



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 04:24 AM BST

PDB ID : 3K1L
Title : Crystal Structure of FANCL
Authors : Cole, A.R.; Walden, H.
Deposited on : 2009-09-28
Resolution : 3.20 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
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) : 2.13.1

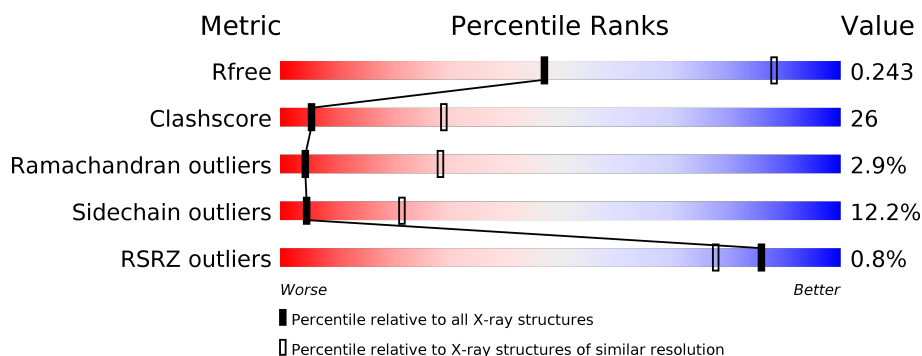
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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 on 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	381	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 52% 37% 9% .. </div> </div>
1	B	381	<div> <div style="width: 100%; height: 10px; background: linear-gradient(to right, red, orange, yellow, green, grey);"></div> <div style="display: flex; justify-content: space-between; margin-top: 5px;"> % 51% 38% 6% .. </div> </div>
2	C	2	<div> <div style="width: 100%; height: 10px; background-color: orange;"></div> <div style="text-align: center; margin-top: 5px;">100%</div> </div>
2	D	2	<div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div style="text-align: center; margin-top: 5px;">100%</div> </div>

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6079 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 FancI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	364	Total	C	N	O	S	0	0	0
			2938	1888	480	546	24			
1	A	376	Total	C	N	O	S	0	0	0
			3026	1943	494	564	25			

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



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

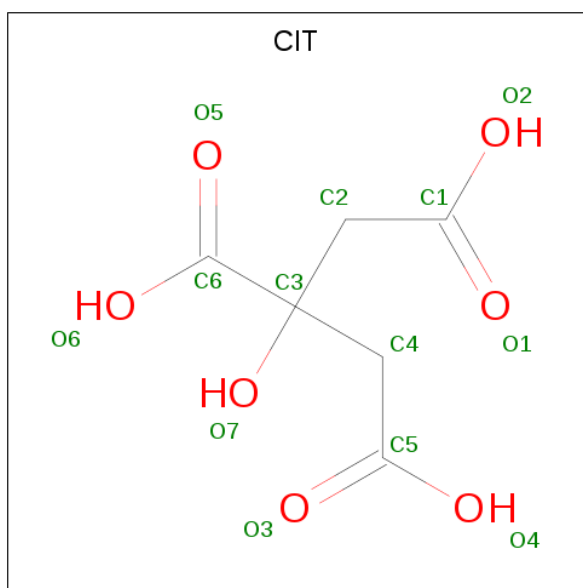
- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Zn	0	0
			2	2		
3	A	2	Total	Zn	0	0
			2	2		

- Molecule 4 is GOLD ION (three-letter code: AU) (formula: Au).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	8	Total	Au	0	0
			8	8		
4	A	5	Total	Au	0	0
			5	5		

- Molecule 5 is CITRIC ACID (three-letter code: CIT) (formula: $C_6H_8O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			13	6	7		
5	A	1	Total	C	O	0	0
			13	6	7		
5	A	1	Total	C	O	0	0
			13	6	7		

