



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 04:15 AM UTC

PDB ID : 2NY2 / pdb_00002ny2
Title : HIV-1 gp120 Envelope Glycoprotein (T123C, T257S, S334A, S375W, G431C)
Complexed with CD4 and Antibody 17b
Authors : Zhou, T.; Xu, L.; Dey, B.; Hessel, A.J.; Van Ryk, D.; Xiang, S.H.; Yang, X.;
Zhang, M.Y.; Zwick, M.B.; Arthos, J.; Burton, D.R.; Dimitrov, D.S.; Sodroski,
J.; Wyatt, R.; Nabel, G.J.; Kwong, P.D.
Deposited on : 2006-11-20
Resolution : 2.00 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

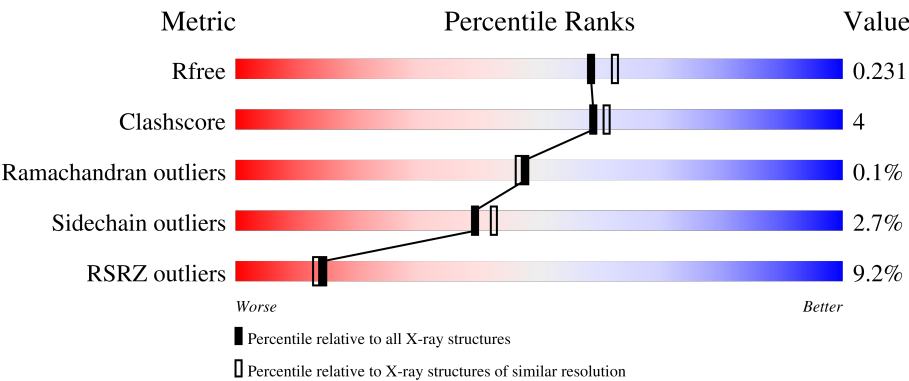
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	317	11% 85% 9% ..
2	B	184	16% 89% 9% ..
3	C	214	6% 89% 10% .
4	D	229	3% 86% 10% .

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Mol	Chain	Length	Quality of chain
5	E	2	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	NAG	A	734	-	-	-	X
6	NAG	A	741	-	-	-	X
6	NAG	A	762	-	-	-	X
6	NAG	A	776	-	-	-	X
6	NAG	A	789	-	-	-	X
6	NAG	A	795	-	-	-	X
6	NAG	A	839	-	-	-	X
6	NAG	A	886	-	-	-	X
6	NAG	A	892	-	-	-	X
6	NAG	A	948	-	-	-	X
8	HEZ	A	4003	-	-	-	X

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 7945 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 ENVELOPE GLYCOPROTEIN GP120.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	304	Total	C	N	O	S	0	0	0
			2358	1482	411	443	22			

- Molecule 2 is a protein called T-cell surface glycoprotein CD4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	181	Total	C	N	O	S	0	0	0
			1412	885	247	276	4			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	1000	MET	-	initiating methionine	UNP P01730

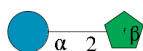
- Molecule 3 is a protein called ANTIBODY 17B, LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	214	Total	C	N	O	S	0	0	0
			1646	1028	282	331	5			

- Molecule 4 is a protein called ANTIBODY 17B, HEAVY CHAIN.

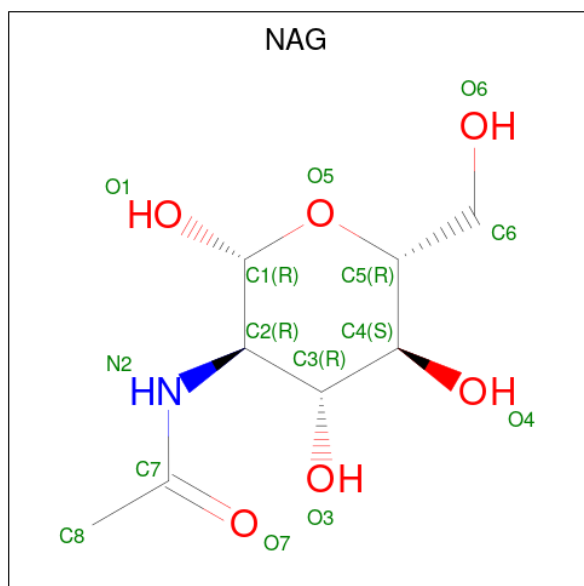
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	221	Total	C	N	O	S	0	1	0
			1671	1057	280	329	5			

- Molecule 5 is an oligosaccharide called beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose.



