



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 10, 2021 – 04:37 PM EDT

PDB ID : 3G84
Title : Crystal structure of the trimeric neck and carbohydrate recognition domain of R343V mutant of human surfactant protein D in complex with alpha 1,2 dimannose.
Authors : Head, J.F.
Deposited on : 2009-02-11
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.23.2
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) : 2.23.2

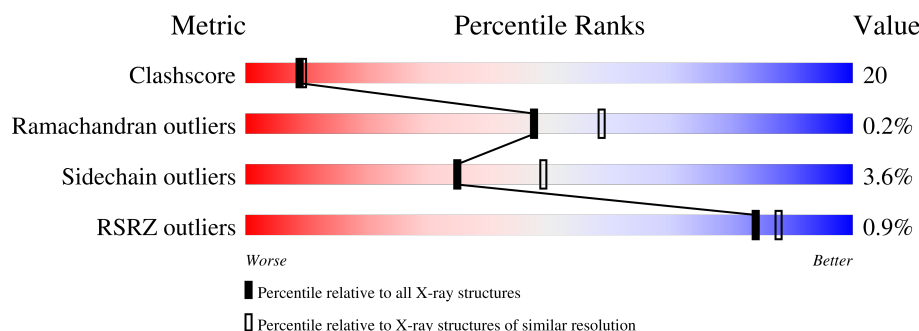
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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	160	<div> <div> <div></div> <div>58%</div> <div>36%</div> <div>• 6%</div> </div> </div>
1	B	160	<div> <div> <div></div> <div>52%</div> <div>41%</div> <div>• 6%</div> </div> </div>
1	C	160	<div> <div> <div></div> <div>56%</div> <div>38%</div> <div>• 6%</div> </div> </div>
2	D	2	<div> <div>100%</div> </div>
2	E	2	<div> <div>100%</div> </div>
2	F	2	<div> <div>50%</div> <div>50%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3657 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 Pulmonary surfactant-associated protein D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	151	Total	C	N	O	S	0	0	0
			1150	722	194	229	5			
1	B	151	Total	C	N	O	S	0	0	0
			1150	722	194	229	5			
1	C	151	Total	C	N	O	S	1	0	0
			1150	722	194	229	5			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	196	ALA	-	expression tag	UNP P35247
A	197	MET	-	expression tag	UNP P35247
A	198	ALA	-	expression tag	UNP P35247
A	199	ASP	-	expression tag	UNP P35247
A	200	ILE	-	expression tag	UNP P35247
A	201	GLY	-	expression tag	UNP P35247
A	202	SER	-	expression tag	UNP P35247
A	343	VAL	ARG	engineered mutation	UNP P35247
B	196	ALA	-	expression tag	UNP P35247
B	197	MET	-	expression tag	UNP P35247
B	198	ALA	-	expression tag	UNP P35247
B	199	ASP	-	expression tag	UNP P35247
B	200	ILE	-	expression tag	UNP P35247
B	201	GLY	-	expression tag	UNP P35247
B	202	SER	-	expression tag	UNP P35247
B	343	VAL	ARG	engineered mutation	UNP P35247
C	196	ALA	-	expression tag	UNP P35247
C	197	MET	-	expression tag	UNP P35247
C	198	ALA	-	expression tag	UNP P35247
C	199	ASP	-	expression tag	UNP P35247
C	200	ILE	-	expression tag	UNP P35247
C	201	GLY	-	expression tag	UNP P35247
C	202	SER	-	expression tag	UNP P35247

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Chain	Residue	Modelled	Actual	Comment	Reference
C	343	VAL	ARG	engineered mutation	UNP P35247

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	D	2	Total	C	O	0	0	0
			23	12	11			
2	E	2	Total	C	O	0	0	0
			23	12	11			
2	F	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Ca	0	0
			3	3		
3	B	3	Total	Ca	0	0
			3	3		
3	C	3	Total	Ca	0	0
			3	3		

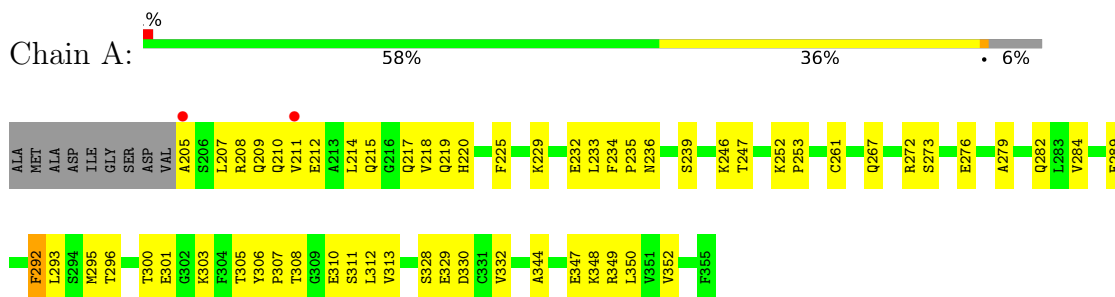
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	37	Total	O	0	0
			37	37		
4	B	46	Total	O	0	0
			46	46		
4	C	46	Total	O	0	0
			46	46		

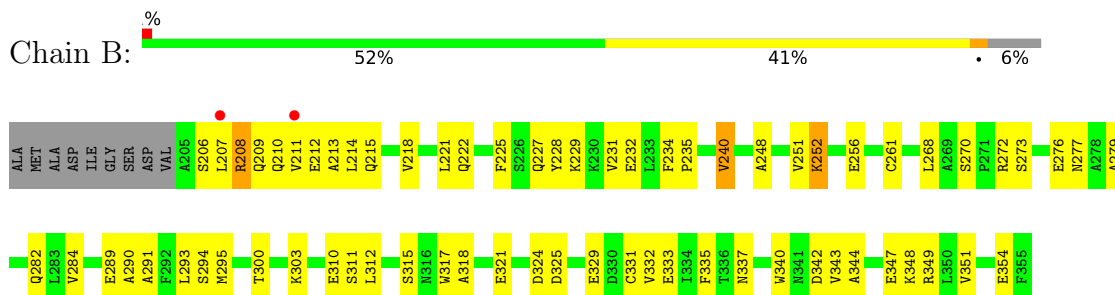
3 Residue-property plots

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.

