



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 12:24 AM UTC

PDB ID : 2FV1 / pdb_00002fv1
Title : UGL_D88N/dGlcA-GlcNAc
Authors : Itoh, T.; Hashimoto, W.; Mikami, B.; Murata, K.
Deposited on : 2006-01-28
Resolution : 1.73 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

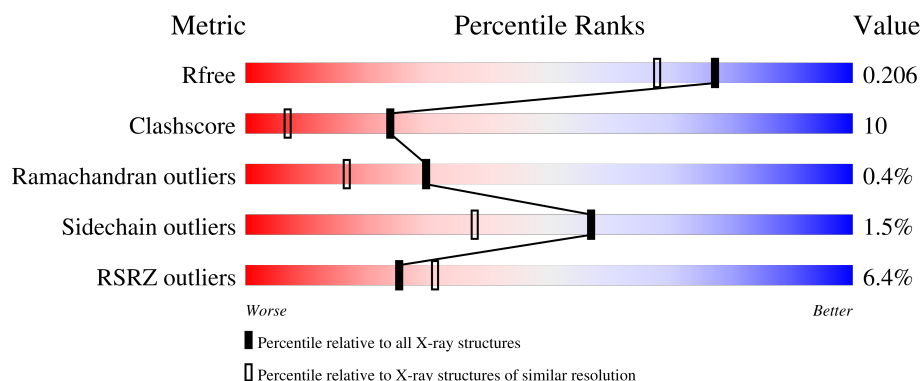
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.73 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	1187 (1.74-1.74)
Clashscore	190562	1207 (1.74-1.74)
Ramachandran outliers	187476	1200 (1.74-1.74)
Sidechain outliers	187428	1200 (1.74-1.74)
RSRZ outliers	180081	1188 (1.74-1.74)

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	377	3% 80% 18% .
1	B	377	9% 79% 19% .
2	C	2	50% 50%
2	D	2	100%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 7016 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.

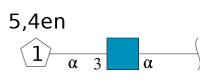
- Molecule 1 is a protein called Unsaturated glucuronyl hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	377	Total	C	N	O	S	0	10	0
			3094	1941	564	583	6			
1	B	377	Total	C	N	O	S	0	7	0
			3081	1934	562	579	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	88	ASN	ASP	engineered mutation	UNP Q9RC92
B	88	ASN	ASP	engineered mutation	UNP Q9RC92

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	N	O	0	0	0
			26	14	1	11			
2	D	2	Total	C	N	O	0	0	0
			26	14	1	11			

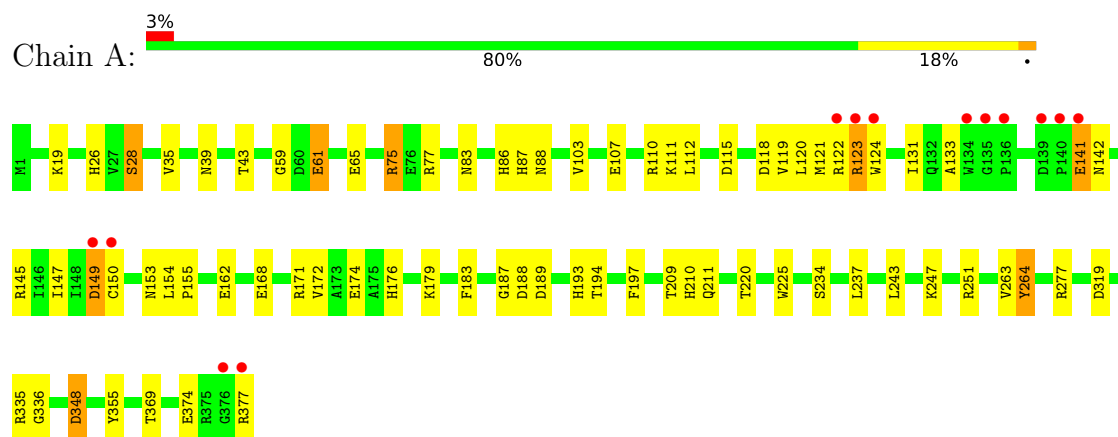
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	437	Total	O	0	0
			437	437		
3	B	352	Total	O	0	0
			352	352		