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

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



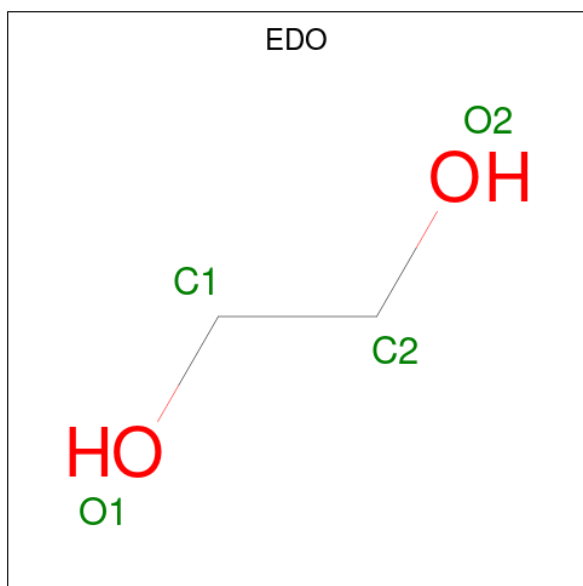
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			14	8	1	5		
6	A	1	Total	C	N	O	0	0
			14	8	1	5		
6	A	1	Total	C	N	O	0	0
			14	8	1	5		
6	A	1	Total	C	N	O	0	0
			14	8	1	5		
6	A	1	Total	C	N	O	0	0
			14	8	1	5		
6	A	1	Total	C	N	O	0	0
			14	8	1	5		
6	A	1	Total	C	N	O	0	0
			14	8	1	5		
6	A	1	Total	C	N	O	0	0
			14	8	1	5		

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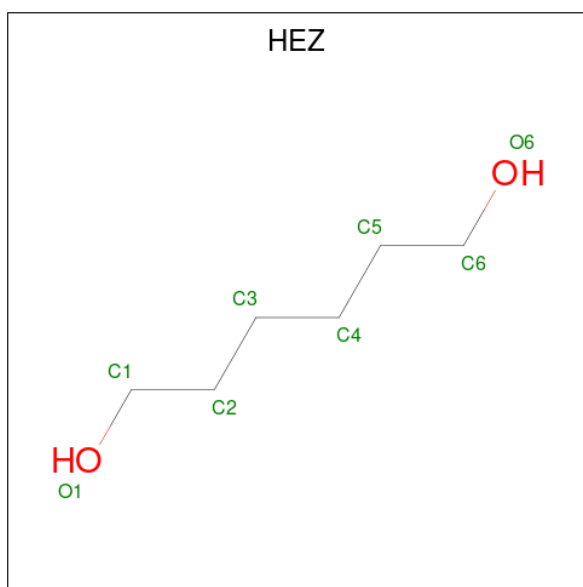
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 7 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			4	2	2		
7	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 8 is HEXANE-1,6-DIOL (CCD ID: HEZ) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			8	6	2		

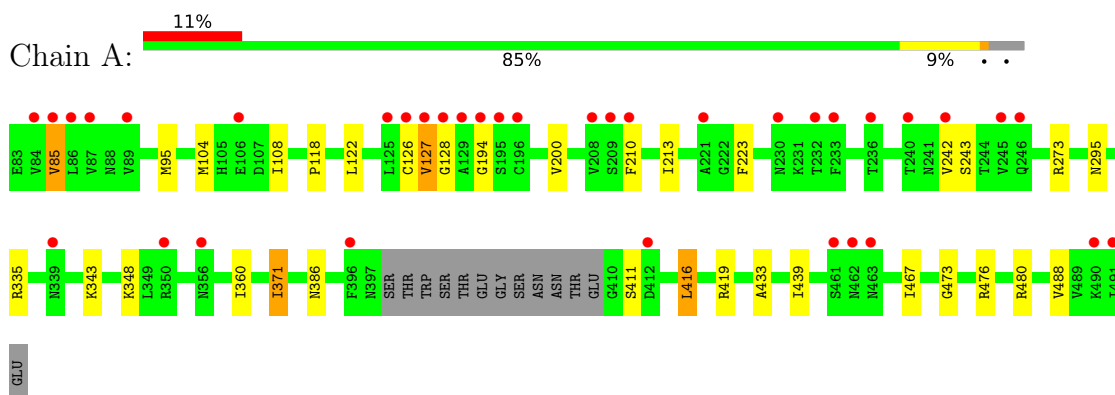
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	204	Total	O	0	0
			204	204		
9	B	98	Total	O	0	0
			98	98		
9	C	164	Total	O	0	0
			164	164		
9	D	199	Total	O	0	0
			199	199		

