



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:57 PM UTC

PDB ID : 5GQD / pdb_00005gqd
Title : Crystal structure of covalent glycosyl-enzyme intermediate of xylanase mutant (T82A, N127S, and E128H) from *Streptomyces olivaceoviridis* E-86
Authors : Suzuki, R.; Fujimoto, Z.; Kaneko, S.; Kuno, A.
Deposited on : 2016-08-07
Resolution : 1.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

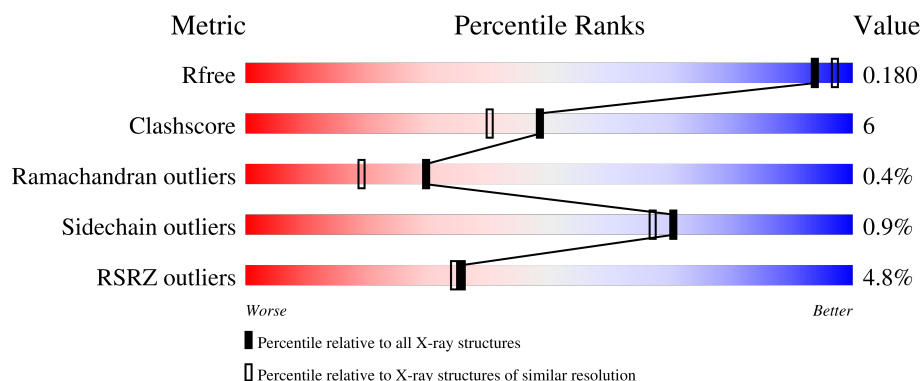
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	436	3% 86% 11% ..
1	B	436	7% 83% 14% ..
2	C	2	50% 50%
2	D	2	50% 50%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7547 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 Beta-xylanase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	427	Total	C	N	O	S	0	0	0
			3231	1987	588	640	16			
1	B	427	Total	C	N	O	S	0	0	0
			3231	1987	588	640	16			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	82	ALA	THR	engineered mutation	UNP Q7SI98
A	127	SER	ASN	engineered mutation	UNP Q7SI98
A	128	HIS	GLU	engineered mutation	UNP Q7SI98
B	582	ALA	THR	engineered mutation	UNP Q7SI98
B	627	SER	ASN	engineered mutation	UNP Q7SI98
B	628	HIS	GLU	engineered mutation	UNP Q7SI98

- Molecule 2 is an oligosaccharide called beta-D-xylopyranose-(1-4)-alpha-D-xylopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			18	10	8			
2	D	2	Total	C	O	0	0	0
			18	10	8			

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

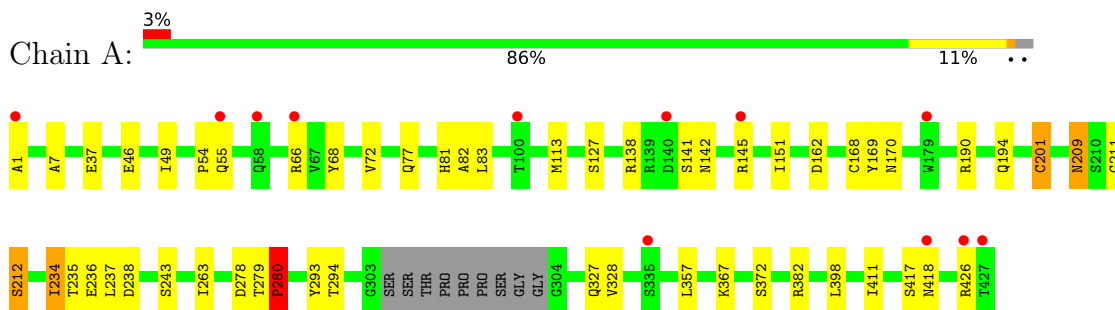
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	488	Total	O	0	0
			488	488		
4	B	513	Total	O	0	0
			513	513		