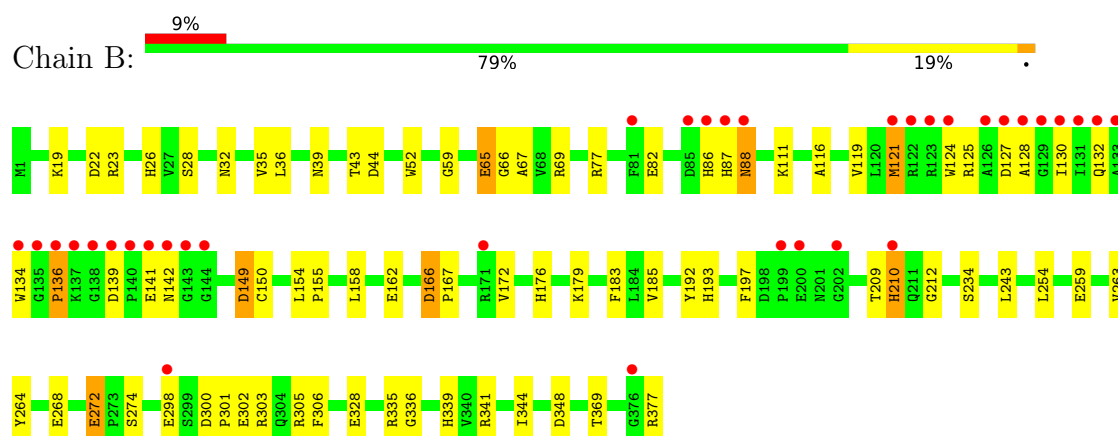
3 Residue-property plots

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: Unsaturated glucuronyl hydrolase



- Molecule 1: Unsaturated glucuronyl hydrolase



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



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

Chain D:

100%

NDJ1
GAD2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.73Å 94.89Å 95.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.95 – 1.73 14.95 – 1.73	Depositor EDS
% Data completeness (in resolution range)	95.6 (14.95-1.73) 95.6 (14.95-1.73)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.98 (at 1.73Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.179 , 0.212 0.173 , 0.206	Depositor DCC
R_{free} test set	7994 reflections (10.02%)	wwPDB-VP
Wilson B-factor (Å ²)	15.9	Xtriage
Anisotropy	0.234	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 56.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.013 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7016	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.00% 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: NDG, 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	0.34	0/3212	0.88	17/4346 (0.4%)
1	B	0.33	0/3188	0.89	13/4316 (0.3%)
All	All	0.33	0/6400	0.89	30/8662 (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	A	0	1

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	263	VAL	N-CA-C	8.80	120.15	110.21
1	B	263	VAL	N-CA-C	8.43	119.74	110.21
1	B	336	GLY	N-CA-C	-8.21	99.93	112.89
1	A	336	GLY	N-CA-C	-7.73	100.68	112.89
1	A	369	THR	N-CA-C	7.06	120.77	112.72
1	B	369	THR	N-CA-C	7.00	120.70	112.72
1	A	335	ARG	N-CA-C	6.14	120.03	111.74
1	A	348	ASP	N-CA-C	5.96	116.43	108.38
1	B	335	ARG	N-CA-C	5.91	119.72	111.74
1	A	264	TYR	N-CA-C	-5.74	103.13	110.53
1	B	264	TYR	N-CA-C	-5.70	103.18	110.53
1	A	220	THR	N-CA-C	5.68	117.47	111.28
1	B	259	GLU	N-CA-C	5.67	118.19	111.33
1	A	75[A]	ARG	N-CA-C	-5.58	105.11	111.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	75[B]	ARG	N-CA-C	-5.58	105.11	111.14
1	B	210	HIS	N-CA-C	-5.52	106.45	114.12
1	B	183	PHE	N-CA-C	5.44	120.19	113.50
1	B	192	TYR	N-CA-C	-5.41	102.51	110.52
1	B	35	VAL	N-CA-C	-5.36	100.70	108.36
1	A	35	VAL	N-CA-C	-5.34	100.72	108.36
1	B	59	GLY	N-CA-C	-5.30	107.94	115.30
1	A	103	VAL	N-CA-C	5.25	115.47	110.53
1	A	28	SER	N-CA-C	-5.22	102.73	110.46
1	A	59	GLY	N-CA-C	-5.21	108.43	115.36
1	B	166	ASP	CA-C-N	5.19	124.91	119.82
1	B	166	ASP	C-N-CA	5.19	124.91	119.82
1	A	183	PHE	N-CA-C	5.18	120.59	113.97
1	A	210	HIS	N-CA-C	-5.15	105.88	113.61
1	A	123	ARG	N-CA-C	-5.13	107.02	113.28
1	A	319	ASP	N-CA-C	5.04	116.46	111.07

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	355	TYR	Sidechain

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	3094	0	2925	56	0
1	B	3081	0	2912	65	0
2	C	26	0	17	3	0
2	D	26	0	17	4	0
3	A	437	0	0	7	0
3	B	352	0	0	7	0
All	All	7016	0	5871	122	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 10.

