



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 21, 2023 – 02:09 PM JST

PDB ID : 7VEW
Title : Crystal structure of bacterial chemotaxis-dependent pectin-binding protein SPH1118 in complex with unsaturated trigalacturonic acid
Authors : Anamizu, K.; Takase, R.; Hio, M.; Watanebe, D.; Mikami, B.; Hashimoto, W.
Deposited on : 2021-09-10
Resolution : 1.92 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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

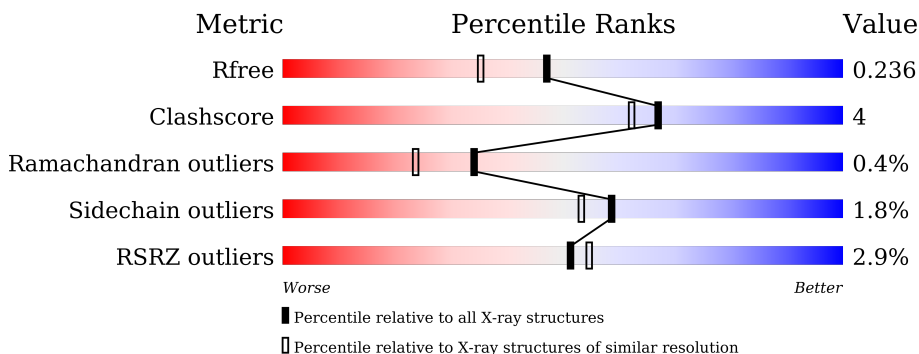
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.92 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	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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 $\leq 5\%$. 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	619	 3% (poor fit), 88% (0-1 outliers), 9% (2 outliers), 0% (3+ outliers), 0% (not modelled)
1	B	619	 3% (poor fit), 89% (0-1 outliers), 9% (2 outliers), 0% (3+ outliers), 0% (not modelled)
2	C	3	 0% (poor fit), 67% (0-1 outliers), 33% (2 outliers), 0% (3+ outliers), 0% (not modelled)
2	D	3	 0% (poor fit), 67% (0-1 outliers), 33% (2 outliers), 0% (3+ outliers), 0% (not modelled)

2 Entry composition [i](#)

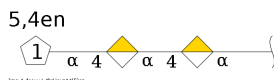
There are 6 unique types of molecules in this entry. The entry contains 10339 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SPH1118.

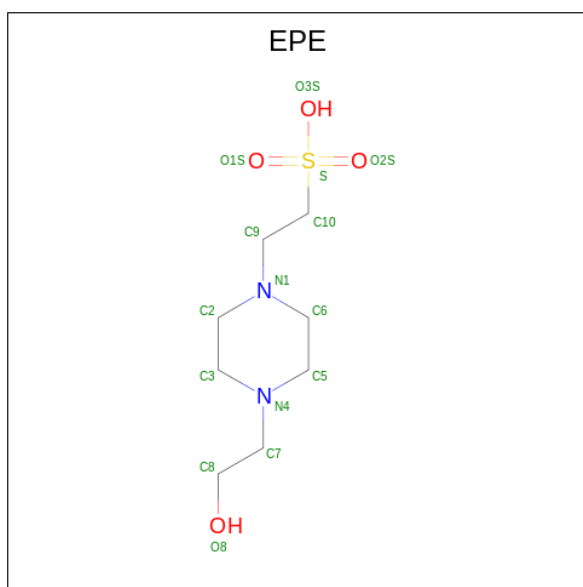
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	609	4854	3099	832	904	19	0	1	0
1	B	609	4854	3100	832	902	20	0	1	0

- Molecule 2 is an oligosaccharide called 2,6-anhydro-3-deoxy-L-threo-hex-2-enonic acid-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-alpha-D-galactopyranuronic acid.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	C	3	36	18	18	0	0	0
2	D	3	36	18	18	0	0	0

- Molecule 3 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	S			
3	A	1	Total	15	8	2	4	1	0	0
3	B	1	Total	15	8	2	4	1	0	0