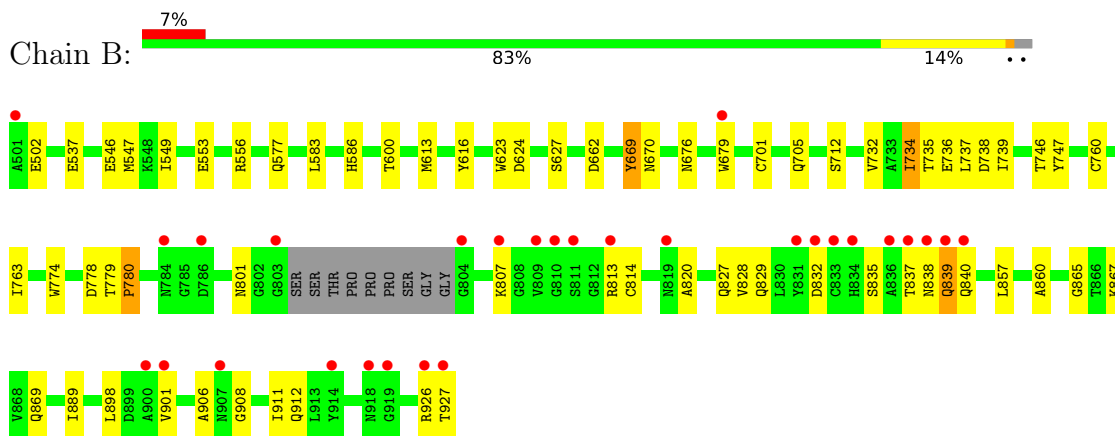
3 Residue-property plots [i](#)

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: Beta-xylanase



- Molecule 1: Beta-xylanase



- Molecule 2: beta-D-xylopyranose-(1-4)-alpha-D-xylopyranose



- Molecule 2: beta-D-xylopyranose-(1-4)-alpha-D-xylopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.19Å 94.05Å 140.08Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.09 – 1.80 40.09 – 1.80	Depositor EDS
% Data completeness (in resolution range)	98.2 (40.09-1.80) 98.4 (40.09-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	11.06 (at 1.79Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.175 , 0.197 0.179 , 0.180	Depositor DCC
R_{free} test set	4741 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	14.6	Xtriage
Anisotropy	0.125	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 58.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7547	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: XYP, GOL, XYS

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	0.35	1/3296 (0.0%)	0.91	14/4470 (0.3%)
1	B	0.35	1/3296 (0.0%)	0.91	17/4470 (0.4%)
All	All	0.35	2/6592 (0.0%)	0.91	31/8940 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	236	GLU	CD-OE2	7.28	1.39	1.25
1	B	736	GLU	CD-OE2	6.85	1.38	1.25

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	235	THR	N-CA-C	8.32	123.56	113.41
1	B	735	THR	N-CA-C	8.08	123.26	113.41
1	A	357	LEU	N-CA-C	-7.11	99.30	109.96
1	A	46	GLU	N-CA-C	7.09	121.02	112.38
1	B	583	LEU	N-CA-C	7.08	120.40	111.69
1	B	537	GLU	N-CA-C	6.83	121.70	113.23
1	A	37	GLU	N-CA-C	6.68	121.52	113.23
1	A	83	LEU	N-CA-C	6.68	119.91	111.69
1	B	546	GLU	N-CA-C	6.65	120.49	112.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	857	LEU	N-CA-C	-6.62	100.09	109.96
1	B	549	ILE	N-CA-C	6.07	116.24	110.42
1	A	398	LEU	N-CA-C	-6.06	101.07	110.10
1	B	839	GLN	N-CA-C	-5.91	98.22	110.80
1	A	49	ILE	N-CA-C	5.75	115.94	110.42
1	A	212	SER	CA-C-N	5.70	125.60	119.85
1	A	212	SER	C-N-CA	5.70	125.60	119.85
1	B	669	TYR	N-CA-C	-5.69	99.91	109.07
1	B	737	LEU	N-CA-C	5.60	118.44	110.10
1	B	898	LEU	N-CA-C	-5.52	101.87	110.10
1	A	77	GLN	N-CA-C	-5.45	103.50	110.53
1	B	577	GLN	N-CA-C	-5.41	102.75	110.59
1	A	237	LEU	N-CA-C	5.31	118.01	110.10
1	B	624	ASP	N-CA-C	-5.19	98.86	107.99
1	A	169	TYR	N-CA-C	-5.18	100.08	108.52
1	B	553	GLU	CA-C-N	5.16	124.95	119.28
1	B	553	GLU	C-N-CA	5.16	124.95	119.28
1	A	280	PRO	N-CA-C	5.13	121.03	114.92
1	B	712	SER	CA-C-N	5.12	125.02	119.85
1	B	712	SER	C-N-CA	5.12	125.02	119.85
1	A	234	ILE	N-CA-C	-5.05	99.84	107.73
1	B	734	ILE	N-CA-C	-5.00	99.93	107.73