All (122) 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:339[A]:HIS:NE2	2:D:1:NDG:H6C2	1.86	0.90
1:A:377:ARG:HD3	1:A:377:ARG:OXT	1.80	0.81
1:A:131:ILE:H	1:A:176:HIS:HD2	1.38	0.72
1:B:22:ASP:HB3	1:B:69[A]:ARG:NH2	2.04	0.71
1:B:193:HIS:HB3	1:B:209:THR:HB	1.73	0.70
1:A:119:VAL:HA	1:A:122[A]:ARG:NH1	2.05	0.70
1:A:86:HIS:HD2	1:A:88:ASN:H	1.40	0.69
1:B:28:SER:OG	1:B:348:ASP:HB3	1.93	0.68
1:B:111:LYS:HD2	3:B:697:HOH:O	1.94	0.68
1:B:19:LYS:HE2	3:B:675:HOH:O	1.94	0.66
1:B:22:ASP:HB3	1:B:69[A]:ARG:HH21	1.61	0.66
1:A:86:HIS:CD2	1:A:88:ASN:HB2	2.32	0.65
1:A:131:ILE:H	1:A:176:HIS:CD2	2.16	0.63
1:B:116:ALA:O	1:B:119:VAL:HG22	1.99	0.63
1:B:298:GLU:H	1:B:298:GLU:CD	2.07	0.62
1:A:87:HIS:CE1	1:A:150:CYS:HB2	2.34	0.62
1:A:133:ALA:CB	1:A:147:ILE:HG12	2.30	0.61
1:B:77:ARG:HG3	1:B:82:GLU:HB3	1.82	0.61
1:B:86:HIS:CD2	1:B:88:ASN:HB2	2.36	0.60
1:B:87:HIS:CD2	1:B:150:CYS:SG	2.95	0.60
1:A:176:HIS:HE1	3:A:572:HOH:O	1.84	0.60
1:A:168:GLU:OE1	1:A:171:ARG:HD3	2.02	0.60
1:B:65:GLU:HB3	1:B:69[B]:ARG:NH2	2.17	0.59
1:B:179:LYS:HE2	1:B:197:PHE:CE2	2.39	0.58
1:B:130:ILE:HG22	1:B:179:LYS:HD3	1.86	0.57
1:A:174:GLU:HG3	1:A:237:LEU:HD13	1.87	0.57
1:B:377:ARG:HH21	1:B:377:ARG:HG2	1.70	0.57
1:B:121:MET:HA	1:B:121:MET:HE2	1.86	0.56
1:A:119:VAL:HA	1:A:122[A]:ARG:HH11	1.71	0.56
1:A:88:ASN:ND2	2:C:2:GAD:O2	2.34	0.55
1:A:111:LYS:HE2	1:A:115:ASP:OD1	2.06	0.55
1:A:75[A]:ARG:HE	1:A:112:LEU:HD13	1.72	0.54
1:A:119:VAL:HG13	1:A:122[A]:ARG:HH12	1.71	0.54
1:A:28:SER:OG	1:A:348:ASP:HB3	2.08	0.54
1:B:339[A]:HIS:HB2	1:B:344:ILE:HB	1.90	0.54
1:B:125:ARG:HH11	1:B:132:GLN:HE21	1.55	0.54
1:A:193:HIS:HB3	1:A:209:THR:HB	1.90	0.53
1:A:39:ASN:ND2	1:A:77:ARG:HH11	2.06	0.53
1:B:176:HIS:HE1	3:B:509:HOH:O	1.90	0.53
1:B:268:GLU:HB2	3:B:416:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:272[A]:GLU:HG3	1:B:274:SER:H	1.72	0.53
1:A:75[A]:ARG:HH21	1:A:112:LEU:HB2	1.74	0.53
1:B:136:PRO:HD2	1:B:139:ASP:HB2	1.91	0.53
1:A:119:VAL:HG22	1:A:122[A]:ARG:NH1	2.25	0.52
1:A:141:GLU:OE2	1:A:142:ASN:ND2	2.43	0.52
1:A:119:VAL:HG22	1:A:122[A]:ARG:HH11	1.74	0.52
1:B:26:HIS:HD2	1:B:43:THR:OG1	1.92	0.52
1:B:39:ASN:ND2	1:B:77:ARG:HH11	2.06	0.52
1:A:147:ILE:HD12	1:A:194:THR:HG22	1.92	0.51
1:B:210:HIS:HB2	3:B:632:HOH:O	2.10	0.51
1:B:124:TRP:CE3	1:B:172:VAL:HG13	2.46	0.51
1:B:125:ARG:HH11	1:B:132:GLN:NE2	2.08	0.51
1:B:339[B]:HIS:HD2	3:B:403:HOH:O	1.94	0.51
1:A:75[A]:ARG:HG2	1:A:75[A]:ARG:HH11	1.75	0.50
1:A:120:LEU:HD12	1:A:123:ARG:HG3	1.93	0.50
1:A:121:MET:SD	1:A:172:VAL:HG21	2.52	0.50
1:A:154:LEU:N	1:A:155:PRO:CD	2.75	0.50
1:B:134:TRP:CD1	2:D:1:NDG:H8C2	2.47	0.50
1:A:147:ILE:HD12	1:A:194:THR:CG2	2.42	0.50
1:B:339[A]:HIS:CE1	2:D:1:NDG:H6C2	2.45	0.49
1:B:136:PRO:HD2	1:B:139:ASP:OD2	2.11	0.49