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



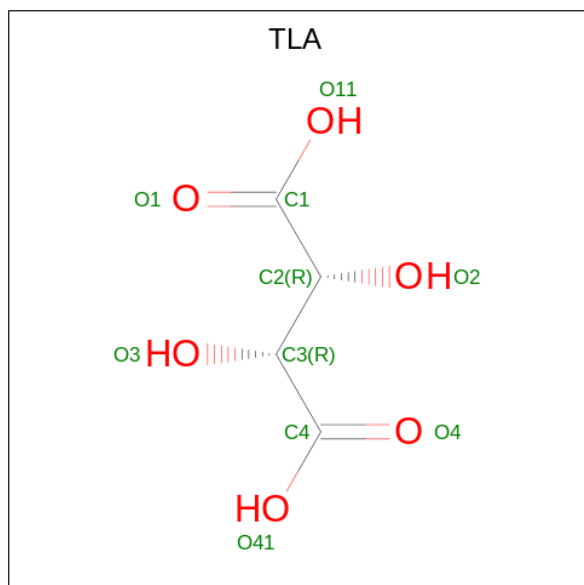
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	O			
4	A	1	Total	6	3	3	0	0
4	B	1	Total	6	3	3	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is L(+)-TARTARIC ACID (three-letter code: TLA) (formula: C₄H₆O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	4	6		
5	B	1	Total	C	O	0	0
			10	4	6		

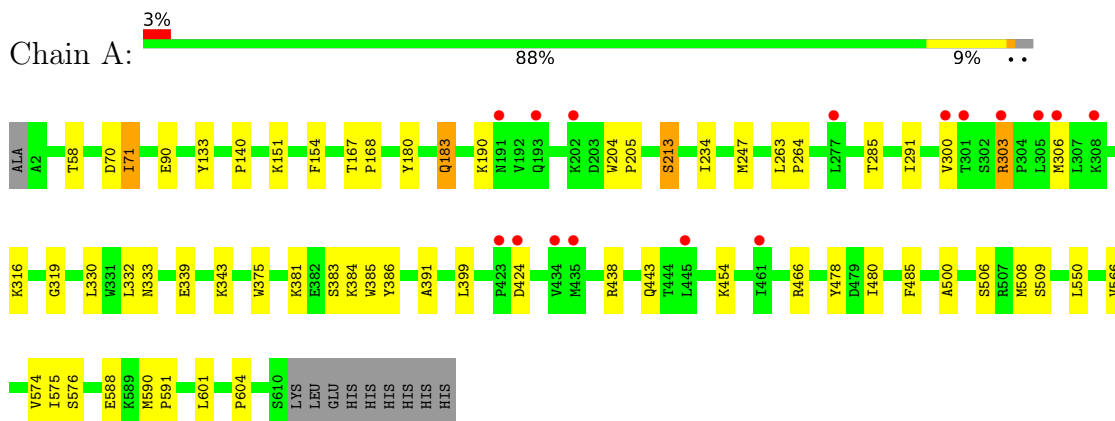
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	252	Total	O	0	0
			252	252		
6	B	227	Total	O	0	0
			227	227		

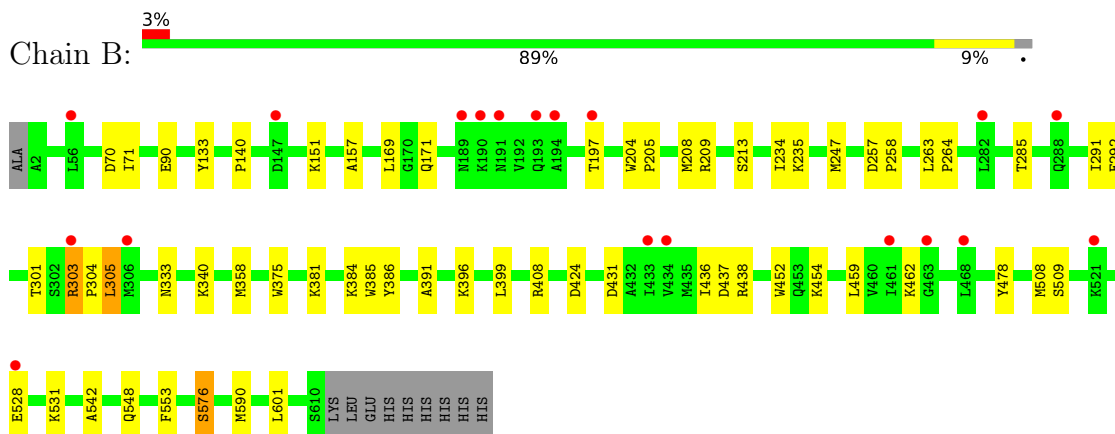
3 Residue-property plots [i](#)

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.