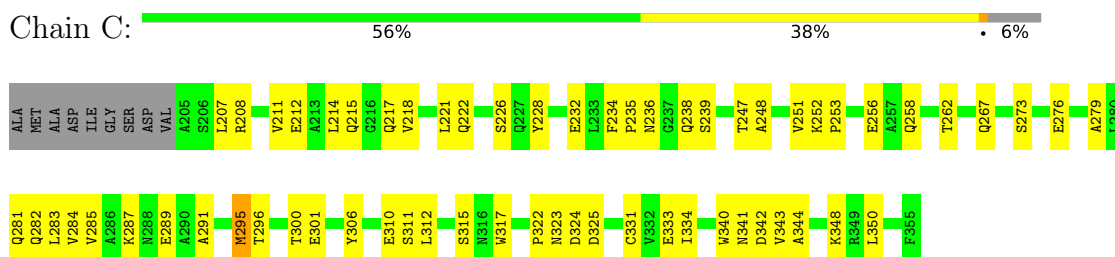
- Molecule 1: Pulmonary surfactant-associated protein D



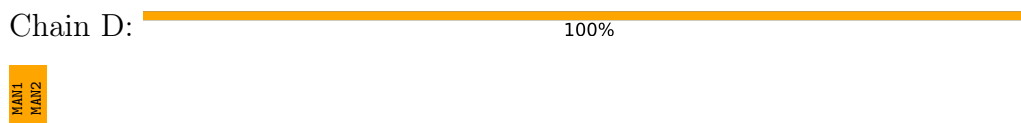
- Molecule 1: Pulmonary surfactant-associated protein D



- Molecule 1: Pulmonary surfactant-associated protein D



- Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose



- Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose

Chain E:  100%



- Molecule 2: α -D-mannopyranose-(1-2)- α -D-mannopyranose

Chain F:  50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	55.60Å 108.07Å 55.59Å 90.00° 91.47° 90.00°	Depositor
Resolution (Å)	49.42 – 2.30 49.43 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.7 (49.42-2.30) 98.6 (49.43-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.06 (at 2.29Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.152 , 0.191 0.154 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	31.4	Xtriage
Anisotropy	0.484	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 32.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.37$, $\langle L^2 \rangle = 0.20$	Xtriage
Estimated twinning fraction	0.135 for -l,k,h 0.135 for -h,-k,l 0.457 for l,-k,h	Xtriage
Reported twinning fraction	0.478 for l,-k,h	Depositor
Outliers	0 of 29053 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3657	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.33% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, 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	0.30	0/1172	0.43	0/1584
1	B	0.29	0/1172	0.45	0/1584
1	C	0.29	0/1172	0.44	0/1584
All	All	0.29	0/3516	0.44	0/4752

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1150	0	1106	55	0
1	B	1150	0	1106	48	0
1	C	1150	0	1106	51	0
2	D	23	0	21	1	0
2	E	23	0	21	1	0
2	F	23	0	21	1	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
4	A	37	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	46	0	0	0	0
4	C	46	0	0	1	0
All	All	3657	0	3381	140	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 20.