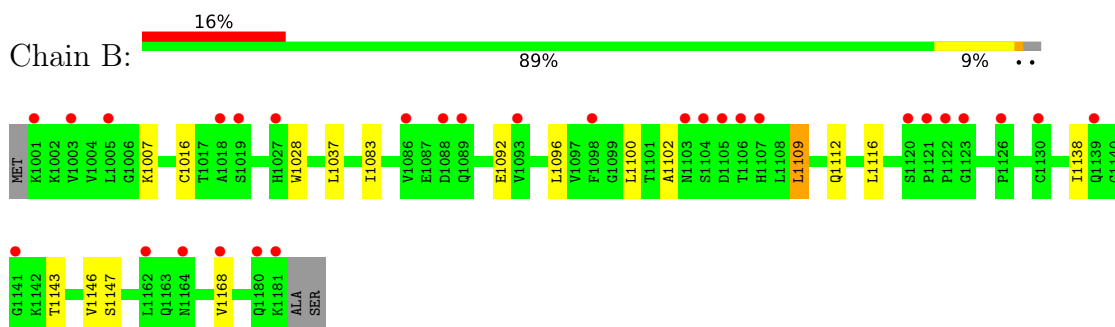
3 Residue-property plots [i](#)

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.

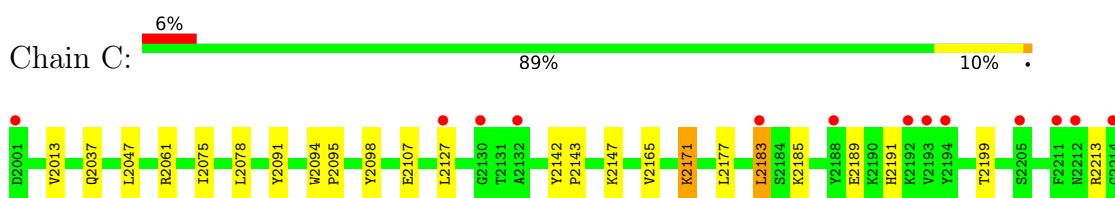
- Molecule 1: ENVELOPE GLYCOPROTEIN GP120



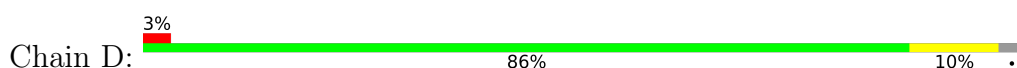
- Molecule 2: T-cell surface glycoprotein CD4

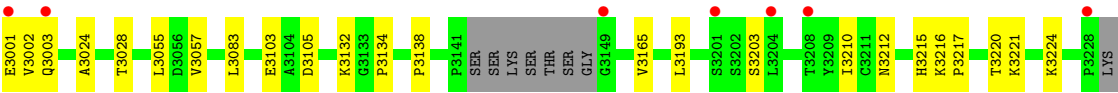


- Molecule 3: ANTIBODY 17B, LIGHT CHAIN



- Molecule 4: ANTIBODY 17B, HEAVY CHAIN





- Molecule 5: beta-D-fructofuranose-(2-1)-alpha-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	71.61Å 88.02Å 196.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.84 – 2.00 42.84 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.6 (42.84-2.00) 94.6 (42.84-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.08 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.194 , 0.223 0.209 , 0.231	Depositor DCC
R_{free} test set	4062 reflections (4.81%)	wwPDB-VP
Wilson B-factor (Å ²)	28.7	Xtriage
Anisotropy	0.082	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 56.4	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	0.94	EDS
Total number of atoms	7945	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.37% 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: FRU, GLC, NAG, EDO, HEZ

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	0.36	0/2406	0.67	0/3263
2	B	0.32	0/1432	0.63	0/1930
3	C	0.34	0/1683	0.68	0/2288
4	D	0.39	0/1716	0.69	0/2339
All	All	0.36	0/7237	0.67	0/9820