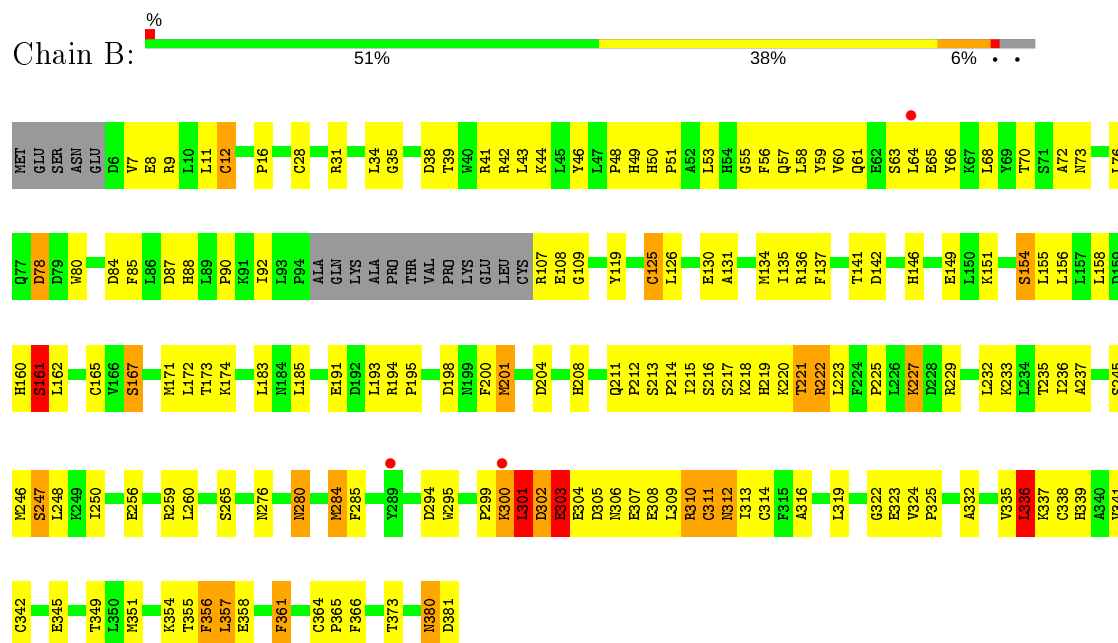
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	6	Total	O	0	0
			6	6		
6	A	7	Total	O	0	0
			7	7		

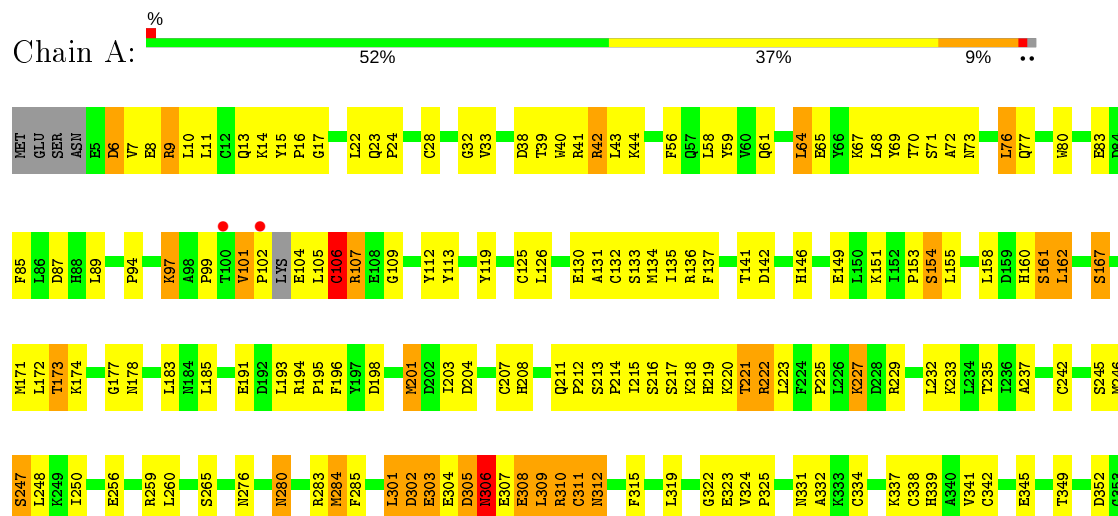
3 Residue-property plots

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: Fanc1



• Molecule 1: Fanc1





- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain C:  100%



- Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain D:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	188.68 Å 188.68 Å 259.36 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	77.93 – 3.20 77.93 – 3.20	Depositor EDS
% Data completeness (in resolution range)	92.9 (77.93-3.20) 87.4 (77.93-3.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 3.19 Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.205 , 0.247 0.202 , 0.243	Depositor DCC
R_{free} test set	1311 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	87.5	Xtriage
Anisotropy	0.081	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 77.3	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	0.93	EDS
Total number of atoms	6079	wwPDB-VP
Average B, all atoms (Å ²)	98.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.02% 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: ZN, GLC, AU, CIT

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	0.40	0/3098	0.56	0/4197
1	B	0.38	1/3008 (0.0%)	0.54	0/4073
All	All	0.39	1/6106 (0.0%)	0.55	0/8270

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	125	CYS	CB-SG	5.54	1.91	1.82

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	3026	0	2981	171	0
1	B	2938	0	2890	156	0
2	C	23	0	21	2	0
2	D	23	0	21	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	5	0	0	0	0
4	B	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	A	26	0	10	2	0
5	B	13	0	5	1	0
6	A	7	0	0	1	0
6	B	6	0	0	0	0
All	All	6079	0	5928	308	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 26.