All (140) 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:A:207:LEU:HA	1:A:210:GLN:HB2	1.53	0.91
1:A:208:ARG:HH21	1:C:207:LEU:HD13	1.38	0.87
1:B:221:LEU:HD21	1:C:222:GLN:HG2	1.66	0.77
1:A:215:GLN:HE22	1:C:217:GLN:HE22	1.32	0.77
1:C:310:GLU:HG2	1:C:311:SER:H	1.49	0.77
1:A:232:GLU:OE1	1:B:232:GLU:HG3	1.84	0.77
1:A:295:MET:HG2	1:A:312:LEU:HD21	1.67	0.75
1:C:295:MET:HG2	1:C:312:LEU:HD21	1.70	0.74
2:F:1:MAN:H1	2:F:2:MAN:H5	1.71	0.73
1:B:295:MET:HG2	1:B:312:LEU:HD21	1.74	0.70
1:A:292:PHE:CD2	1:A:349:ARG:HB2	2.30	0.66
1:A:247:THR:HB	1:A:352:VAL:HG22	1.78	0.65
1:A:308:THR:OG1	1:A:310:GLU:HG2	1.97	0.64
1:A:232:GLU:HB2	1:B:232:GLU:OE1	1.98	0.63
1:B:270:SER:O	1:B:272:ARG:HG3	1.98	0.63
1:C:310:GLU:HG2	1:C:311:SER:N	2.14	0.62
1:B:248:ALA:HB3	1:B:351:VAL:HB	1.83	0.61
1:B:284:VAL:HG13	1:B:289:GLU:O	2.00	0.61
1:B:206:SER:HB2	1:B:209:GLN:OE1	2.01	0.61
1:B:225:PHE:CE2	1:B:229:LYS:HE3	2.35	0.61
1:B:344:ALA:HB3	1:B:347:GLU:HG2	1.82	0.61
1:C:214:LEU:O	1:C:218:VAL:HG23	2.01	0.60
1:A:214:LEU:O	1:A:218:VAL:HG23	2.01	0.59
1:A:300:THR:HB	1:A:303:LYS:HE2	1.86	0.57
1:C:284:VAL:HG21	1:C:334:ILE:HG23	1.86	0.57
1:A:306:TYR:CD2	1:A:310:GLU:HG3	2.40	0.56
1:B:293:LEU:HD12	1:B:332:VAL:CG1	2.35	0.56
1:B:293:LEU:HD12	1:B:332:VAL:HG12	1.88	0.55
1:B:208:ARG:NH2	1:B:209:GLN:HE21	2.06	0.54
1:C:317:TRP:CZ2	1:C:322:PRO:HG3	2.41	0.54
1:B:214:LEU:O	1:B:218:VAL:HG23	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:TYR:CE2	1:A:312:LEU:HA	2.42	0.54
1:B:300:THR:HG21	1:B:303:LYS:HD2	1.89	0.54
1:A:261:CYS:HB3	1:A:267:GLN:CA	2.37	0.54
1:B:240:VAL:HG13	1:B:240:VAL:O	2.09	0.53
1:B:324:ASP:HA	1:B:329:GLU:HB2	1.90	0.53
1:A:217:GLN:HB3	1:B:222:GLN:NE2	2.24	0.52
2:E:1:MAN:H1	2:E:2:MAN:H5	1.91	0.52
1:C:315:SER:HA	1:C:340:TRP:CZ3	2.45	0.52
1:A:211:VAL:HG11	1:C:211:VAL:HG22	1.91	0.52
1:A:261:CYS:HB3	1:A:267:GLN:HA	1.92	0.52
1:B:331:CYS:O	1:B:342:ASP:HA	2.09	0.52
1:B:317:TRP:HA	1:B:340:TRP:HB2	1.92	0.52
2:D:1:MAN:H1	2:D:2:MAN:H5	1.92	0.52
1:A:295:MET:HG2	1:A:312:LEU:CD2	2.39	0.51
1:C:291:ALA:O	1:C:333:GLU:HA	2.10	0.51
1:B:210:GLN:O	1:B:213:ALA:HB3	2.11	0.51
1:A:214:LEU:HD23	1:B:214:LEU:HB2	1.93	0.51
1:A:305:THR:HG22	1:A:311:SER:HA	1.93	0.50
1:A:208:ARG:O	1:A:212:GLU:HB2	2.11	0.50
1:B:261:CYS:SG	1:B:268:LEU:HD23	2.52	0.50
1:C:212:GLU:O	1:C:215:GLN:HB2	2.10	0.50
1:B:279:ALA:O	1:B:282:GLN:HB2	2.11	0.50
1:B:318:ALA:HB3	1:B:321:GLU:HG3	1.93	0.50
1:B:315:SER:HA	1:B:340:TRP:CZ3	2.47	0.49
1:A:215:GLN:O	1:A:219:GLN:HG2	2.12	0.49
1:A:306:TYR:HD2	1:A:310:GLU:HG3	1.76	0.49
1:C:253:PRO:HD2	1:C:256:GLU:OE1	2.12	0.49
1:C:251:VAL:HG23	1:C:348:LYS:HB3	1.94	0.49
1:A:208:ARG:HE	1:C:207:LEU:HD22	1.77	0.49
1:A:205:ALA:HA	1:A:208:ARG:HB2	1.95	0.49
1:A:273:SER:OG	1:A:276:GLU:HG3	2.13	0.49
1:A:293:LEU:HD12	1:A:332:VAL:HG11	1.93	0.49
1:B:256:GLU:OE1	1:B:256:GLU:N	2.46	0.49
1:A:217:GLN:HB3	1:B:222:GLN:HE22	1.77	0.49
1:B:235:PRO:HG3	1:C:248:ALA:HA	1.96	0.48
1:C:317:TRP:HA	1:C:340:TRP:HB2	1.94	0.48
1:A:301:GLU:HA	1:A:301:GLU:OE1	2.12	0.48
1:A:236:ASN:O	1:A:246:LYS:HG3	2.14	0.48
1:A:284:VAL:HG13	1:A:289:GLU:O	2.13	0.48
1:A:300:THR:HB	1:A:303:LYS:CE	2.43	0.48
1:A:279:ALA:HA	1:A:282:GLN:HE21	1.79	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:306:TYR:OH	1:A:313:VAL:HG22	2.14	0.47
1:B:252:LYS:HG3	1:B:351:VAL:HG21	1.95	0.47
1:B:234:PHE:CZ	1:C:248:ALA:HB2	2.50	0.47
1:A:234:PHE:HA	1:A:235:PRO:HA	1.76	0.47
1:C:333:GLU:OE2	1:C:343:VAL:HG21	2.15	0.47
1:A:253:PRO:HA	1:A:347:GLU:O	2.15	0.47
1:C:236:ASN:HB3	1:C:247:THR:HG22	1.96	0.46
1:A:247:THR:CB	1:A:352:VAL:HG22	2.43	0.46
1:C:222:GLN:O	1:C:226:SER:HB2	2.15	0.46
1:A:293:LEU:HD12	1:A:332:VAL:CG1	2.44	0.46
1:C:295:MET:SD	1:C:295:MET:N	2.88	0.46
1:A:289:GLU:HB2	1:A:350:LEU:HD22	1.98	0.46
1:A:289:GLU:HB3	1:A:350:LEU:HB2	1.96	0.46
1:C:273:SER:HB3	1:C:276:GLU:HG3	1.97	0.46
1:B:214:LEU:HD22	1:C:218:VAL:HG21	1.98	0.46
1:C:251:VAL:O	1:C:252:LYS:HG2	2.16	0.46
1:C:284:VAL:HG13	1:C:289:GLU:O	2.15	0.46
1:C:333:GLU:HG2	1:C:343:VAL:CG2	2.46	0.45
1:A:344:ALA:HA	4:A:106:HOH:O	2.15	0.45
1:C:281:GLN:O	1:C:285:VAL:HG23	2.15	0.45
1:A:233:LEU:HD11	1:A:239:SER:HB2	1.97	0.45
1:B:234:PHE:HA	1:B:235:PRO:HA	1.83	0.45
1:B:268:LEU:O	1:B:294:SER:HB3	2.17	0.45
1:B:310:GLU:HG2	1:B:311:SER:N	2.32	0.45
1:C:315:SER:HA	1:C:340:TRP:CH2	2.52	0.45
1:C:267:GLN:HB2	4:C:95:HOH:O	2.17	0.45
1:B:227:GLN:O	1:B:231:VAL:HG23	2.17	0.44
1:B:300:THR:CG2	1:B:303:LYS:HD2	2.47	0.44
1:B:354:GLU:O	1:B:354:GLU:HG3	2.18	0.44
1:C:221:LEU:HD23	1:C:221:LEU:HA	1.86	0.44
1:B:228:TYR:HA	1:B:231:VAL:HB	1.99	0.44
1:A:205:ALA:O	1:A:209:GLN:HG2	2.18	0.44
1:A:234:PHE:CZ	1:B:248:ALA:HB2	2.52	0.44
1:A:247:THR:HA	1:A:352:VAL:HA	2.00	0.44
1:C:258:GLN:O	1:C:262:THR:HG23	2.18	0.44
1:C:343:VAL:HG12	1:C:344:ALA:N	2.32	0.44
1:B:291:ALA:O	1:B:333:GLU:HA	2.17	0.44
1:A:232:GLU:OE1	1:C:232:GLU:OE1	2.36	0.43
1:C:331:CYS:O	1:C:342:ASP:HA	2.18	0.43
1:C:291:ALA:HB2	1:C:350:LEU:HB3	2.00	0.43
1:B:251:VAL:C	1:B:252:LYS:HG2	2.39	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:341:ASN:ND2	1:C:343:VAL:HG22	2.34	0.43
1:C:279:ALA:O	1:C:282:GLN:HB2	2.19	0.43
1:C:323:ASN:OD1	1:C:325:ASP:HB2	2.19	0.43
1:A:306:TYR:HB3	1:A:307:PRO:HD2	2.01	0.42
1:C:208:ARG:O	1:C:212:GLU:HB2	2.20	0.42
1:A:225:PHE:CZ	1:A:229:LYS:HD3	2.54	0.42
1:C:341:ASN:O	1:C:343:VAL:HG23	2.19	0.42
1:B:290:ALA:O	1:B:349:ARG:HB3	2.18	0.42
1:C:301:GLU:OE1	1:C:301:GLU:HA	2.20	0.42
1:A:229:LYS:NZ	1:C:228:TYR:HB2	2.35	0.42
1:A:296:THR:HA	1:A:330:ASP:O	2.19	0.41
1:A:328:SER:O	1:A:329:GLU:HG3	2.20	0.41
1:B:273:SER:OG	1:B:276:GLU:HB2	2.20	0.41
1:A:252:LYS:O	1:A:348:LYS:HA	2.20	0.41
1:C:287:LYS:HA	1:C:287:LYS:HD3	1.89	0.41
1:B:335:PHE:HB2	1:B:337:ASN:OD1	2.21	0.41
1:C:295:MET:CE	1:C:306:TYR:HE1	2.34	0.41
1:B:211:VAL:O	1:B:215:GLN:HG3	2.21	0.41
1:C:238:GLN:HG3	1:C:279:ALA:O	2.21	0.41
1:C:234:PHE:HA	1:C:235:PRO:HA	1.80	0.40
1:A:235:PRO:HG3	1:B:248:ALA:HA	2.03	0.40
1:C:238:GLN:HG2	1:C:283:LEU:HG	2.03	0.40
1:C:343:VAL:HG12	1:C:344:ALA:H	1.86	0.40
1:B:207:LEU:HA	1:B:210:GLN:OE1	2.21	0.40
1:A:205:ALA:N	1:A:208:ARG:HG3	2.37	0.40
1:A:272:ARG:NH1	1:A:313:VAL:HG21	2.36	0.40
1:C:324:ASP:O	1:C:325:ASP:C	2.60	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	149/160 (93%)	141 (95%)	8 (5%)	0	100	100
1	B	149/160 (93%)	143 (96%)	5 (3%)	1 (1%)	22	26
1	C	149/160 (93%)	142 (95%)	7 (5%)	0	100	100
All	All	447/480 (93%)	426 (95%)	20 (4%)	1 (0%)	47	58