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	2358	0	2304	24	0
2	B	1412	0	1441	7	0
3	C	1646	0	1590	15	0
4	D	1671	0	1631	16	0
5	E	23	0	21	1	0
6	A	154	0	143	5	0
7	A	4	0	6	0	0
7	D	4	0	6	0	0
8	A	8	0	14	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	A	204	0	0	5	1
9	B	98	0	0	0	0
9	C	164	0	0	2	0
9	D	199	0	0	6	0
All	All	7945	0	7156	63	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (63) 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)
4:D:3212:ASN:ND2	9:D:240:HOH:O	1.89	1.00
3:C:2191:HIS:O	3:C:2213:ARG:NH1	1.97	0.98
4:D:3215:HIS:HB3	4:D:3220:THR:HG22	1.54	0.89
4:D:3210:ILE:HG22	9:D:240:HOH:O	1.83	0.79
3:C:2185:LYS:HE2	3:C:2189:GLU:OE2	1.89	0.72
4:D:3215:HIS:HB3	4:D:3220:THR:CG2	2.20	0.72
6:A:839:NAG:O7	9:A:4205:HOH:O	2.10	0.69
1:A:122:LEU:HD22	1:A:200:VAL:HG22	1.75	0.67
1:A:122:LEU:HD21	4:D:3057:VAL:HG21	1.76	0.67
1:A:411:SER:OG	9:A:4123:HOH:O	2.14	0.65
1:A:371:ILE:HD11	1:A:473:GLY:CA	2.29	0.63
1:A:419:ARG:NH1	9:A:4079:HOH:O	2.32	0.63
2:B:1138:ILE:HG21	2:B:1146:VAL:HG22	1.82	0.60
4:D:3134:PRO:HD2	4:D:3220:THR:HG21	1.83	0.60
2:B:1083:ILE:HG13	2:B:1092:GLU:HG3	1.83	0.59
3:C:2094:TRP:HA	3:C:2095:PRO:C	2.29	0.57
1:A:419:ARG:NH1	4:D:3103:GLU:OE2	2.33	0.57
3:C:2183:LEU:N	3:C:2183:LEU:HD23	2.20	0.55
3:C:2171:LYS:HA	3:C:2171:LYS:HE3	1.87	0.55
1:A:386:ASN:O	1:A:416:LEU:HG	2.07	0.54
1:A:335:ARG:NH2	9:A:4103:HOH:O	2.21	0.54
1:A:371:ILE:HD11	1:A:473:GLY:HA3	1.90	0.54
1:A:223:PHE:HB3	1:A:488:VAL:HG13	1.91	0.52
3:C:2147:LYS:HB3	3:C:2199:THR:HB	1.91	0.51
3:C:2165:VAL:HG22	3:C:2177:LEU:HD12	1.93	0.51
4:D:3002:VAL:HG21	4:D:3024:ALA:HB1	1.94	0.50
1:A:128:GLY:C	1:A:194:GLY:H	2.20	0.50
3:C:2091:TYR:HA	3:C:2098:TYR:CD1	2.47	0.49
1:A:210:PHE:HB3	9:A:4179:HOH:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:CYS:O	1:A:127:VAL:HB	2.14	0.47
1:A:360:ILE:HB	1:A:467:ILE:HD13	1.96	0.46
9:D:45:HOH:O	5:E:2:FRU:O1	2.18	0.46
1:A:104:MET:O	1:A:108:ILE:HG12	2.16	0.46
3:C:2075:ILE:HG22	9:C:340:HOH:O	2.15	0.46
4:D:3105:ASP:OD1	4:D:3105:ASP:N	2.49	0.46
3:C:2037:GLN:HB2	3:C:2047:LEU:HD11	1.98	0.46
1:A:371:ILE:HD11	1:A:473:GLY:HA2	1.95	0.45
1:A:295:ASN:HD22	6:A:795:NAG:H83	1.81	0.45
3:C:2061:ARG:HB2	9:C:340:HOH:O	2.15	0.45
1:A:343:LYS:HE3	6:A:839:NAG:H82	1.97	0.45
6:A:839:NAG:O7	6:A:839:NAG:C1	2.65	0.45
4:D:3193:LEU:C	4:D:3193:LEU:HD12	2.42	0.45
3:C:2183:LEU:HD23	3:C:2183:LEU:H	1.82	0.44
4:D:3216:LYS:N	4:D:3217:PRO:CD	2.81	0.44
1:A:85:VAL:HA	1:A:243:SER:HB3	2.00	0.44
2:B:1102:ALA:HA	2:B:1116:LEU:HD23	2.01	0.43
1:A:343:LYS:HE3	6:A:839:NAG:C8	2.49	0.43
2:B:1109:LEU:HB2	2:B:1112:GLN:CD	2.44	0.42
2:B:1007:LYS:HG3	2:B:1168:VAL:HG21	2.01	0.42
3:C:2142:TYR:CG	3:C:2143:PRO:HA	2.55	0.42
3:C:2183:LEU:N	3:C:2183:LEU:CD2	2.83	0.42
1:A:95:MET:SD	1:A:273:ARG:HD3	2.60	0.41
1:A:348:LYS:HD3	1:A:348:LYS:HA	1.90	0.41
1:A:476:ARG:O	1:A:480:ARG:HB2	2.20	0.41
2:B:1016:CYS:HB2	2:B:1028:TRP:CZ2	2.56	0.41
4:D:3003:GLN:NE2	9:D:279:HOH:O	2.54	0.41
4:D:3003:GLN:NE2	9:D:372:HOH:O	2.53	0.41
4:D:3132:LYS:HE3	4:D:3132:LYS:HB2	1.95	0.41
1:A:118:PRO:HB3	1:A:433:ALA:HB1	2.03	0.41
4:D:3138:PRO:HD3	4:D:3224:LYS:HE2	2.03	0.41
2:B:1138:ILE:HD13	2:B:1146:VAL:HG23	2.03	0.40
3:C:2013:VAL:CG1	3:C:2078:LEU:HD12	2.52	0.40
4:D:3221:LYS:NZ	9:D:614:HOH:O	2.53	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:A:4171:HOH:O	9:A:4171:HOH:O[4_556]	1.30	0.90