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	669	TYR	Sidechain

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	3231	0	3041	28	0
1	B	3231	0	3038	41	0
2	C	18	0	7	0	0
2	D	18	0	7	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	18	0	24	1	0
3	B	30	0	40	3	0
4	A	488	0	0	3	0
4	B	513	0	0	2	0
All	All	7547	0	6157	70	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 6.

All (70) 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:926:ARG:HA	1:B:927:THR:HB	1.56	0.87
1:B:807:LYS:HG2	1:B:814:CYS:SG	2.21	0.80
1:B:807:LYS:HG3	1:B:927:THR:HG21	1.62	0.80
1:A:327:GLN:HB2	1:A:367:LYS:HE3	1.64	0.79
1:B:547:MET:HE3	1:B:616:TYR:CE2	2.16	0.79
1:B:827:GLN:HB2	1:B:867:LYS:HE3	1.76	0.67
1:B:820:ALA:HA	1:B:838:ASN:HB3	1.76	0.67
1:B:547:MET:HE2	1:B:623:TRP:CH2	2.34	0.61
1:B:828:VAL:HG23	1:B:911:ILE:HB	1.82	0.61
3:A:503:GOL:H31	4:A:626:HOH:O	2.02	0.59
1:A:142:ASN:HA	1:A:145:ARG:HD2	1.84	0.59
1:B:865:GLY:HA2	1:B:912:GLN:OE1	2.03	0.58
1:B:829:GLN:OE1	1:B:908:GLY:HA2	2.04	0.58
1:A:234:ILE:HD12	1:A:263:ILE:HG12	1.86	0.57
1:B:926:ARG:N	1:B:926:ARG:HD2	2.19	0.57
1:B:734:ILE:HD12	1:B:763:ILE:HG12	1.87	0.57
1:B:837:THR:C	1:B:839:GLN:H	2.13	0.56
1:A:328:VAL:HG23	1:A:411:ILE:HB	1.88	0.56
1:B:778:ASP:O	1:B:779:THR:C	2.48	0.56
1:B:738:ASP:HB2	1:B:780:PRO:HB2	1.87	0.56
1:A:278:ASP:O	1:A:279:THR:C	2.48	0.55
1:B:676:ASN:HB3	1:B:679:TRP:CD2	2.40	0.55
1:B:547:MET:HE2	1:B:623:TRP:HH2	1.71	0.54
1:B:813:ARG:HG2	1:B:813:ARG:HH11	1.73	0.54
1:A:190:ARG:O	1:A:194:GLN:HG3	2.08	0.54
1:B:926:ARG:CA	1:B:927:THR:HB	2.33	0.54
1:B:705:GLN:OE1	3:B:1007:GOL:H31	2.08	0.53
1:A:209:ASN:C	1:A:209:ASN:HD22	2.16	0.53
1:A:238:ASP:HB2	1:A:280:PRO:HB2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:SER:HA	1:A:170:ASN:O	2.12	0.49
1:A:372:SER:HB2	4:A:705:HOH:O	2.11	0.49
1:B:813:ARG:NH1	1:B:832:ASP:OD1	2.45	0.49
1:B:779:THR:N	1:B:780:PRO:HD3	2.28	0.48
1:A:294:THR:OG1	1:A:382:ARG:HD2	2.12	0.48
1:B:926:ARG:HA	1:B:927:THR:OXT	2.13	0.48
1:B:889:ILE:HD11	1:B:926:ARG:HH21	1.79	0.48
1:A:417:SER:O	1:A:418:ASN:HB2	2.15	0.47
1:A:66:ARG:HH11	1:A:66:ARG:HG3	1.79	0.46
1:B:807:LYS:CG	1:B:927:THR:HG21	2.40	0.46