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	240	VAL

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	121/127 (95%)	119 (98%)	2 (2%)	60	76
1	B	121/127 (95%)	114 (94%)	7 (6%)	20	27
1	C	121/127 (95%)	117 (97%)	4 (3%)	38	53
All	All	363/381 (95%)	350 (96%)	13 (4%)	35	49

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	220	HIS
1	A	292	PHE
1	B	208	ARG
1	B	212	GLU
1	B	252	LYS
1	B	277	ASN
1	B	325	ASP
1	B	343	VAL
1	B	348	LYS
1	C	239	SER
1	C	295	MET
1	C	296	THR

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Mol	Chain	Res	Type
1	C	300	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	210	GLN
1	A	215	GLN
1	A	222	GLN
1	A	263	GLN
1	A	282	GLN
1	B	209	GLN
1	B	222	GLN
1	B	282	GLN
1	C	217	GLN
1	C	263	GLN
1	C	267	GLN
1	C	281	GLN
1	C	282	GLN

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 ⓘ

6 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MAN	D	1	2	12,12,12	0.44	0	17,17,17	1.40	1 (5%)
2	MAN	D	2	2	11,11,12	0.57	0	15,15,17	1.31	1 (6%)
2	MAN	E	1	2	12,12,12	0.54	0	17,17,17	0.88	1 (5%)
2	MAN	E	2	2	11,11,12	0.55	0	15,15,17	1.10	1 (6%)
2	MAN	F	1	2	12,12,12	0.55	0	17,17,17	0.91	0
2	MAN	F	2	2	11,11,12	0.51	0	15,15,17	1.42	1 (6%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MAN	D	1	2	-	2/2/22/22	0/1/1/1
2	MAN	D	2	2	-	2/2/19/22	0/1/1/1
2	MAN	E	1	2	-	0/2/22/22	0/1/1/1
2	MAN	E	2	2	-	0/2/19/22	0/1/1/1
2	MAN	F	1	2	-	1/2/22/22	0/1/1/1
2	MAN	F	2	2	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	2	MAN	C1-O5-C5	4.02	117.64	112.19
2	D	2	MAN	C1-O5-C5	3.83	117.38	112.19
2	D	1	MAN	O2-C2-C3	-3.31	102.70	110.35
2	E	2	MAN	C1-O5-C5	2.74	115.90	112.19
2	E	1	MAN	O2-C2-C3	-2.13	105.42	110.35