All (308) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:101:VAL:HG22	1:A:102:PRO:HD2	1.37	1.04
1:A:213:SER:HB3	1:A:214:PRO:HD3	1.40	1.02
1:B:213:SER:HB3	1:B:214:PRO:HD3	1.41	1.00
1:A:42:ARG:HD2	1:A:43:LEU:H	1.27	0.97
1:A:23:GLN:HB3	1:A:24:PRO:HD2	1.45	0.95
1:A:70:THR:HG22	1:A:72:ALA:H	1.40	0.86
1:B:312:ASN:ND2	1:B:336:LEU:HD23	1.91	0.85
1:B:302:ASP:O	1:B:303:GLU:HB3	1.76	0.85
1:A:102:PRO:HD3	1:A:173:THR:O	1.78	0.84
1:A:216:SER:H	1:A:219:HIS:HD2	1.25	0.84
1:A:43:LEU:HD11	1:A:58:LEU:HD11	1.60	0.83
1:B:216:SER:H	1:B:219:HIS:HD2	1.27	0.82
1:A:355:THR:HG23	1:A:358:GLU:OE2	1.82	0.79
2:C:1:GLC:H61	2:C:2:GLC:H5	1.64	0.79
1:B:295:TRP:CH2	1:A:132:CYS:HB2	2.17	0.78
1:B:50:HIS:HD2	1:B:51:PRO:HA	1.48	0.78
1:B:46:TYR:HB2	1:B:57:GLN:HB2	1.65	0.78
1:A:216:SER:H	1:A:219:HIS:CD2	2.02	0.76
1:B:300:LYS:NZ	1:A:133:SER:HA	2.03	0.74
1:B:216:SER:H	1:B:219:HIS:CD2	2.03	0.74
1:A:215:ILE:HG23	1:A:222:ARG:HH12	1.53	0.74
1:B:156:LEU:HD21	1:A:310:ARG:HH11	1.53	0.74
1:B:355:THR:HG22	1:B:358:GLU:OE2	1.87	0.74
1:B:167:SER:O	1:B:171:MET:HG3	1.88	0.74
1:B:311:CYS:O	1:B:312:ASN:HB2	1.87	0.73
1:A:311:CYS:O	1:A:312:ASN:HB2	1.88	0.72
1:B:12:CYS:SG	1:A:349:THR:HG21	2.29	0.72
1:A:70:THR:N	1:A:73:ASN:HD22	1.88	0.72
1:A:204:ASP:OD1	1:A:222:ARG:HD3	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:324:VAL:HG13	1:A:325:PRO:HD2	1.70	0.72
1:B:312:ASN:CG	1:B:336:LEU:HD23	2.10	0.72
1:B:204:ASP:OD1	1:B:222:ARG:HD3	1.90	0.71
1:A:167:SER:O	1:A:171:MET:HG3	1.90	0.71
1:B:227:LYS:NZ	1:B:259:ARG:HH22	1.89	0.71
1:A:65:GLU:OE2	5:A:2504:CIT:O1	2.09	0.70
1:A:227:LYS:NZ	1:A:259:ARG:HH22	1.88	0.70
1:A:304:GLU:HG2	1:A:306:ASN:ND2	2.07	0.69
1:B:70:THR:H	1:B:73:ASN:HB2	1.58	0.69
1:B:324:VAL:HG13	1:B:325:PRO:HD2	1.74	0.69
1:A:198:ASP:O	1:A:201:MET:HB2	1.93	0.69
1:A:9:ARG:HG3	1:A:10:LEU:N	2.07	0.68
1:A:42:ARG:HD2	1:A:43:LEU:N	2.06	0.68
1:A:357:LEU:HD23	1:A:357:LEU:H	1.59	0.68
1:A:97:LYS:HE3	1:A:97:LYS:H	1.59	0.67
1:B:215:ILE:HG23	1:B:222:ARG:HH12	1.59	0.67
1:A:310:ARG:O	1:A:311:CYS:C	2.33	0.66
1:B:355:THR:HG23	1:B:357:LEU:HG	1.77	0.66
1:B:50:HIS:CD2	1:B:51:PRO:HA	2.31	0.66
1:A:227:LYS:HZ3	1:A:259:ARG:HH22	1.43	0.65
1:B:53:LEU:O	1:B:56:PHE:HB3	1.96	0.65
1:A:106:CYS:SG	1:A:107:ARG:N	2.71	0.64