1:A:149:ASP:OD2	1:A:225:TRP:NE1	2.46	0.49
1:B:87:HIS:H	1:B:87:HIS:HD1	1.59	0.48
1:A:86:HIS:HD2	1:A:88:ASN:N	2.09	0.48
1:B:302:GLU:HG2	1:B:306:PHE:CE2	2.49	0.48
1:B:179:LYS:HE2	1:B:197:PHE:CZ	2.49	0.48
1:B:377:ARG:HG2	1:B:377:ARG:NH2	2.28	0.47
1:B:65:GLU:O	1:B:69[B]:ARG:HG3	2.13	0.47
1:A:149:ASP:OD1	1:A:149:ASP:N	2.47	0.47
1:B:128:ALA:O	1:B:130:ILE:HG23	2.14	0.47
1:B:158:LEU:O	1:B:162:GLU:HG3	2.14	0.47
1:B:149:ASP:OD2	1:B:149:ASP:N	2.48	0.47
1:A:26:HIS:HD2	1:A:43:THR:OG1	1.98	0.46
1:B:254:LEU:HD12	1:B:305[A]:ARG:HH22	1.80	0.46
1:B:305[A]:ARG:NH2	1:B:306:PHE:CE1	2.83	0.46
1:B:212:GLY:O	1:B:341:ARG:HD2	2.16	0.46
1:B:254:LEU:HD12	1:B:305[A]:ARG:NH2	2.30	0.45
1:A:75[A]:ARG:NH2	3:A:411:HOH:O	2.49	0.45
1:B:132:GLN:HG3	1:B:142:ASN:O	2.17	0.45
1:A:188:ASP:O	1:A:189:ASP:HB2	2.16	0.45
1:B:234:SER:OG	1:B:243:LEU:HB2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:147:ILE:HG22	1:A:149:ASP:OD1	2.16	0.45
1:A:264:TYR:HA	1:A:277:ARG:HA	1.99	0.45
1:B:130:ILE:CG2	1:B:179:LYS:HD3	2.47	0.44
1:A:61[B]:GLU:OE2	1:A:65:GLU:HG2	2.17	0.44
1:A:83:ASN:HB3	3:A:732:HOH:O	2.17	0.44
1:A:86:HIS:CD2	1:A:88:ASN:H	2.28	0.44
1:A:374:GLU:HB2	1:A:377:ARG:HB2	1.99	0.44
1:B:23[A]:ARG:HH21	1:B:23[A]:ARG:HG3	1.83	0.44
1:A:119:VAL:HG13	1:A:122[A]:ARG:NH1	2.32	0.43
1:A:150:CYS:HB3	3:A:572:HOH:O	2.19	0.43
1:B:52:TRP:CE2	1:B:67:ALA:HB1	2.54	0.43
1:A:234[B]:SER:OG	1:A:243:LEU:HB2	2.18	0.43
2:C:1:NDG:H2	2:C:1:NDG:H8C1	1.70	0.42
1:A:87:HIS:O	1:A:153:ASN:CG	2.62	0.42
1:A:107[B]:GLU:OE2	1:A:110:ARG:HD3	2.20	0.42
1:A:87:HIS:ND1	1:A:150:CYS:HB2	2.33	0.42
1:B:23[A]:ARG:HE	1:B:36:LEU:HD12	1.85	0.42
1:B:88:ASN:ND2	2:D:1:NDG:O7	2.53	0.42
1:A:247:LYS:O	1:A:251[A]:ARG:HG3	2.20	0.42
1:B:166:ASP:HA	1:B:167:PRO:HD2	1.91	0.42
1:A:19:LYS:NZ	3:A:414:HOH:O	2.52	0.42
1:B:39:ASN:HA	1:B:44:ASP:OD1	2.20	0.42
1:B:66:GLY:HA2	1:B:69[B]:ARG:HD2	2.01	0.42
1:A:124:TRP:HA	1:A:131:ILE:HD13	2.02	0.41
1:A:162:GLU:HB2	3:A:647:HOH:O	2.20	0.41
3:A:431:HOH:O	2:C:2:GAD:H2	2.19	0.41
1:B:125:ARG:C	1:B:127:ASP:N	2.77	0.41
1:B:125:ARG:NH1	1:B:132:GLN:NE2	2.67	0.41
1:A:187:GLY:HA3	1:B:185:VAL:HB	2.01	0.41
1:A:179:LYS:HE2	1:A:197:PHE:CE2	2.55	0.41
1:B:176:HIS:CE1	3:B:509:HOH:O	2.71	0.41
1:B:32:ASN:ND2	1:B:328:GLU:HG2	2.36	0.41
1:B:298:GLU:HG3	1:B:303[A]:ARG:NH1	2.36	0.41
1:A:145:ARG:NH2	1:A:194:THR:HG21	2.35	0.40
1:B:305[A]:ARG:NH2	1:B:306:PHE:HE1	2.19	0.40
1:B:32:ASN:HD21	1:B:328:GLU:HG2	1.86	0.40
1:B:154:LEU:N	1:B:155:PRO:CD	2.84	0.40
1:B:124:TRP:CD2	1:B:172:VAL:HG13	2.56	0.40
1:B:300:ASP:HA	1:B:301:PRO:HD2	1.97	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	385/377 (102%)	372 (97%)	13 (3%)	0	100	100
1	B	382/377 (101%)	360 (94%)	19 (5%)	3 (1%)	16	4
All	All	767/754 (102%)	732 (95%)	32 (4%)	3 (0%)	30	16