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	1	MAN	O5-C5-C6-O6
2	D	1	MAN	C4-C5-C6-O6
2	F	2	MAN	O5-C5-C6-O6
2	F	2	MAN	C4-C5-C6-O6
2	D	2	MAN	C4-C5-C6-O6

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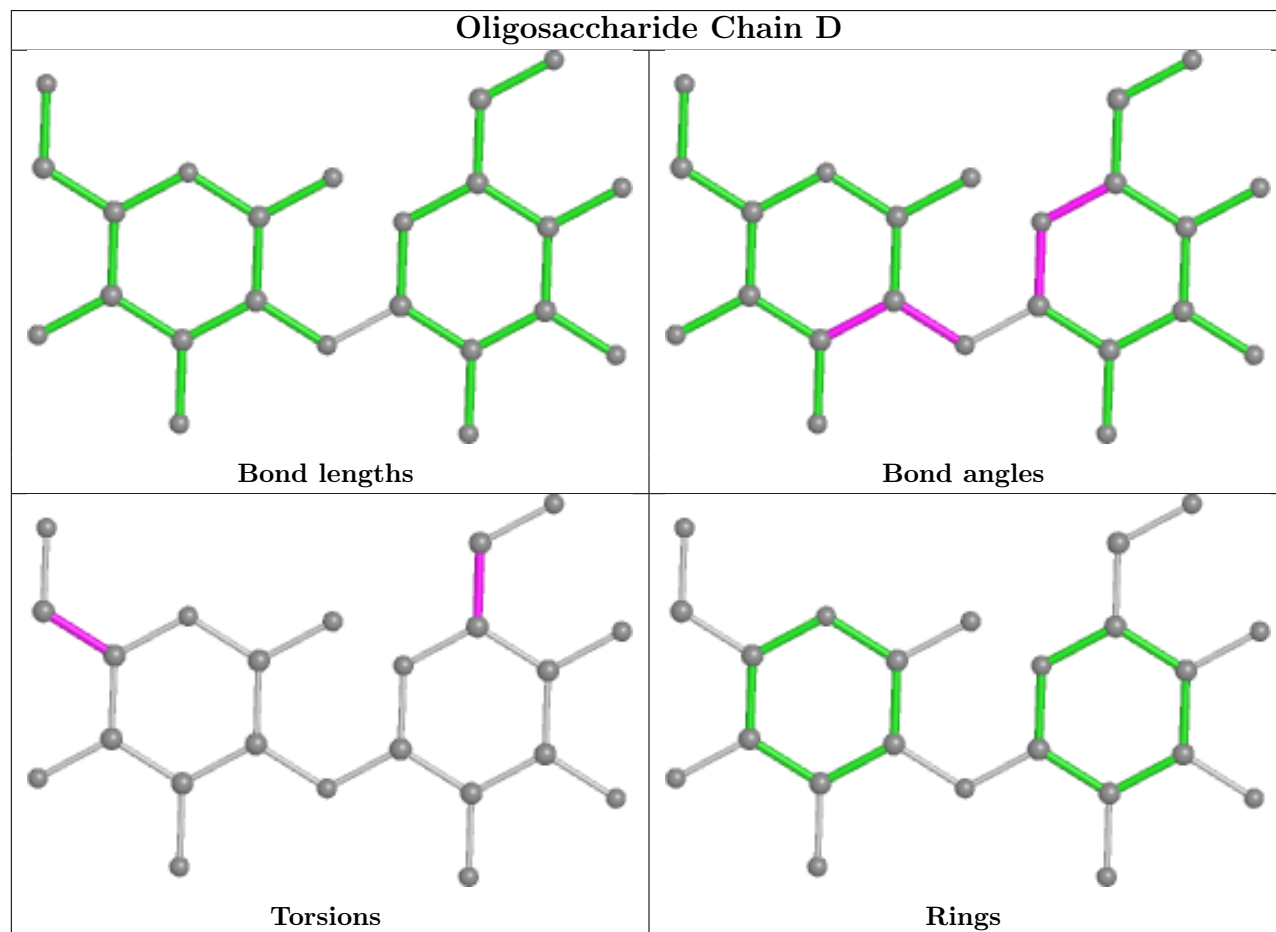
Mol	Chain	Res	Type	Atoms
2	F	1	MAN	C4-C5-C6-O6
2	D	2	MAN	O5-C5-C6-O6