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	300/317 (95%)	286 (95%)	13 (4%)	1 (0%)	36	35
2	B	179/184 (97%)	173 (97%)	6 (3%)	0	100	100
3	C	212/214 (99%)	205 (97%)	7 (3%)	0	100	100
4	D	218/229 (95%)	214 (98%)	4 (2%)	0	100	100
All	All	909/944 (96%)	878 (97%)	30 (3%)	1 (0%)	48	46

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	127	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	269/281 (96%)	263 (98%)	6 (2%)	45	50
2	B	164/166 (99%)	158 (96%)	6 (4%)	30	30
3	C	184/184 (100%)	180 (98%)	4 (2%)	45	50
4	D	187/193 (97%)	181 (97%)	6 (3%)	34	35
All	All	804/824 (98%)	782 (97%)	22 (3%)	39	42

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	85	VAL
1	A	213	ILE
1	A	242	VAL
1	A	371	ILE
1	A	416	LEU
1	A	439	ILE
2	B	1037	LEU
2	B	1096	LEU
2	B	1100	LEU
2	B	1109	LEU
2	B	1143	THR
2	B	1147	SER
3	C	2107	GLU
3	C	2127	LEU
3	C	2171	LYS
3	C	2183	LEU
4	D	3001	GLU
4	D	3028	THR
4	D	3055	LEU
4	D	3083	LEU
4	D	3165	VAL
4	D	3203	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	ASN
1	A	352	GLN
2	B	1025	GLN
2	B	1064	GLN
2	B	1073	ASN
2	B	1163	GLN
2	B	1164	ASN
3	C	2038	GLN
3	C	2102	GLN
3	C	2212	ASN
4	D	3039	GLN
4	D	3085	ASN

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 ⓘ

2 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	$\# Z > 2$
5	GLC	E	1	5	11,11,12	0.60	0	15,15,17	0.57	0
5	FRU	E	2	5	11,12,12	0.47	0	10,18,18	0.62	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
5	GLC	E	1	5	-	1/2/19/22	0/1/1/1
5	FRU	E	2	5	-	5/5/24/24	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

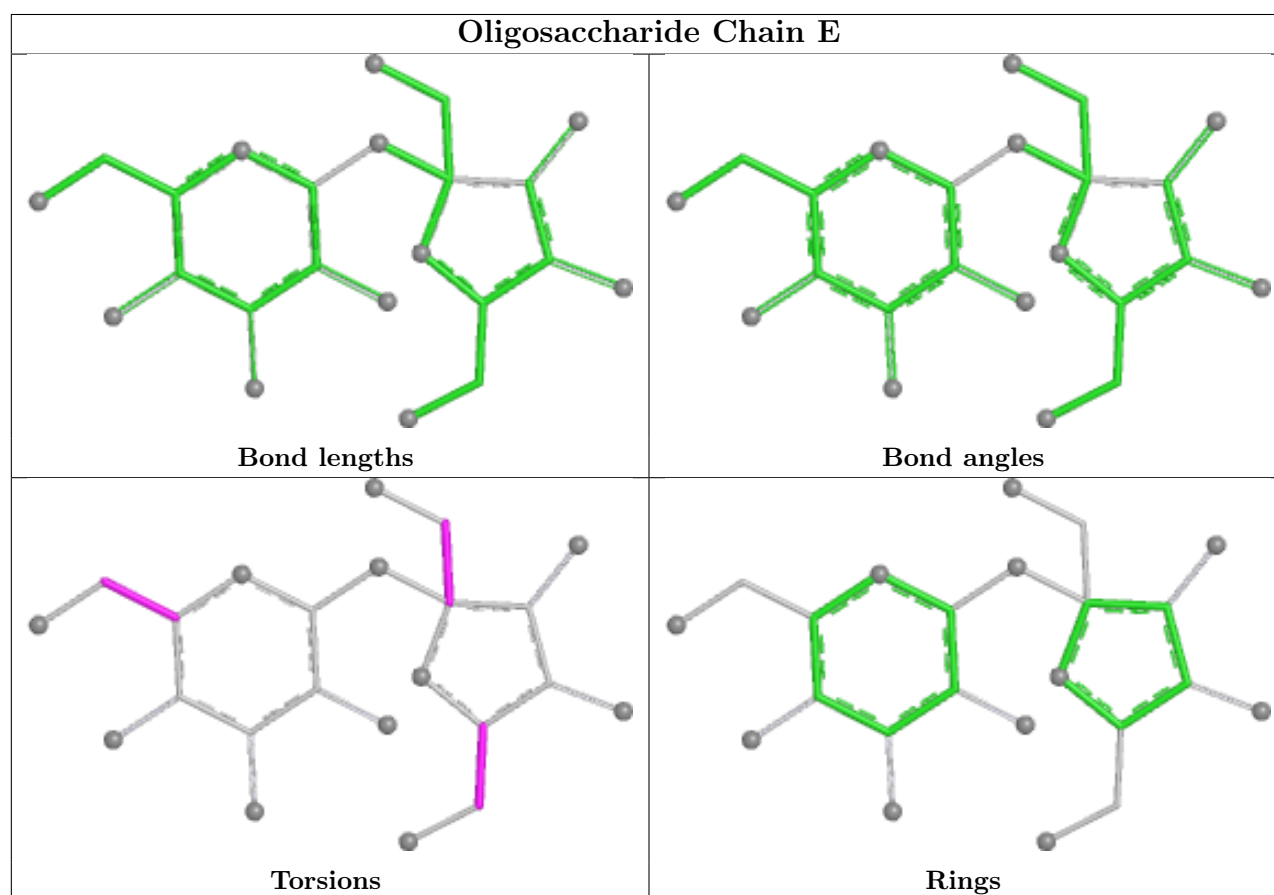
Mol	Chain	Res	Type	Atoms
5	E	2	FRU	O1-C1-C2-C3
5	E	2	FRU	O1-C1-C2-O2
5	E	2	FRU	O1-C1-C2-O5
5	E	2	FRU	C4-C5-C6-O6
5	E	2	FRU	O5-C5-C6-O6
5	E	1	GLC	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	E	2	FRU	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](#)