1:A:243:SER:HB2	4:A:689:HOH:O	2.14	0.46
1:A:279:THR:N	1:A:280:PRO:HD3	2.30	0.46
1:B:774:TRP:CH2	3:B:1007:GOL:H32	2.51	0.46
1:B:732:VAL:HG23	1:B:760:CYS:HA	1.99	0.45
1:B:774:TRP:HH2	3:B:1007:GOL:H32	1.81	0.45
1:B:613:MET:HE2	1:B:662:ASP:HB3	1.99	0.45
1:B:860:ALA:HB2	1:B:869:GLN:HE22	1.83	0.44
1:B:837:THR:C	1:B:839:GLN:N	2.74	0.44
1:A:426:ARG:HB2	1:A:426:ARG:NH1	2.33	0.44
1:B:739:ILE:HD13	1:B:746:THR:HG22	2.00	0.44
1:B:807:LYS:HE2	1:B:927:THR:OG1	2.18	0.44
1:A:293:TYR:CD1	1:A:293:TYR:C	2.97	0.43
1:B:556:ARG:HH12	1:B:600:THR:CG2	2.32	0.43
1:B:813:ARG:HE	1:B:906:ALA:HA	1.84	0.43
1:A:1:ALA:HB3	1:A:7:ALA:HB1	2.01	0.42
1:B:820:ALA:HA	1:B:838:ASN:CB	2.45	0.42
1:B:586:HIS:HD2	4:B:1241:HOH:O	2.01	0.42
1:B:502:GLU:HB2	1:B:801:ASN:OD1	2.19	0.42
1:B:627:SER:HA	1:B:670:ASN:O	2.19	0.42
1:B:926:ARG:HB3	1:B:927:THR:OXT	2.20	0.42
1:A:209:ASN:HD22	1:A:212:SER:H	1.67	0.42
1:B:586:HIS:HE1	4:B:1525:HOH:O	2.02	0.41
1:A:168:CYS:HA	1:A:201:CYS:O	2.20	0.41
1:A:54:PRO:HB2	1:A:55:GLN:NE2	2.36	0.41
1:A:81:HIS:HA	1:A:82:ALA:HA	1.86	0.41
1:A:141:SER:O	1:A:145:ARG:HG3	2.21	0.41
1:A:209:ASN:ND2	1:A:212:SER:H	2.20	0.40
1:A:209:ASN:HD22	1:A:211:GLY:H	1.69	0.40
1:A:113:MET:HE2	1:A:162:ASP:HB3	2.03	0.40
1:A:68:TYR:CZ	1:A:72:VAL:HG21	2.57	0.40
1:A:138:ARG:CZ	1:A:151:ILE:HD12	2.52	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	423/436 (97%)	414 (98%)	9 (2%)	0	100	100
1	B	423/436 (97%)	409 (97%)	11 (3%)	3 (1%)	18	8
All	All	846/872 (97%)	823 (97%)	20 (2%)	3 (0%)	30	19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	835	SER
1	B	840	GLN
1	B	901	VAL

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	333/340 (98%)	330 (99%)	3 (1%)	70	67
1	B	333/340 (98%)	330 (99%)	3 (1%)	70	67
All	All	666/680 (98%)	660 (99%)	6 (1%)	70	67

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	201	CYS
1	A	209	ASN
1	A	280	PRO
1	B	701	CYS
1	B	747	TYR
1	B	780	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	55	GLN
1	A	74	ASN
1	A	194	GLN
1	A	209	ASN
1	A	298	ASN
1	A	334	HIS
1	B	511	GLN
1	B	558	GLN
1	B	574	ASN
1	B	586	HIS
1	B	673	ASN
1	B	683	GLN
1	B	723	GLN
1	B	819	ASN