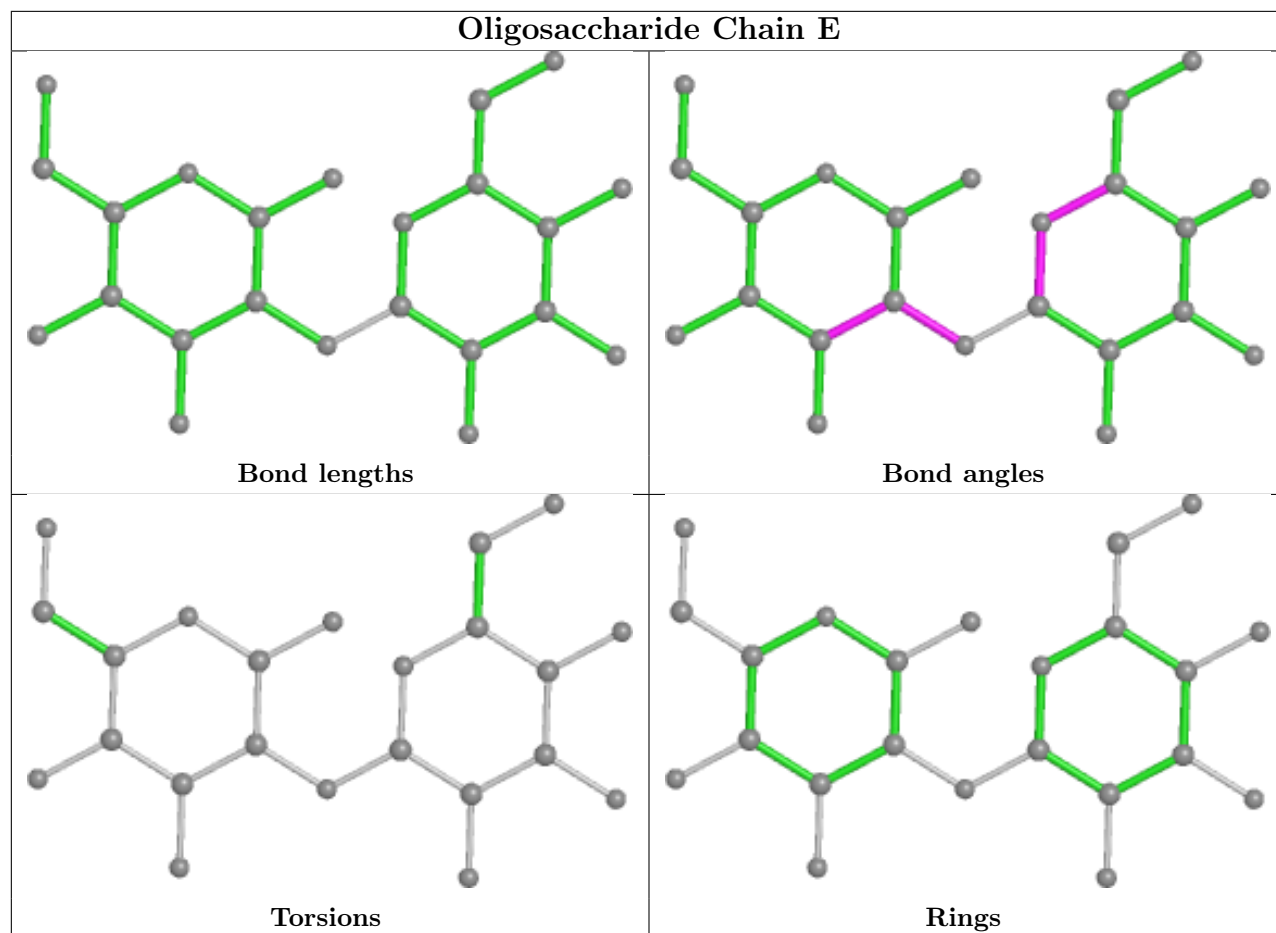
There are no ring outliers.

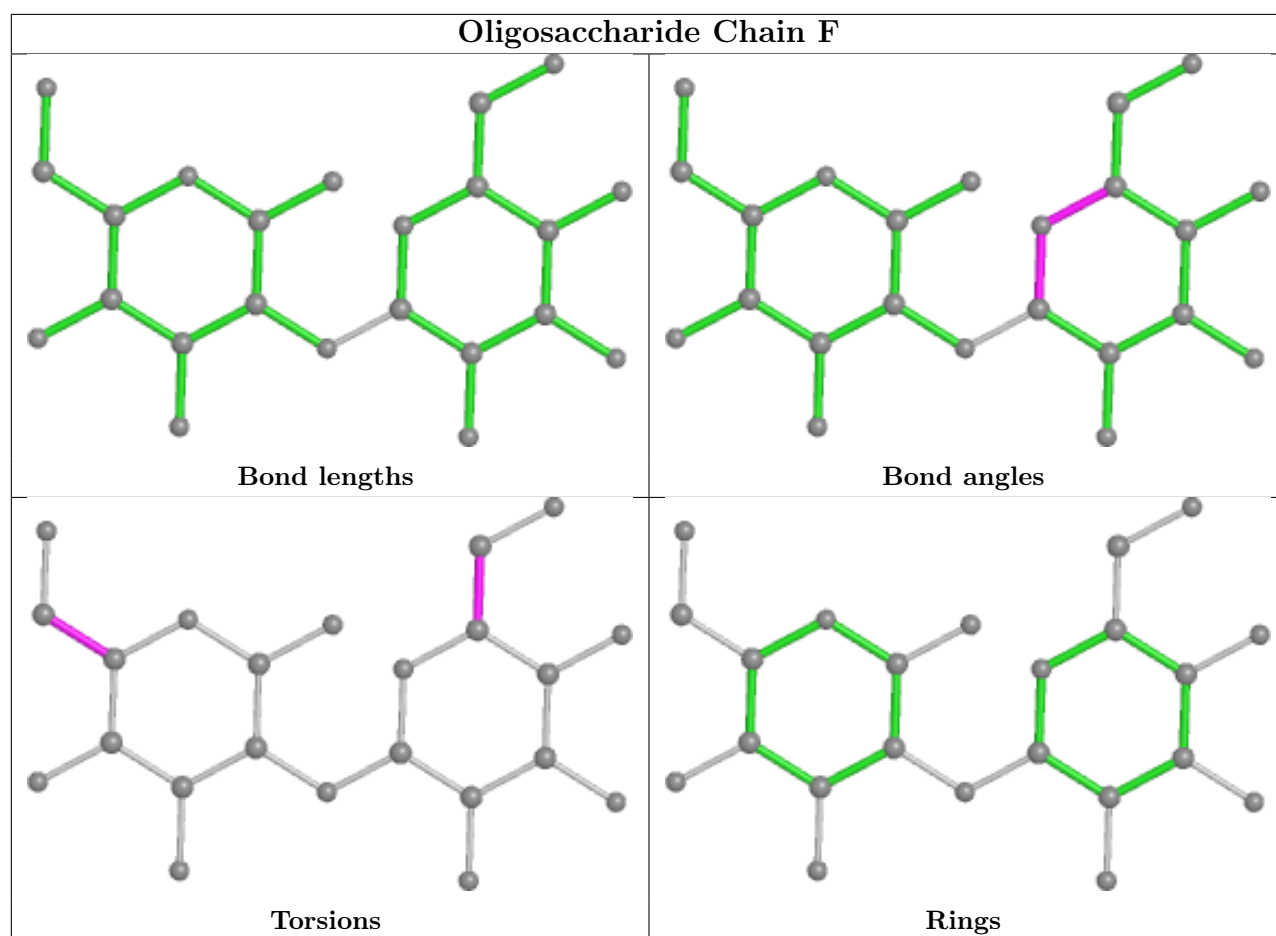
6 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	2	MAN	1	0
2	F	1	MAN	1	0
2	F	2	MAN	1	0
2	D	2	MAN	1	0
2	E	1	MAN	1	0
2	D	1	MAN	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 9 ligands modelled in this entry, 9 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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	151/160 (94%)	-0.41	2 (1%) 77 81	26, 33, 61, 82	0
1	B	151/160 (94%)	-0.47	2 (1%) 77 81	23, 31, 60, 72	0
1	C	151/160 (94%)	-0.49	0 100 100	24, 33, 57, 71	1 (0%)
All	All	453/480 (94%)	-0.46	4 (0%) 84 88	23, 32, 58, 82	1 (0%)