14 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
6	NAG	A	776	1	14,14,15	0.53	0	17,19,21	0.76	0
6	NAG	A	734	1	14,14,15	0.49	0	17,19,21	0.83	0
6	NAG	A	839	1	14,14,15	0.62	0	17,19,21	0.88	1 (5%)
7	EDO	D	4002	-	3,3,3	0.41	0	2,2,2	0.41	0
8	HEZ	A	4003	-	7,7,7	0.27	0	6,6,6	0.51	0
6	NAG	A	730	1	14,14,15	0.59	0	17,19,21	0.79	0
6	NAG	A	892	1	14,14,15	0.51	0	17,19,21	0.69	0
6	NAG	A	762	1	14,14,15	0.57	0	17,19,21	0.78	0
6	NAG	A	789	1	14,14,15	0.54	0	17,19,21	0.83	0
6	NAG	A	886	1	14,14,15	0.54	0	17,19,21	0.69	0
6	NAG	A	948	1	14,14,15	0.53	0	17,19,21	0.76	0
6	NAG	A	795	1	14,14,15	0.48	0	17,19,21	1.08	1 (5%)
6	NAG	A	741	1	14,14,15	0.50	0	17,19,21	0.84	1 (5%)
7	EDO	A	4001	-	3,3,3	0.39	0	2,2,2	0.41	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
6	NAG	A	776	1	-	0/6/23/26	0/1/1/1
6	NAG	A	734	1	-	0/6/23/26	0/1/1/1
6	NAG	A	839	1	-	2/6/23/26	0/1/1/1
7	EDO	D	4002	-	-	0/1/1/1	-
8	HEZ	A	4003	-	-	2/5/5/5	-
6	NAG	A	730	1	-	0/6/23/26	0/1/1/1
6	NAG	A	892	1	-	2/6/23/26	0/1/1/1
6	NAG	A	762	1	-	2/6/23/26	0/1/1/1
6	NAG	A	789	1	-	0/6/23/26	0/1/1/1
6	NAG	A	886	1	-	2/6/23/26	0/1/1/1
6	NAG	A	948	1	-	0/6/23/26	0/1/1/1
6	NAG	A	795	1	-	2/6/23/26	0/1/1/1
6	NAG	A	741	1	-	2/6/23/26	0/1/1/1
7	EDO	A	4001	-	-	1/1/1/1	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	795	NAG	C2-N2-C7	2.46	126.20	122.90
6	A	741	NAG	C2-N2-C7	-2.05	120.15	122.90
6	A	839	NAG	O5-C1-C2	2.04	114.45	111.29