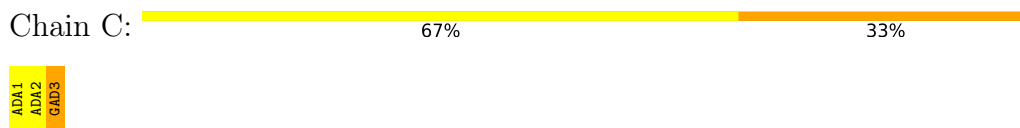
- Molecule 1: SPH1118



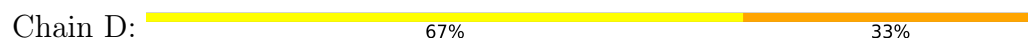
- Molecule 1: SPH1118



- Molecule 2: 2,6-anhydro-3-deoxy-L-threo-hex-2-enonic acid-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-alpha-D-galactopyranuronic acid



- Molecule 2: 2,6-anhydro-3-deoxy-L-threo-hex-2-enonic acid-(1-4)-alpha-D-galactopyranuronic acid-(1-4)-alpha-D-galactopyranuronic acid



ADA1
ADA2
GAD3

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	57.69Å 83.66Å 148.52Å 90.00° 101.04° 90.00°	Depositor
Resolution (Å)	48.59 – 1.92 48.59 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.59-1.92) 99.6 (48.59-1.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.69 (at 1.91Å)	Xtrriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.210 , 0.232 0.212 , 0.236	Depositor DCC
R_{free} test set	5421 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	29.6	Xtrriage
Anisotropy	0.684	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 41.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.013 for h,-k,-h-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	10339	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.46% 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: EPE, ADA, TLA, GOL, GAD

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	0.37	0/4976	0.53	0/6760
1	B	0.37	0/4976	0.52	0/6759
All	All	0.37	0/9952	0.53	0/13519

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	4854	0	4819	37	0
1	B	4854	0	4824	37	0
2	C	36	0	20	2	0
2	D	36	0	20	2	0
3	A	15	0	17	3	0
3	B	15	0	17	1	0
4	A	6	0	8	1	0
4	B	24	0	32	4	0
5	A	10	0	4	1	0
5	B	10	0	4	0	0
6	A	252	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	227	0	0	1	0
All	All	10339	0	9765	76	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 4.