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	205	ALA	4.0
1	A	211	VAL	2.9
1	B	207	LEU	2.4
1	B	211	VAL	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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

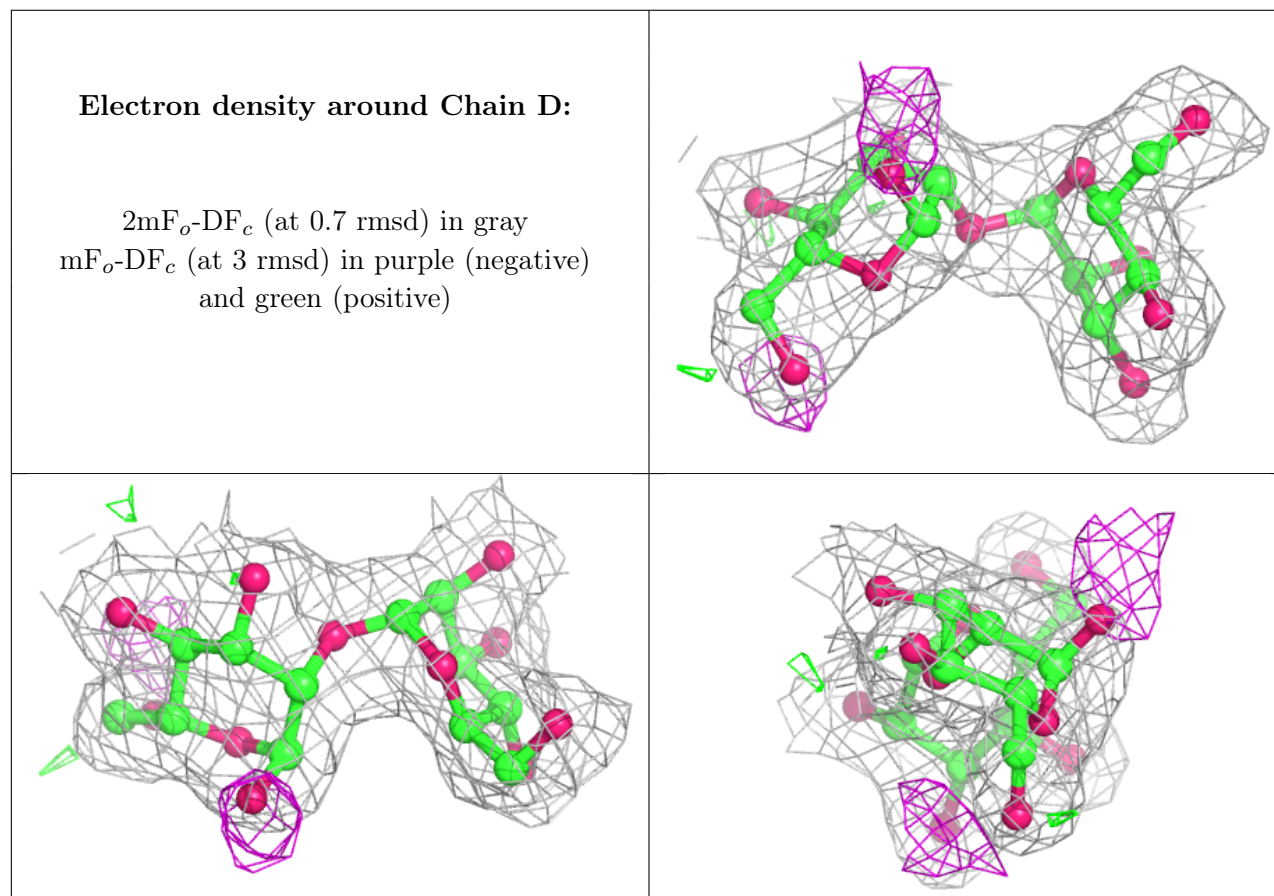
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MAN	D	1	12/12	0.96	0.11	33,37,41,44	0
2	MAN	D	2	11/12	0.97	0.11	33,35,42,42	0
2	MAN	E	2	11/12	0.97	0.10	28,34,39,40	0
2	MAN	F	1	12/12	0.97	0.11	29,34,40,41	0