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	839	NAG	C1-C2-N2-C7
6	A	892	NAG	O5-C5-C6-O6
6	A	886	NAG	C8-C7-N2-C2
6	A	795	NAG	C8-C7-N2-C2
6	A	795	NAG	O7-C7-N2-C2
6	A	886	NAG	O7-C7-N2-C2
6	A	892	NAG	C4-C5-C6-O6
6	A	741	NAG	C4-C5-C6-O6
8	A	4003	HEZ	C3-C4-C5-C6
6	A	741	NAG	O5-C5-C6-O6
8	A	4003	HEZ	O1-C1-C2-C3
6	A	762	NAG	C4-C5-C6-O6
6	A	762	NAG	O5-C5-C6-O6
7	A	4001	EDO	O1-C1-C2-O2
6	A	839	NAG	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	839	NAG	4	0
6	A	795	NAG	1	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	304/317 (95%)	0.57	36 (11%) 9 8	11, 34, 72, 122	0
2	B	181/184 (98%)	0.87	29 (16%) 5 4	18, 46, 78, 90	0
3	C	214/214 (100%)	0.22	13 (6%) 27 26	9, 30, 70, 83	0
4	D	221/229 (96%)	-0.04	7 (3%) 50 49	7, 20, 61, 84	1 (0%)
All	All	920/944 (97%)	0.40	85 (9%) 14 13	7, 32, 73, 122	1 (0%)

All (85) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	129	ALA	7.3
1	A	210	PHE	4.6
1	A	127	VAL	3.8
2	B	1088	ASP	3.8
1	A	125	LEU	3.8
1	A	128	GLY	3.8
3	C	2192	LYS	3.7
2	B	1105	ASP	3.6
2	B	1181	LYS	3.5
2	B	1104	SER	3.4
2	B	1106	THR	3.4
2	B	1162	LEU	3.4
1	A	490	LYS	3.3
2	B	1123	GLY	3.3
1	A	195	SER	3.2
2	B	1121	PRO	3.2
1	A	462	ASN	3.2
1	A	491	ILE	3.1
1	A	232	THR	3.0
3	C	2183	LEU	3.0
1	A	126	CYS	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	87	VAL	3.0
1	A	396	PHE	3.0
1	A	86	LEU	2.9
3	C	2188	TYR	2.9
1	A	85	VAL	2.9
4	D	3001	GLU	2.9
1	A	194	GLY	2.9
1	A	196	CYS	2.9
3	C	2214	GLY	2.9
4	D	3204	LEU	2.8
2	B	1089	GLN	2.7
2	B	1120	SER	2.7
3	C	2001	ASP	2.7
2	B	1098	PHE	2.7
1	A	230	ASN	2.7
1	A	84	VAL	2.7
1	A	221	ALA	2.6
1	A	350	ARG	2.6
1	A	356	ASN	2.6
4	D	3149	GLY	2.6
2	B	1019	SER	2.6
1	A	89	VAL	2.5
2	B	1003	VAL	2.5
3	C	2193	VAL	2.5
1	A	209	SER	2.5
4	D	3228	PRO	2.5
2	B	1141	GLY	2.5
3	C	2212	ASN	2.5
4	D	3201	SER	2.4
2	B	1103	ASN	2.4
2	B	1139	GLN	2.4
3	C	2211	PHE	2.4
1	A	242	VAL	2.4
3	C	2194	TYR	2.4
2	B	1168	VAL	2.4
4	D	3003	GLN	2.3
1	A	106	GLU	2.3
1	A	246	GLN	2.3
1	A	245	VAL	2.2
2	B	1093	VAL	2.2
1	A	412	ASP	2.2
4	D	3208	THR	2.2

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Mol	Chain	Res	Type	RSRZ
3	C	2132	ALA	2.2
2	B	1018	ALA	2.2
2	B	1122	PRO	2.1
1	A	233	PHE	2.1
2	B	1180	GLN	2.1
1	A	236	THR	2.1
1	A	240	THR	2.1
1	A	208	VAL	2.1
2	B	1086	VAL	2.1
1	A	461	SER	2.1
1	A	339	ASN	2.1
2	B	1164	ASN	2.1
3	C	2130	GLY	2.1
3	C	2127	LEU	2.1
3	C	2205	SER	2.1
2	B	1130	CYS	2.1
1	A	463	ASN	2.1
2	B	1126	PRO	2.1
2	B	1005	LEU	2.0
2	B	1027	HIS	2.0
2	B	1107	HIS	2.0
2	B	1001	LYS	2.0