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](#)

4 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	XYS	C	1	1,2	9,9,10	0.59	0	10,12,14	0.91	1 (10%)
2	XYP	C	2	2	9,9,10	0.58	0	10,12,14	0.73	0
2	XYS	D	1	1,2	9,9,10	0.60	0	10,12,14	0.95	1 (10%)
2	XYP	D	2	2	9,9,10	0.61	0	10,12,14	0.73	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
2	XYS	C	1	1,2	-	-	0/1/1/1
2	XYP	C	2	2	-	-	0/1/1/1
2	XYS	D	1	1,2	-	-	0/1/1/1
2	XYP	D	2	2	-	-	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1	XYS	C4-C3-C2	-2.33	108.15	110.92
2	C	1	XYS	C4-C3-C2	-2.21	108.29	110.92

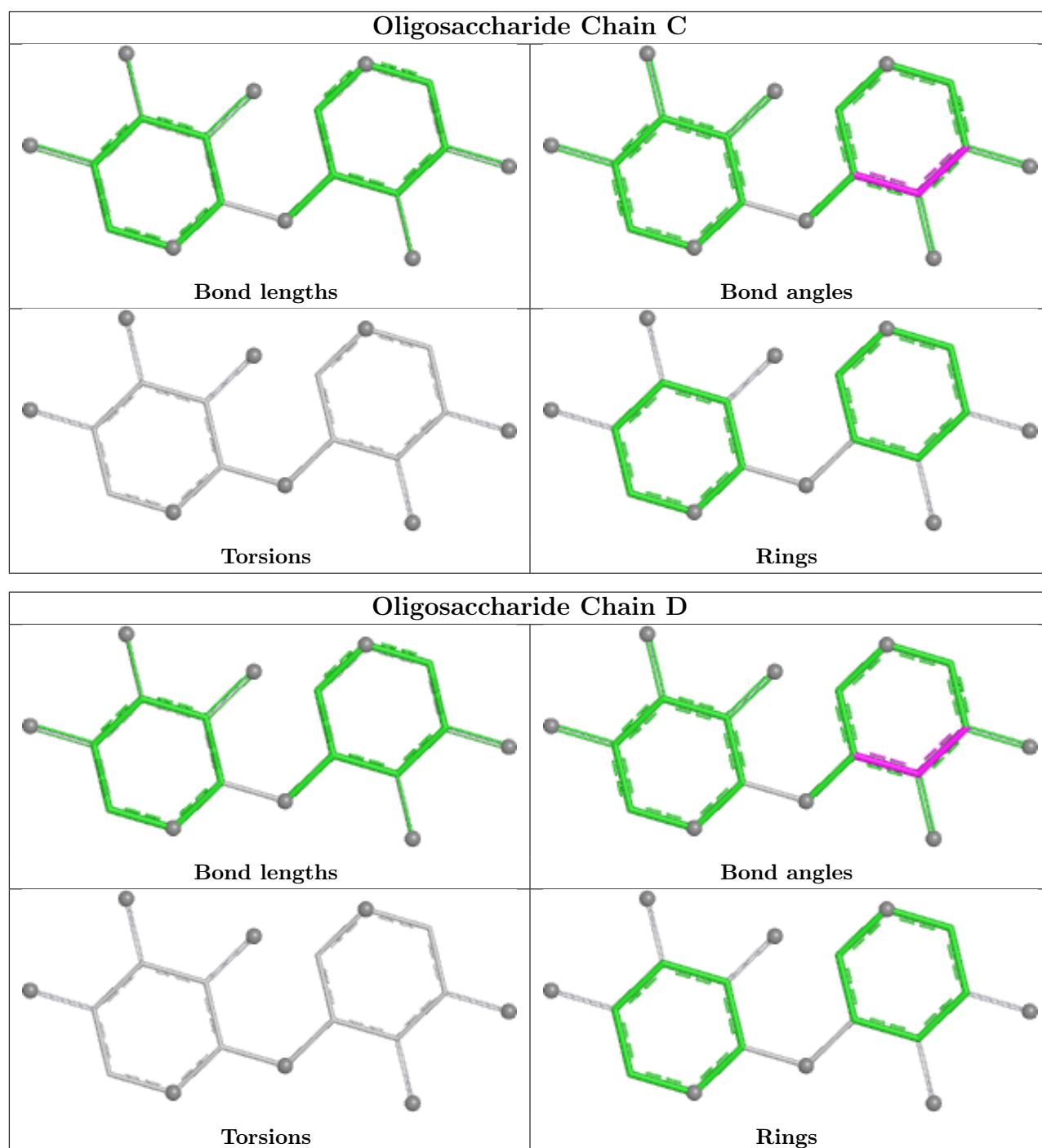
There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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](#)