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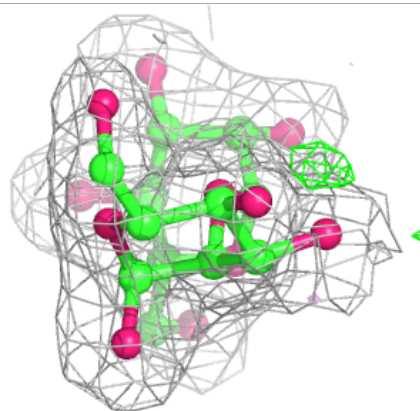
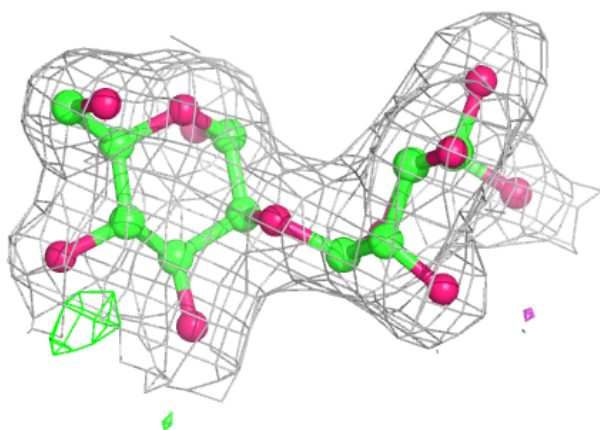
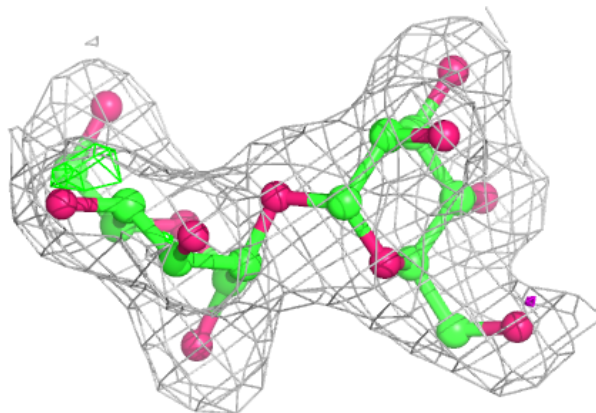
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	MAN	F	2	11/12	0.97	0.11	36,39,44,49	0
2	MAN	E	1	12/12	0.98	0.10	30,33,38,43	0

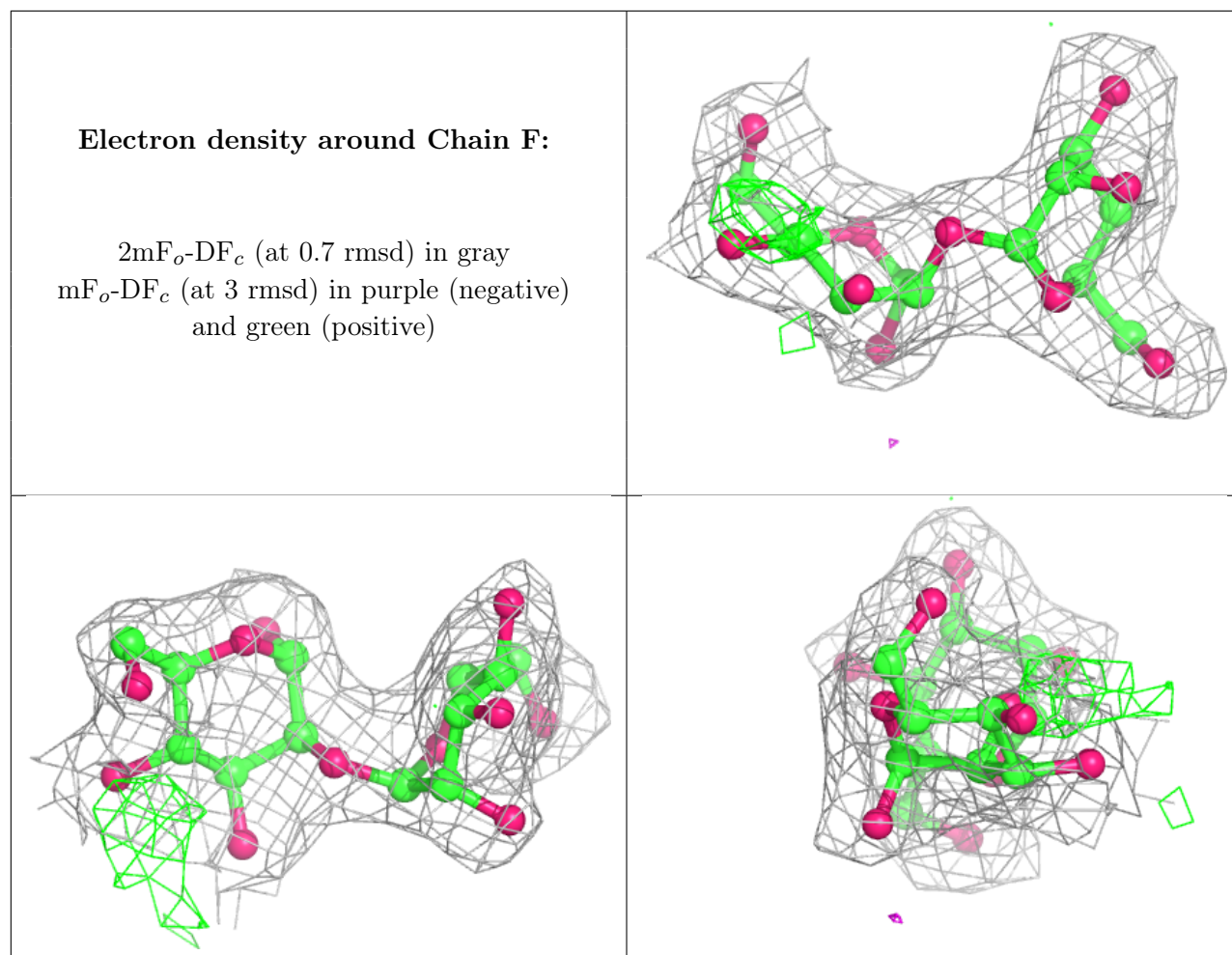
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around Chain 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
3	CA	B	403	1/1	0.98	0.11	41,41,41,41	0
3	CA	C	401	1/1	0.98	0.15	28,28,28,28	0
3	CA	C	403	1/1	0.98	0.12	41,41,41,41	0
3	CA	B	402	1/1	0.99	0.13	34,34,34,34	0
3	CA	A	401	1/1	0.99	0.13	33,33,33,33	0
3	CA	A	402	1/1	0.99	0.13	34,34,34,34	0
3	CA	C	402	1/1	0.99	0.13	30,30,30,30	0
3	CA	A	403	1/1	0.99	0.13	36,36,36,36	0
3	CA	B	401	1/1	1.00	0.16	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.