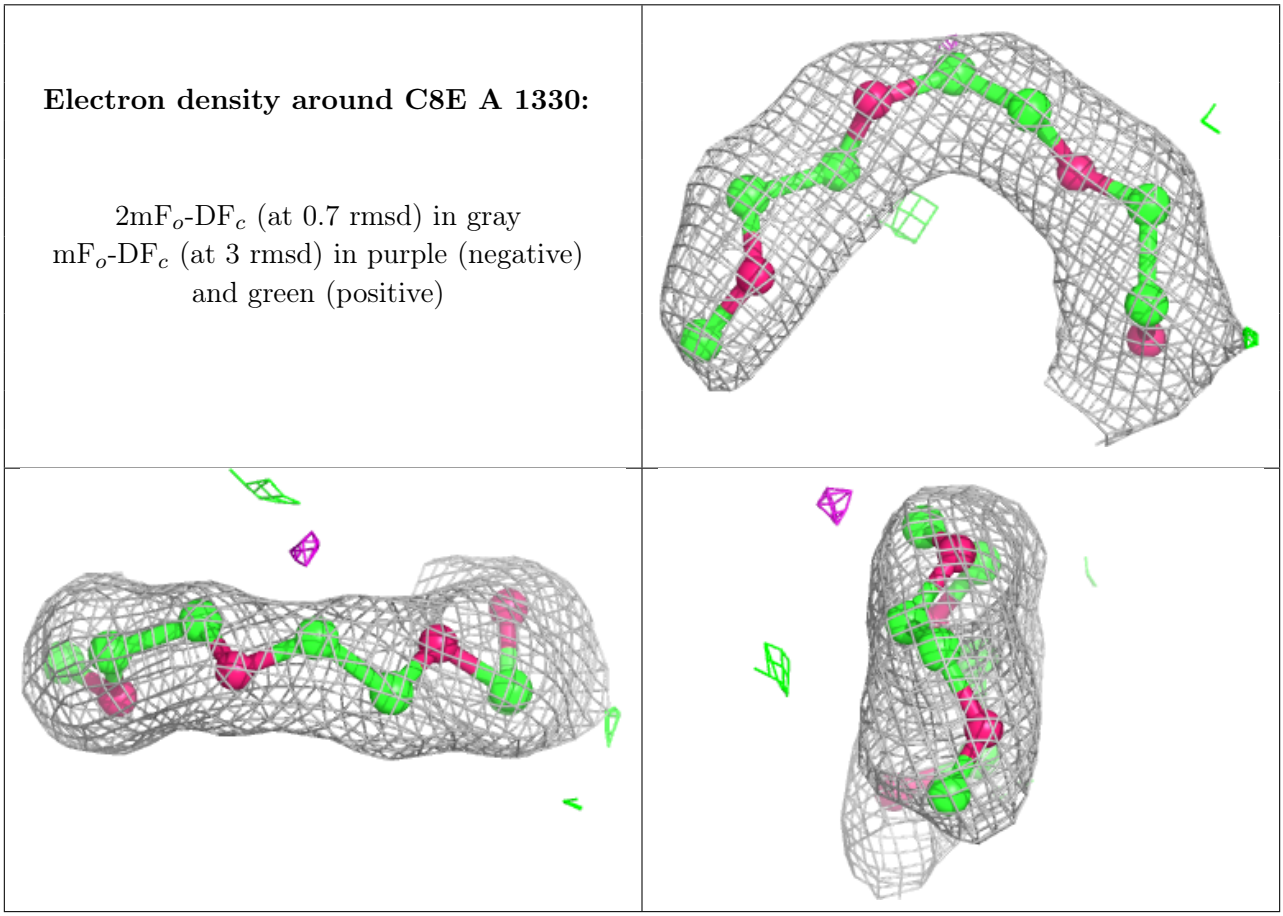
Electron density around C8E B 1327:

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



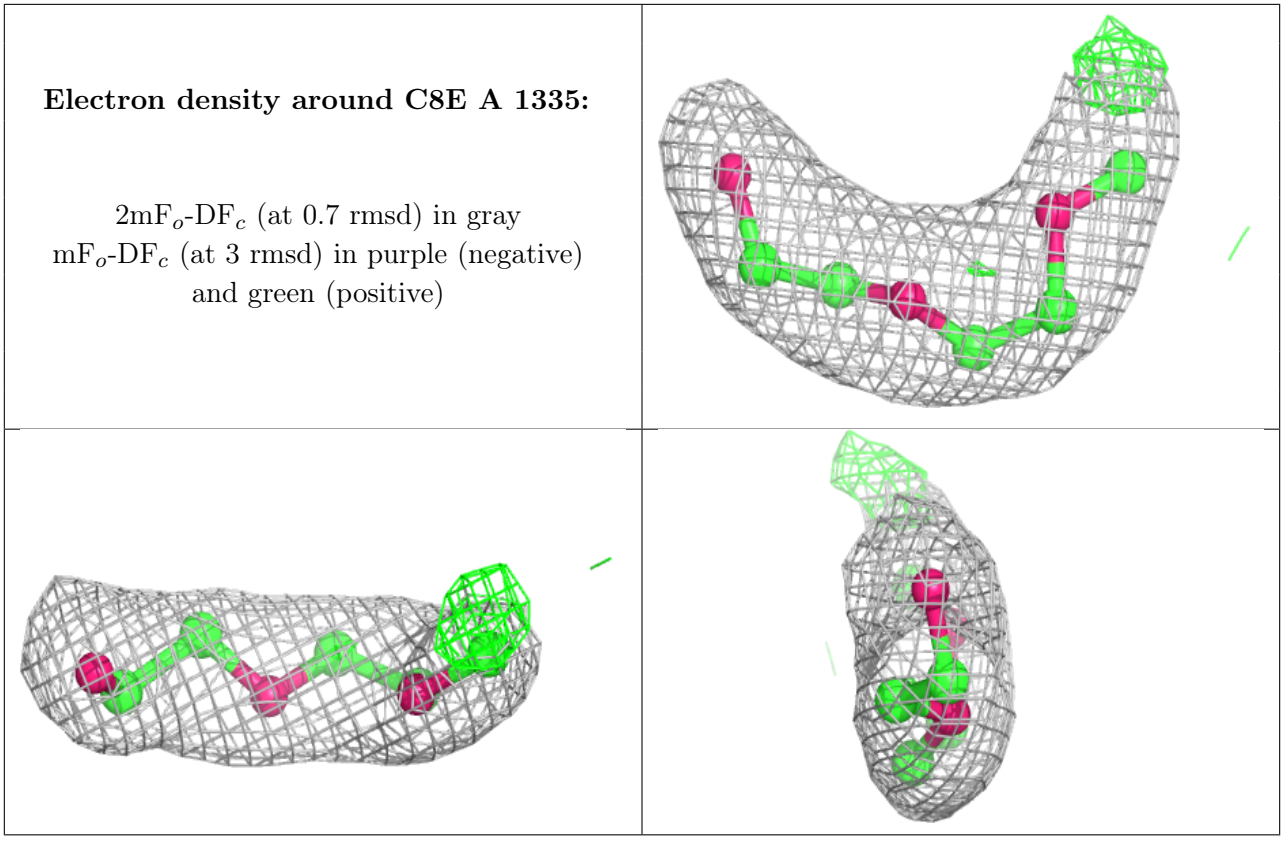
Electron density around C8E A 1330:

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



Electron density around C8E A 1335:

$2mF_o$-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.