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	136	PRO
1	B	88	ASN
1	B	141	GLU

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	314/304 (103%)	307 (98%)	7 (2%)	45	22
1	B	311/304 (102%)	306 (98%)	5 (2%)	55	35
All	All	625/608 (103%)	613 (98%)	12 (2%)	57	28

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	61[A]	GLU
1	A	61[B]	GLU
1	A	118	ASP

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Mol	Chain	Res	Type
1	A	141	GLU
1	A	149	ASP
1	A	211[A]	GLN
1	A	211[B]	GLN
1	B	65	GLU
1	B	121	MET
1	B	149	ASP
1	B	272[A]	GLU
1	B	272[B]	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	HIS
1	A	39	ASN
1	A	83	ASN
1	A	86	HIS
1	A	142	ASN
1	A	176	HIS
1	A	193	HIS
1	A	223	GLN
1	A	304	GLN
1	B	16	ASN
1	B	26	HIS
1	B	39	ASN
1	B	86	HIS
1	B	100	GLN
1	B	132	GLN
1	B	176	HIS
1	B	210	HIS
1	B	223	GLN

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

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	NDG	C	1	2	15,15,15	0.63	0	21,21,21	0.77	0
2	GAD	C	2	2	10,11,11	1.69	2 (20%)	12,15,15	1.55	3 (25%)
2	NDG	D	1	2	15,15,15	0.62	0	21,21,21	0.73	0
2	GAD	D	2	2	10,11,11	1.71	2 (20%)	12,15,15	1.68	3 (25%)

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	NDG	C	1	2	-	3/6/26/26	0/1/1/1
2	GAD	C	2	2	-	4/4/17/17	0/1/1/1
2	NDG	D	1	2	-	2/6/26/26	0/1/1/1
2	GAD	D	2	2	-	0/4/17/17	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	GAD	C4-C5	3.77	1.39	1.33
2	C	2	GAD	C4-C5	3.69	1.39	1.33
2	D	2	GAD	C3-C4	2.24	1.53	1.50
2	C	2	GAD	C3-C4	2.22	1.53	1.50

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	GAD	O5-C5-C6	3.97	119.15	111.85
2	C	2	GAD	O5-C5-C6	3.64	118.54	111.85
2	D	2	GAD	O5-C5-C4	-2.60	122.55	124.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	GAD	C4-C5-C6	-2.43	118.29	123.56
2	C	2	GAD	C4-C5-C6	-2.32	118.53	123.56
2	C	2	GAD	O5-C5-C4	-2.18	122.93	124.94

There are no chirality outliers.

All (9) torsion outliers are listed below:

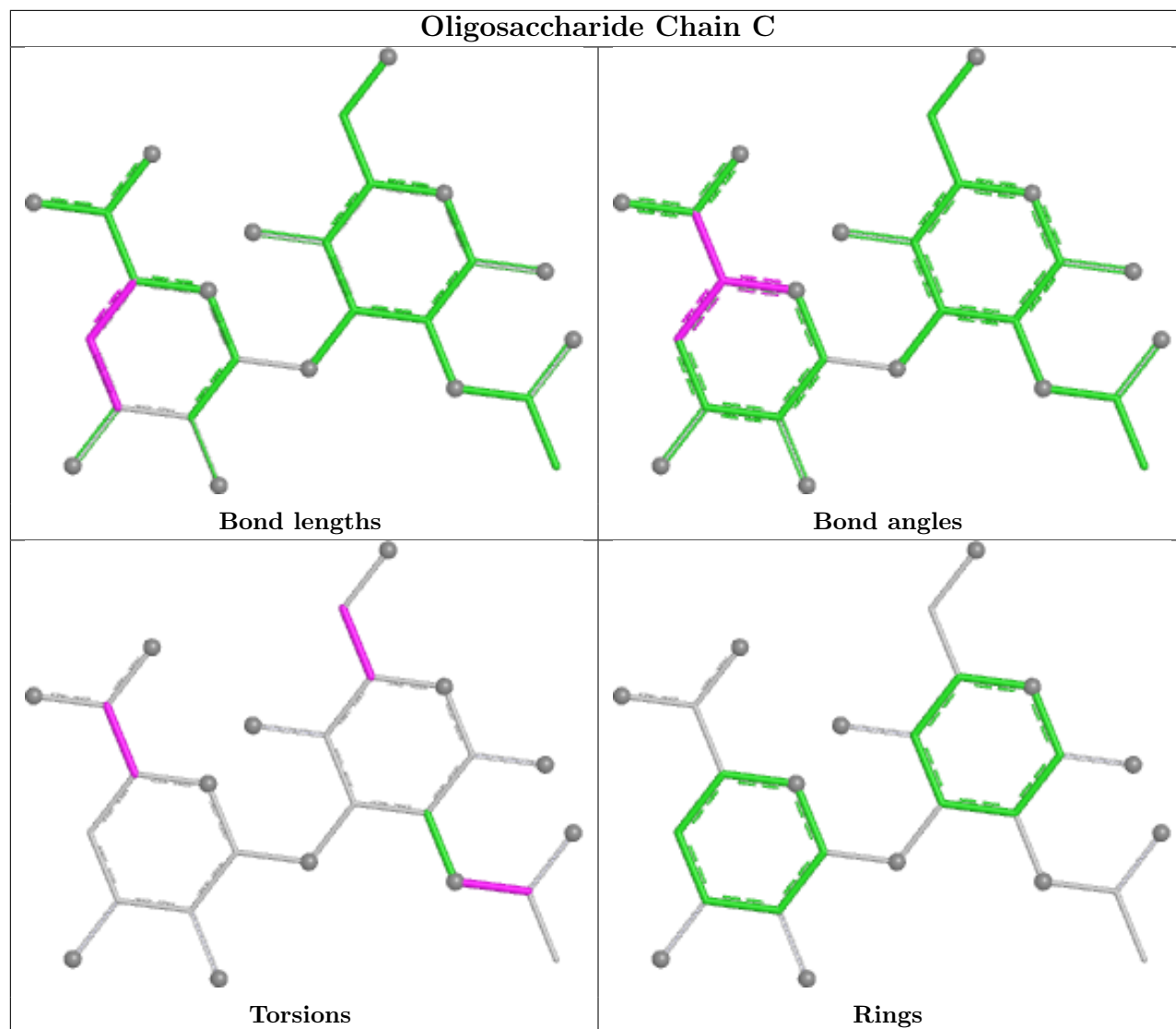
Mol	Chain	Res	Type	Atoms
2	C	1	NDG	C8-C7-N2-C2
2	C	1	NDG	O7-C7-N2-C2
2	D	1	NDG	C8-C7-N2-C2
2	D	1	NDG	O7-C7-N2-C2
2	C	1	NDG	O5-C5-C6-O6
2	C	2	GAD	C4-C5-C6-O6A
2	C	2	GAD	C4-C5-C6-O6B
2	C	2	GAD	O5-C5-C6-O6A
2	C	2	GAD	O5-C5-C6-O6B