All (76) 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:303:ARG:HH11	1:B:576:SER:HB2	1.42	0.82
1:A:90:GLU:OE2	1:A:151:LYS:HD2	1.84	0.76
1:B:303:ARG:NH1	1:B:576:SER:HB2	2.04	0.73
1:A:180:TYR:O	1:A:183:GLN:HG3	1.90	0.72
1:B:303:ARG:HD2	1:B:576:SER:CB	2.21	0.70
1:B:303:ARG:HD2	1:B:576:SER:OG	1.92	0.69
1:B:358[B]:MET:HG3	1:B:452:TRP:HE1	1.58	0.68
1:A:466:ARG:HE	4:A:802:GOL:H11	1.59	0.67
1:B:169:LEU:HD12	4:B:805:GOL:H12	1.77	0.66
1:A:384:LYS:H	3:A:801:EPE:H32	1.62	0.64
1:B:384:LYS:HG2	3:B:801:EPE:O1S	1.99	0.62
1:A:438:ARG:HD2	2:C:3:GAD:O6B	2.00	0.62
1:B:133:TYR:O	1:B:140:PRO:HD3	2.00	0.61
1:A:383:SER:HA	3:A:801:EPE:H81	1.85	0.59
1:B:263:LEU:HB3	1:B:264:PRO:HA	1.85	0.58
1:B:436:ILE:HD12	1:B:437:ASP:N	2.18	0.58
5:A:803:TLA:O11	5:A:803:TLA:O3	2.22	0.57
1:B:438:ARG:HD2	2:D:3:GAD:O6A	2.05	0.57
1:B:438:ARG:NH2	2:D:3:GAD:O6A	2.38	0.57
1:A:385:TRP:CZ3	1:A:550:LEU:HD12	2.40	0.56
1:A:133:TYR:O	1:A:140:PRO:HD3	2.04	0.56
1:A:438:ARG:NH1	2:C:3:GAD:O6B	2.40	0.55
1:B:303:ARG:HD2	1:B:576:SER:HB2	1.88	0.54
1:B:205:PRO:O	1:B:209:ARG:HG3	2.08	0.54
1:A:204:TRP:CD1	1:A:205:PRO:HD3	2.43	0.53
1:B:169:LEU:CD1	4:B:805:GOL:H12	2.38	0.53
1:A:234:ILE:HD13	1:A:247:MET:HG2	1.91	0.52
1:A:70:ASP:O	1:A:71:ILE:HG13	2.10	0.51
1:B:375:TRP:CZ2	1:B:391:ALA:HB1	2.45	0.51
1:A:303:ARG:HA	1:A:306:MET:HB2	1.91	0.51
1:B:385:TRP:CZ3	1:B:553:PHE:HB3	2.47	0.50
1:A:285:THR:HG22	1:A:306:MET:SD	2.52	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:399:LEU:HD21	1:A:454:LYS:HB3	1.93	0.49
1:B:399:LEU:HD21	1:B:454:LYS:HB3	1.95	0.49
1:B:90:GLU:CD	1:B:151:LYS:HD2	2.32	0.49
1:A:167:THR:HB	1:A:168:PRO:HD2	1.95	0.48
1:A:303:ARG:NE	1:A:576:SER:OG	2.46	0.48
1:B:431:ASP:HB3	1:B:462:LYS:HD2	1.96	0.48
1:A:303:ARG:NH2	1:A:574:VAL:O	2.47	0.47
1:B:528:GLU:OE2	1:B:531:LYS:NZ	2.40	0.47
1:B:333:ASN:HB2	1:B:478:TYR:CE2	2.49	0.47
1:A:339:GLU:O	1:A:343:LYS:NZ	2.46	0.47
4:B:805:GOL:H2	6:B:1086:HOH:O	2.16	0.46
1:A:300:VAL:HG13	1:A:303:ARG:HD2	1.97	0.46
1:B:381:LYS:HA	1:B:386:TYR:CD1	2.51	0.46
1:A:333:ASN:HB2	1:A:478:TYR:CE2	2.51	0.46
1:A:384:LYS:H	3:A:801:EPE:C3	2.28	0.46
1:A:375:TRP:CZ2	1:A:391:ALA:HB1	2.51	0.45
1:B:303:ARG:N	1:B:304:PRO:HD2	2.32	0.45
1:A:319:GLY:HA2	1:A:575:ILE:O	2.17	0.45
1:B:204:TRP:CZ3	1:B:208:MET:HG3	2.52	0.44
1:A:140:PRO:HB3	1:A:154:PHE:CD1	2.52	0.44
1:B:70:ASP:O	1:B:71:ILE:HG13	2.17	0.44
1:B:396:LYS:HD3	1:B:396:LYS:HA	1.64	0.44
1:B:285:THR:HG21	1:B:305:LEU:HD13	2.00	0.43
1:B:234:ILE:HD13	1:B:247:MET:HG2	2.00	0.43
1:A:263:LEU:HB3	1:A:264:PRO:HA	1.99	0.43
1:A:316:LYS:HE3	1:A:588:GLU:OE1	2.19	0.43
1:A:590:MET:HB2	1:A:601:LEU:HD23	2.01	0.42
1:A:485:PHE:CZ	1:A:566:VAL:HB	2.54	0.42
1:A:424:ASP:OD1	1:A:424:ASP:N	2.53	0.42
1:B:171:GLN:O	1:B:171:GLN:HG2	2.20	0.42
1:A:204:TRP:CG	1:A:205:PRO:HD3	2.55	0.42
1:A:381:LYS:HA	1:A:386:TYR:CD1	2.55	0.42
1:B:340:LYS:HE2	1:B:424:ASP:OD2	2.19	0.42
1:B:204:TRP:CD1	1:B:205:PRO:HD3	2.55	0.42
1:A:500:ALA:HA	1:A:506:SER:HB3	2.01	0.41
1:A:58:THR:O	1:A:604:PRO:HG3	2.19	0.41
1:B:257:ASP:HB2	1:B:258:PRO:CD	2.51	0.41
1:B:590:MET:HB2	1:B:601:LEU:HD23	2.03	0.41
1:A:306:MET:HA	1:A:306:MET:HE3	2.03	0.41
1:A:71:ILE:HG12	1:A:591:PRO:HG3	2.03	0.40
1:A:332:LEU:HD22	1:A:480:ILE:HG12	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:408:ARG:HE	4:B:804:GOL:C3	2.34	0.40
1:B:157:ALA:O	1:B:542:ALA:HA	2.22	0.40
1:B:358[B]:MET:HG3	1:B:452:TRP:NE1	2.33	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	608/619 (98%)	583 (96%)	22 (4%)	3 (0%)	29 18
1	B	608/619 (98%)	580 (95%)	26 (4%)	2 (0%)	41 31
All	All	1216/1238 (98%)	1163 (96%)	48 (4%)	5 (0%)	34 24

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	509	SER
1	B	213	SER
1	A	213	SER
1	B	509	SER
1	A	71	ILE

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	522/530 (98%)	514 (98%)	8 (2%)	65	61
1	B	522/530 (98%)	511 (98%)	11 (2%)	53	46
All	All	1044/1060 (98%)	1025 (98%)	19 (2%)	59	53