8 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	$\# Z > 2$
3	GOL	A	503	-	5,5,5	0.18	0	5,5,5	0.34	0
3	GOL	A	505	-	5,5,5	0.18	0	5,5,5	0.31	0
3	GOL	A	504	-	5,5,5	0.15	0	5,5,5	0.28	0
3	GOL	B	1006	-	5,5,5	0.16	0	5,5,5	0.34	0
3	GOL	B	1003	-	5,5,5	0.14	0	5,5,5	0.35	0
3	GOL	B	1004	-	5,5,5	0.19	0	5,5,5	0.34	0
3	GOL	B	1005	-	5,5,5	0.14	0	5,5,5	0.33	0
3	GOL	B	1007	-	5,5,5	0.15	0	5,5,5	0.35	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
3	GOL	A	503	-	-	0/4/4/4	-
3	GOL	A	505	-	-	0/4/4/4	-
3	GOL	A	504	-	-	0/4/4/4	-
3	GOL	B	1006	-	-	0/4/4/4	-
3	GOL	B	1003	-	-	0/4/4/4	-
3	GOL	B	1004	-	-	0/4/4/4	-
3	GOL	B	1005	-	-	0/4/4/4	-
3	GOL	B	1007	-	-	0/4/4/4	-

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.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	503	GOL	1	0
3	B	1007	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	427/436 (97%)	-0.04	12 (2%) 55 55	7, 15, 29, 44	0
1	B	427/436 (97%)	-0.07	29 (6%) 23 22	6, 13, 33, 52	0
All	All	854/872 (97%)	-0.05	41 (4%) 35 34	6, 14, 31, 52	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	810	GLY	10.0
1	B	927	THR	6.6
1	A	1	ALA	5.7
1	B	501	ALA	5.2
1	B	839	GLN	5.1
1	B	926	ARG	4.9
1	A	427	THR	4.5
1	B	918	ASN	4.0
1	B	840	GLN	3.5
1	B	809	VAL	3.2
1	A	335	SER	3.2
1	A	145	ARG	3.2
1	B	811	SER	3.0
1	B	813	ARG	3.0
1	B	807	LYS	2.9
1	B	837	THR	2.9
1	B	784	ASN	2.8
1	B	834	HIS	2.7
1	B	919	GLY	2.6
1	A	418	ASN	2.6
1	A	58	GLN	2.5
1	A	179	TRP	2.5
1	A	140	ASP	2.4
1	B	833	CYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	55	GLN	2.3
1	A	100	THR	2.3
1	B	819	ASN	2.3
1	B	786	ASP	2.3
1	A	426	ARG	2.2
1	B	679	TRP	2.2
1	B	803	GLY	2.2
1	B	914	TYR	2.2
1	B	836	ALA	2.2
1	A	66	ARG	2.2
1	B	901	VAL	2.2
1	B	838	ASN	2.1
1	B	832	ASP	2.1
1	B	907	ASN	2.1
1	B	831	TYR	2.0
1	B	804	GLY	2.0
1	B	900	ALA	2.0