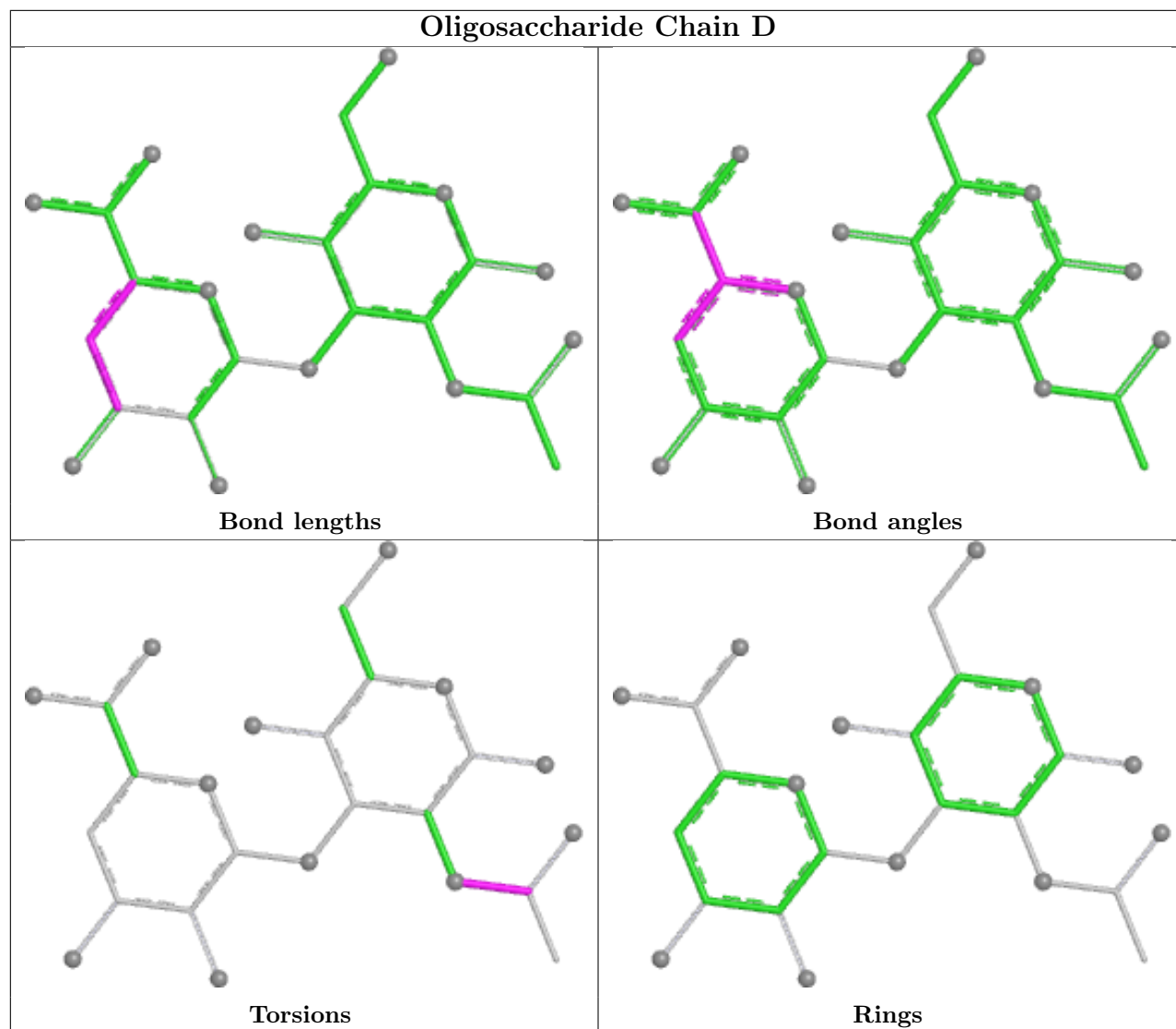
There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	NDG	4	0
2	C	1	NDG	1	0
2	C	2	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](#)

There are no ligands in this entry.

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	377/377 (100%)	-0.21	13 (3%)	48 56	7, 14, 31, 46	10 (2%)
1	B	377/377 (100%)	0.18	35 (9%)	14 18	7, 17, 45, 52	7 (1%)
All	All	754/754 (100%)	-0.02	48 (6%)	25 31	7, 15, 41, 52	17 (2%)

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	134	TRP	6.4
1	A	134	TRP	6.1
1	B	136	PRO	5.9
1	B	131	ILE	5.8
1	B	124	TRP	5.7
1	B	135	GLY	5.5
1	A	376	GLY	5.3
1	B	140	PRO	5.1
1	A	136	PRO	4.8
1	B	126	ALA	4.5
1	B	87	HIS	4.5
1	B	133	ALA	4.2
1	B	129	GLY	4.0
1	B	138	GLY	3.9
1	A	140	PRO	3.9
1	B	137	LYS	3.9
1	B	128	ALA	3.8
1	B	141	GLU	3.7
1	B	376	GLY	3.7
1	B	86	HIS	3.5
1	B	127	ASP	3.4
1	A	149	ASP	3.3
1	A	377	ARG	3.3
1	B	139	ASP	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	210	HIS	3.2
1	B	142	ASN	3.1
1	B	143	GLY	3.1
1	B	123	ARG	3.1
1	A	141	GLU	3.1
1	B	85	ASP	3.1
1	B	199	PRO	2.8
1	B	144	GLY	2.7
1	B	122	ARG	2.7
1	A	135	GLY	2.6
1	B	132	GLN	2.5
1	B	200	GLU	2.3
1	A	123	ARG	2.3
1	B	88	ASN	2.2
1	B	81	PHE	2.2
1	A	124	TRP	2.1
1	A	139	ASP	2.1
1	B	130	ILE	2.1
1	B	298	GLU	2.0
1	B	121	MET	2.0
1	A	122[A]	ARG	2.0
1	A	150	CYS	2.0
1	B	202	GLY	2.0
1	B	171	ARG	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](#)