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	183	GLN
1	A	190	LYS
1	A	213	SER
1	A	291	ILE
1	A	303	ARG
1	A	330	LEU
1	A	443	GLN
1	A	508	MET
1	B	197	THR
1	B	235	LYS
1	B	291	ILE
1	B	292	GLU
1	B	301	THR
1	B	303	ARG
1	B	305	LEU
1	B	459	LEU
1	B	508	MET
1	B	548	GLN
1	B	576	SER

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

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates i

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	ADA	C	1	2	13,13,13	1.08	1 (7%)	18,19,19	0.93	0
2	ADA	C	2	2	12,12,13	1.09	1 (8%)	14,17,19	0.75	0
2	GAD	C	3	2	10,11,11	3.12	4 (40%)	13,15,15	1.36	3 (23%)
2	ADA	D	1	2	13,13,13	1.18	2 (15%)	18,19,19	0.98	1 (5%)
2	ADA	D	2	2	12,12,13	1.20	1 (8%)	14,17,19	0.72	0
2	GAD	D	3	2	10,11,11	3.27	3 (30%)	13,15,15	1.35	2 (15%)

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	ADA	C	1	2	-	0/4/24/24	0/1/1/1
2	ADA	C	2	2	-	0/4/21/24	0/1/1/1
2	GAD	C	3	2	-	0/4/17/17	0/1/1/1
2	ADA	D	1	2	-	0/4/24/24	0/1/1/1
2	ADA	D	2	2	-	1/4/21/24	0/1/1/1
2	GAD	D	3	2	-	0/4/17/17	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	3	GAD	C3-C4	-8.53	1.39	1.50
2	C	3	GAD	C3-C4	-7.97	1.40	1.50
2	D	3	GAD	C4-C5	4.05	1.39	1.33
2	C	3	GAD	C4-C5	3.81	1.39	1.33
2	D	3	GAD	C5-C6	-3.69	1.39	1.48
2	C	3	GAD	C5-C6	-3.40	1.40	1.48
2	D	1	ADA	C4-C5	2.67	1.57	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	ADA	O5-C5	2.40	1.47	1.43
2	C	1	ADA	C4-C5	2.27	1.57	1.53
2	C	2	ADA	O5-C5	2.24	1.47	1.43
2	D	1	ADA	C1-C2	2.04	1.57	1.52
2	C	3	GAD	O6B-C6	-2.02	1.24	1.30

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3	GAD	O6B-C6-C5	2.79	121.17	114.20
2	D	3	GAD	O5-C5-C6	2.59	115.41	111.52
2	C	3	GAD	C2-C3-C4	-2.30	109.18	112.32
2	C	3	GAD	O5-C5-C6	2.23	114.86	111.52
2	C	3	GAD	O6B-C6-C5	2.18	119.63	114.20
2	D	1	ADA	O4-C4-C5	2.03	114.30	109.74

There are no chirality outliers.

All (1) torsion outliers are listed below:

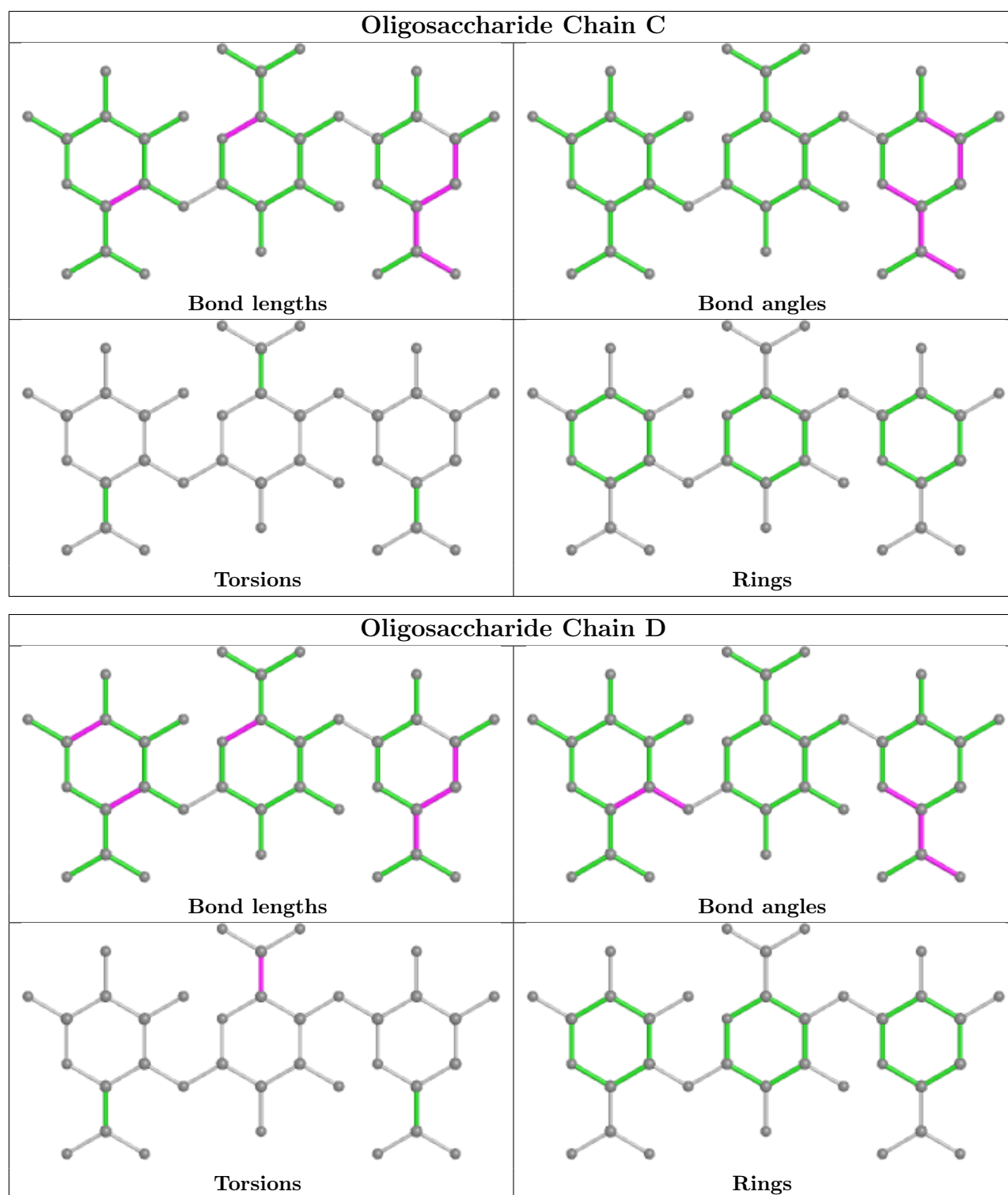
Mol	Chain	Res	Type	Atoms
2	D	2	ADA	C4-C5-C6-O6B

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	3	GAD	2	0
2	C	3	GAD	2	0

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



5.6 Ligand geometry [i](#)

9 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
5	TLA	B	806	-	9,9,9	1.11	0	12,12,12	1.01	0
4	GOL	B	802	-	5,5,5	0.95	0	5,5,5	1.49	1 (20%)
5	TLA	A	803	-	9,9,9	1.18	0	12,12,12	1.14	1 (8%)
3	EPE	A	801	-	15,15,15	0.89	1 (6%)	18,20,20	1.73	5 (27%)
4	GOL	B	803	-	5,5,5	0.76	0	5,5,5	1.20	0
3	EPE	B	801	-	15,15,15	0.90	1 (6%)	18,20,20	2.18	6 (33%)
4	GOL	A	802	-	5,5,5	0.42	0	5,5,5	0.91	0
4	GOL	B	804	-	5,5,5	0.81	0	5,5,5	1.12	0
4	GOL	B	805	-	5,5,5	0.79	0	5,5,5	1.20	1 (20%)

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
5	TLA	B	806	-	-	7/12/12/12	-
4	GOL	B	802	-	-	2/4/4/4	-
5	TLA	A	803	-	-	10/12/12/12	-
3	EPE	A	801	-	-	3/9/19/19	0/1/1/1
4	GOL	B	803	-	-	4/4/4/4	-
3	EPE	B	801	-	-	3/9/19/19	0/1/1/1
4	GOL	A	802	-	-	1/4/4/4	-
4	GOL	B	804	-	-	0/4/4/4	-
4	GOL	B	805	-	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	EPE	C10-S	2.93	1.81	1.77
3	B	801	EPE	C10-S	2.84	1.81	1.77

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	801	EPE	C5-N4-C3	4.42	118.77	108.83
3	B	801	EPE	C5-N4-C3	4.38	118.68	108.83
3	B	801	EPE	O2S-S-C10	4.19	111.97	106.92
3	B	801	EPE	C7-N4-C3	3.78	120.89	111.23
3	B	801	EPE	O3S-S-C10	3.18	110.91	105.77
3	B	801	EPE	C7-N4-C5	3.00	118.90	111.23
3	B	801	EPE	O1S-S-C10	-2.97	103.34	106.92
3	A	801	EPE	C7-N4-C3	2.96	118.81	111.23
4	B	802	GOL	C3-C2-C1	-2.54	101.81	111.70
3	A	801	EPE	O1S-S-C10	2.37	109.77	106.92
5	A	803	TLA	O1-C1-C2	-2.23	115.77	121.63
3	A	801	EPE	O2S-S-C10	2.18	109.54	106.92
3	A	801	EPE	C7-N4-C5	2.15	116.72	111.23
4	B	805	GOL	C3-C2-C1	-2.03	103.81	111.70

There are no chirality outliers.