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 ⓘ

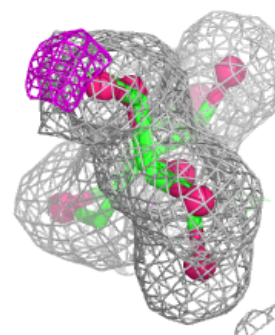
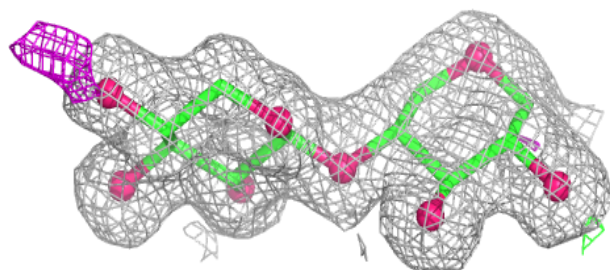
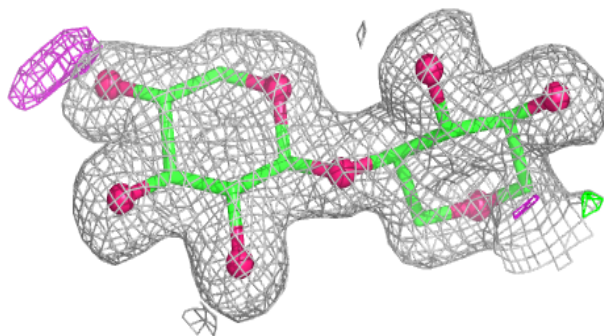
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	XYS	C	1	9/10	0.95	0.06	14,15,16,16	0
2	XYP	C	2	9/10	0.95	0.07	13,15,16,18	0
2	XYP	D	2	9/10	0.96	0.06	9,10,10,11	0
2	XYS	D	1	9/10	0.97	0.05	10,10,11,12	0

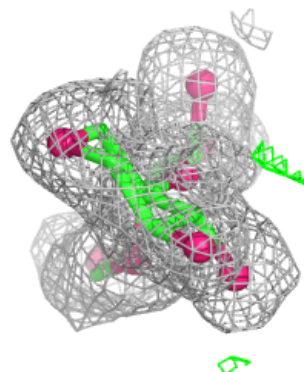
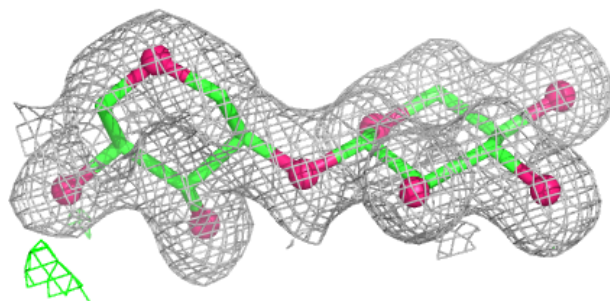
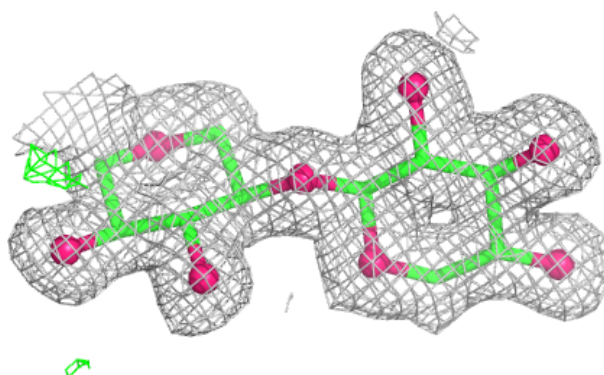
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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($\text{\AA}^2$)	Q<0.9
3	GOL	B	1006	6/6	0.71	0.22	36,38,39,43	0
3	GOL	B	1007	6/6	0.76	0.27	48,48,49,50	0
3	GOL	A	503	6/6	0.78	0.22	25,30,33,34	0
3	GOL	B	1005	6/6	0.78	0.21	44,44,45,45	0
3	GOL	A	505	6/6	0.83	0.19	24,30,31,35	0
3	GOL	B	1004	6/6	0.85	0.15	28,30,31,33	0
3	GOL	A	504	6/6	0.89	0.13	21,25,26,27	0
3	GOL	B	1003	6/6	0.91	0.10	25,25,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.