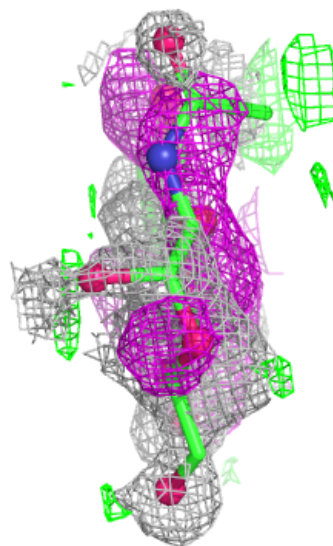
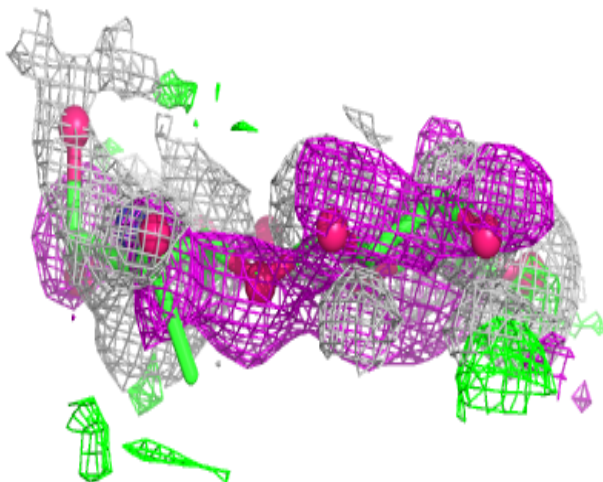
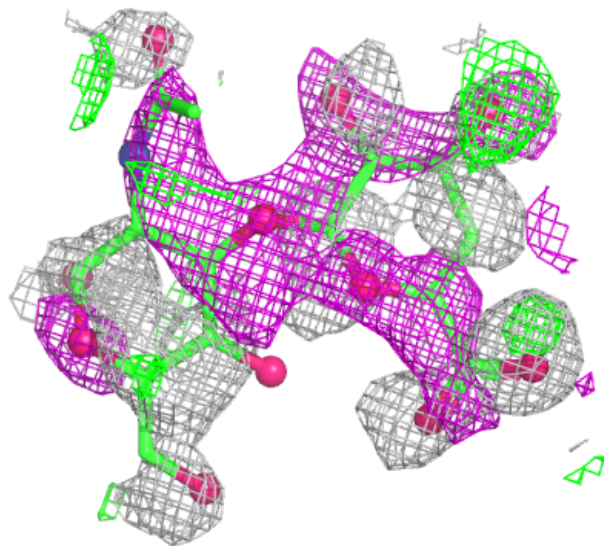
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	NDG	C	1	15/15	0.34	0.18	46,50,54,55	0
2	NDG	D	1	15/15	0.37	0.15	49,50,55,55	0
2	GAD	C	2	11/11	0.70	0.20	17,43,54,56	0
2	GAD	D	2	11/11	0.70	0.20	18,44,54,57	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.

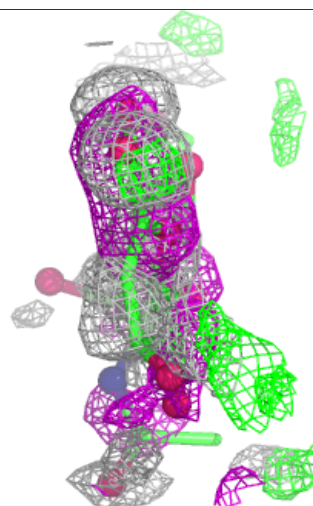
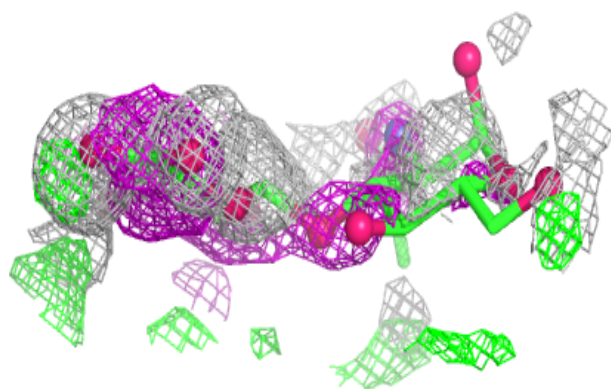
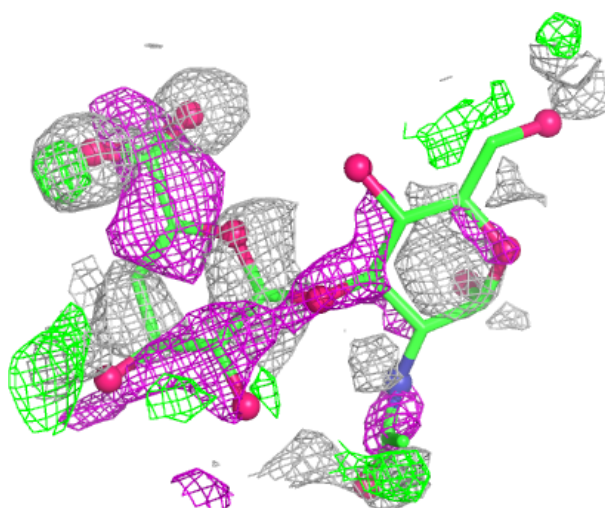
Electron density around Chain C:

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.