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

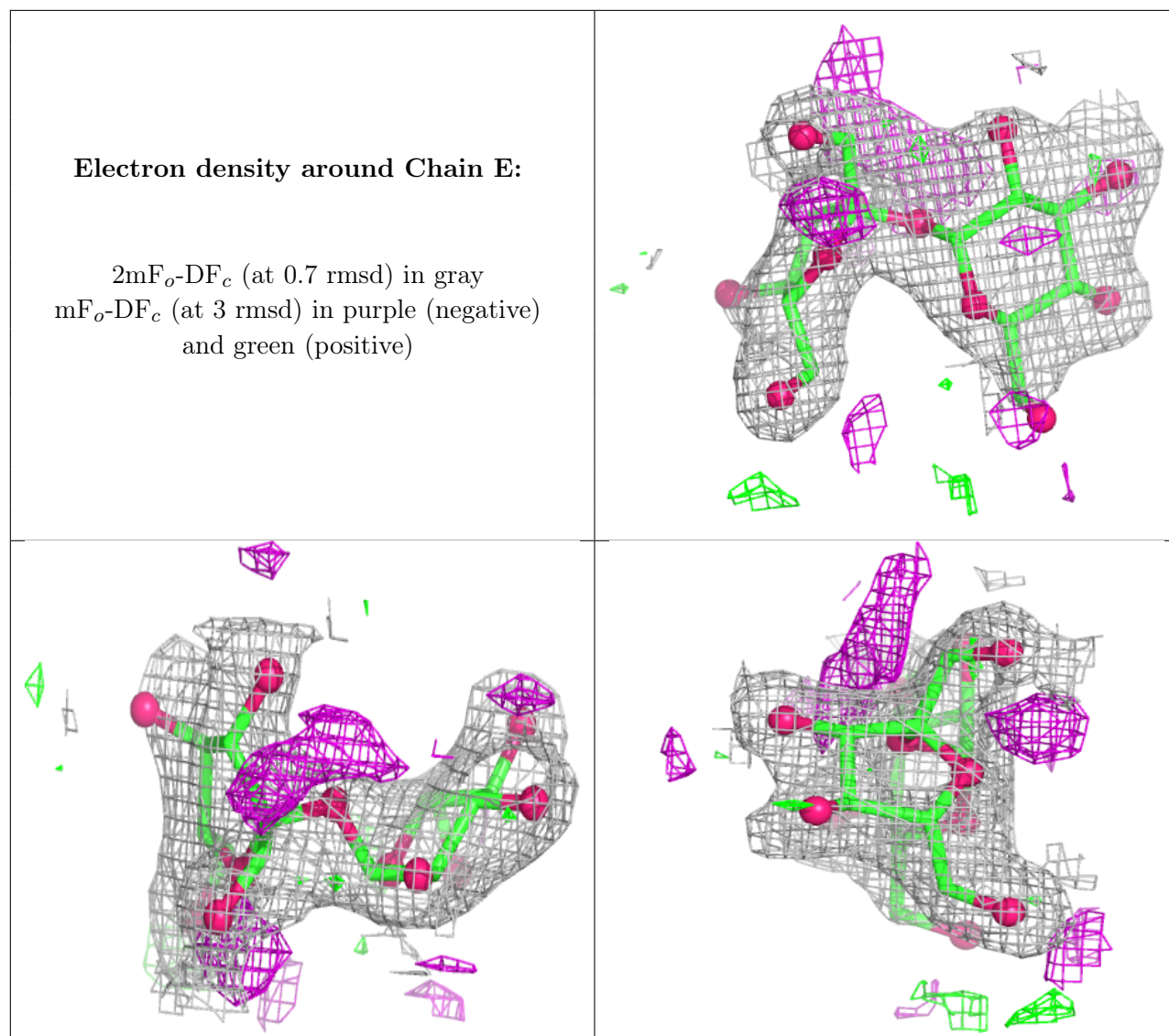
There are no non-standard protein/DNA/RNA residues in this entry.

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
5	FRU	E	2	12/12	0.72	0.21	71,75,79,81	0
5	GLC	E	1	11/12	0.79	0.18	60,62,68,69	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.



6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 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
6	NAG	A	741	14/15	0.21	0.50	570,617,651,654	0
6	NAG	A	839	14/15	0.31	0.48	503,529,544,546	0
6	NAG	A	892	14/15	0.34	0.64	359,402,439,443	0
6	NAG	A	730	14/15	0.37	0.40	585,602,630,632	0
6	NAG	A	948	14/15	0.49	0.79	386,435,470,472	0
6	NAG	A	776	14/15	0.54	0.74	404,423,430,432	0
6	NAG	A	886	14/15	0.54	0.68	260,289,303,311	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	NAG	A	789	14/15	0.56	0.56	532,552,571,579	0
6	NAG	A	734	14/15	0.59	0.64	502,528,545,549	0
6	NAG	A	762	14/15	0.65	0.79	307,331,362,363	0
6	NAG	A	795	14/15	0.68	0.74	386,398,423,424	0
8	HEZ	A	4003	8/8	0.79	0.54	371,376,382,385	0
7	EDO	D	4002	4/4	0.86	0.14	53,54,55,55	0
7	EDO	A	4001	4/4	0.86	0.46	323,324,330,332	0

6.5 Other polymers [i](#)

There are no such residues in this entry.