All (30) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	803	GOL	C1-C2-C3-O3
5	A	803	TLA	O1-C1-C2-O2
5	A	803	TLA	O11-C1-C2-O2
5	A	803	TLA	C1-C2-C3-C4
5	A	803	TLA	C1-C2-C3-O3
5	A	803	TLA	O2-C2-C3-C4
5	A	803	TLA	O2-C2-C3-O3
4	B	803	GOL	O2-C2-C3-O3
4	B	802	GOL	C1-C2-C3-O3
4	B	803	GOL	O1-C1-C2-C3
5	A	803	TLA	C2-C3-C4-O4
5	A	803	TLA	C2-C3-C4-O41
5	B	806	TLA	C2-C3-C4-O4
3	B	801	EPE	C8-C7-N4-C5
4	B	802	GOL	O1-C1-C2-C3
5	B	806	TLA	C2-C3-C4-O41
3	A	801	EPE	C10-C9-N1-C2
3	A	801	EPE	C10-C9-N1-C6
3	B	801	EPE	C10-C9-N1-C6
3	A	801	EPE	N4-C7-C8-O8
5	A	803	TLA	O3-C3-C4-O41
5	B	806	TLA	O1-C1-C2-O2
5	A	803	TLA	O3-C3-C4-O4
5	B	806	TLA	O11-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
5	B	806	TLA	O1-C1-C2-C3
5	B	806	TLA	O11-C1-C2-O2
4	B	803	GOL	O1-C1-C2-O2
3	B	801	EPE	C10-C9-N1-C2
4	A	802	GOL	C1-C2-C3-O3
5	B	806	TLA	O3-C3-C4-O4

There are no ring outliers.

6 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	803	TLA	1	0
3	A	801	EPE	3	0
3	B	801	EPE	1	0
4	A	802	GOL	1	0
4	B	804	GOL	1	0
4	B	805	GOL	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	609/619 (98%)	0.24	16 (2%) 56 59	25, 35, 51, 62	0
1	B	609/619 (98%)	0.26	19 (3%) 49 52	28, 36, 52, 65	0
All	All	1218/1238 (98%)	0.25	35 (2%) 51 55	25, 36, 52, 65	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	303	ARG	3.5
1	A	305	LEU	3.1
1	B	190	LYS	3.0
1	A	193	GLN	2.9
1	A	308	LYS	2.9
1	A	461	ILE	2.8
1	B	191	ASN	2.7
1	B	303	ARG	2.6
1	B	193	GLN	2.6
1	B	194	ALA	2.5
1	A	300	VAL	2.4
1	B	288	GLN	2.4
1	A	277	LEU	2.4
1	A	306	MET	2.4
1	B	434	VAL	2.4
1	B	282	LEU	2.3
1	B	468	LEU	2.3
1	A	434	VAL	2.3
1	B	528	GLU	2.3
1	B	461	ILE	2.3
1	A	202	LYS	2.3
1	A	445	LEU	2.2
1	A	423	PRO	2.2
1	A	301	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	191	ASN	2.2
1	B	197	THR	2.2
1	B	147	ASP	2.2
1	B	521	LYS	2.2
1	A	424	ASP	2.1
1	A	435	MET	2.1
1	B	306	MET	2.1
1	B	433	ILE	2.1
1	B	463	GLY	2.1
1	B	189	ASN	2.0
1	B	56	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

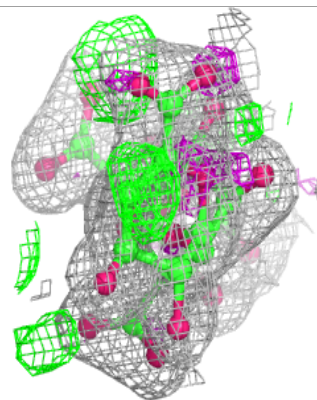
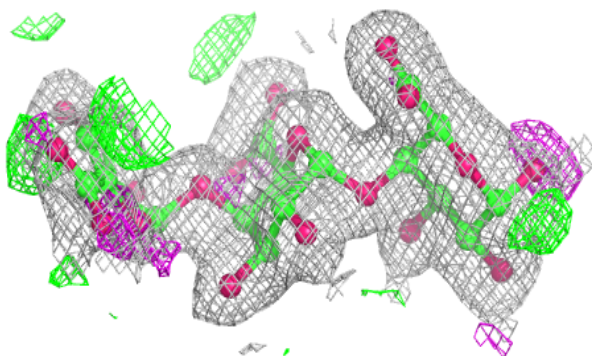
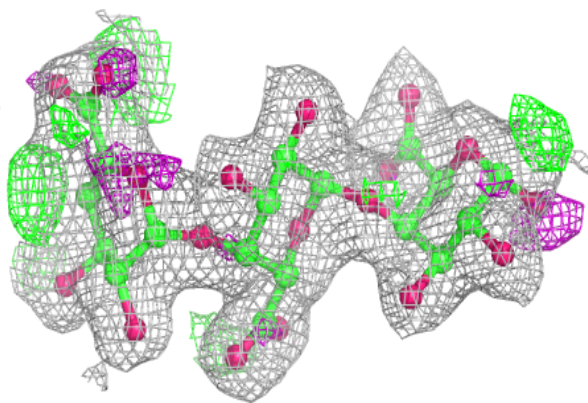
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GAD	D	3	11/11	0.84	0.13	34,38,45,47	0
2	GAD	C	3	11/11	0.88	0.14	32,36,43,45	0
2	ADA	D	2	12/13	0.92	0.13	25,30,32,32	0
2	ADA	C	2	12/13	0.93	0.12	24,28,29,30	0
2	ADA	C	1	13/13	0.94	0.14	23,26,29,34	0
2	ADA	D	1	13/13	0.94	0.15	26,28,30,36	0

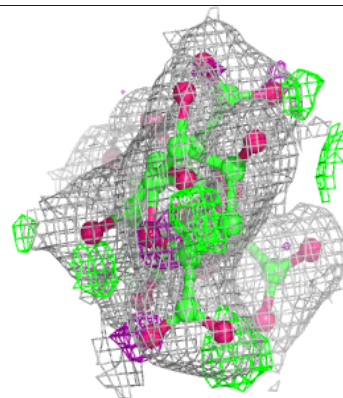
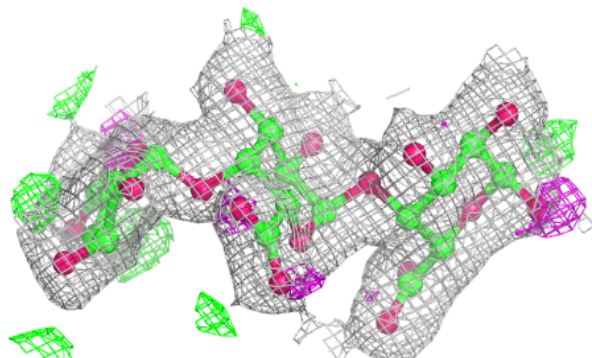
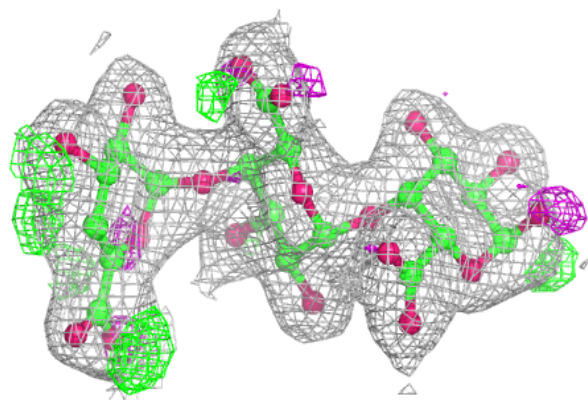
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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 D:**

$2mF_o-DF_c$ (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
4	GOL	A	802	6/6	0.59	0.30	41,42,45,49	0
3	EPE	A	801	15/15	0.64	0.27	39,56,85,90	0
4	GOL	B	804	6/6	0.72	0.31	48,54,57,60	0
4	GOL	B	803	6/6	0.78	0.25	47,48,51,53	0
5	TLA	A	803	10/10	0.80	0.27	44,51,55,59	0
4	GOL	B	805	6/6	0.83	0.20	41,42,45,46	0
3	EPE	B	801	15/15	0.84	0.23	42,58,69,69	0
4	GOL	B	802	6/6	0.85	0.17	42,47,49,50	0
5	TLA	B	806	10/10	0.88	0.27	49,51,54,55	0

6.5 Other polymers [i](#)

There are no such residues in this entry.