1:B:198:ASP:O	1:B:201:MET:HB2	1.98	0.64
1:B:310:ARG:O	1:B:311:CYS:C	2.36	0.63
1:A:23:GLN:HB3	1:A:24:PRO:CD	2.26	0.63
1:B:366:PHE:CD1	1:A:38:ASP:HA	2.34	0.63
1:B:308:GLU:OE1	1:B:308:GLU:HA	1.98	0.62
1:B:194:ARG:HG3	1:B:195:PRO:HD3	1.82	0.62
1:B:43:LEU:HD11	1:B:58:LEU:HD11	1.82	0.62
1:A:106:CYS:O	1:A:107:ARG:HD2	1.99	0.62
1:B:310:ARG:NH1	1:A:173:THR:HG21	2.15	0.61
1:B:357:LEU:H	1:B:357:LEU:HD23	1.65	0.61
1:A:76:LEU:HD22	1:A:85:PHE:HE1	1.65	0.61
1:A:105:LEU:N	1:A:105:LEU:HD23	2.16	0.60
1:B:316:ALA:HB2	1:A:64:LEU:HD11	1.84	0.60
1:A:191:GLU:O	1:A:194:ARG:HG2	2.01	0.60
1:A:194:ARG:HG3	1:A:195:PRO:HD3	1.83	0.60
1:B:221:THR:HG23	1:B:222:ARG:N	2.17	0.59
1:A:221:THR:HG23	1:A:222:ARG:N	2.18	0.59
1:B:191:GLU:O	1:B:194:ARG:HG2	2.02	0.59
1:B:300:LYS:HZ1	1:A:133:SER:HA	1.68	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:CYS:SG	1:A:107:ARG:HG3	2.43	0.59
1:A:361:PHE:HA	1:A:373:THR:HG23	1.85	0.59
1:B:355:THR:CG2	1:B:358:GLU:HG3	2.33	0.59
1:B:130:GLU:HG3	1:B:131:ALA:N	2.18	0.58
1:A:232:LEU:HD11	1:A:248:LEU:HD22	1.85	0.58
1:A:43:LEU:HD12	1:A:44:LYS:N	2.19	0.58
1:B:42:ARG:NH1	1:A:315:PHE:HE1	2.01	0.58
1:A:216:SER:N	1:A:219:HIS:HD2	1.98	0.58
1:A:309:LEU:HD23	1:A:309:LEU:N	2.19	0.58
1:B:154:SER:O	1:B:155:LEU:HB2	2.03	0.58
1:B:216:SER:N	1:B:219:HIS:HD2	1.99	0.58
1:B:361:PHE:HA	1:B:373:THR:HG23	1.86	0.57
1:B:301:LEU:N	1:B:301:LEU:HD23	2.19	0.57
1:B:34:LEU:HD21	1:B:90:PRO:HD3	1.86	0.57
1:B:276:ASN:O	1:B:280:ASN:HB2	2.05	0.57
1:A:154:SER:O	1:A:155:LEU:HB2	2.04	0.57
1:A:160:HIS:ND1	1:A:161:SER:O	2.37	0.57
1:A:213:SER:HB3	1:A:214:PRO:CD	2.26	0.57
1:A:101:VAL:CG2	1:A:102:PRO:HD2	2.24	0.57
1:B:151:LYS:HG3	1:B:158:LEU:HD21	1.87	0.57
1:A:130:GLU:HG3	1:A:131:ALA:N	2.19	0.57
1:A:76:LEU:HD22	1:A:85:PHE:CE1	2.39	0.56
1:B:213:SER:HB3	1:B:214:PRO:CD	2.28	0.56
1:A:134:MET:HG3	1:A:158:LEU:HD12	1.87	0.56
1:B:134:MET:HG3	1:B:158:LEU:HD12	1.87	0.56
1:B:160:HIS:ND1	1:B:161:SER:O	2.39	0.56
1:A:308:GLU:HA	1:A:308:GLU:OE1	2.05	0.56
1:B:227:LYS:HZ3	1:B:259:ARG:HH22	1.52	0.56
1:B:172:LEU:HD23	1:B:185:LEU:HD23	1.88	0.56
1:B:305:ASP:O	1:B:306:ASN:C	2.42	0.56
1:A:208:HIS:O	1:A:225:PRO:HD3	2.06	0.56
1:A:216:SER:N	1:A:219:HIS:CD2	2.74	0.55
1:A:276:ASN:O	1:A:280:ASN:HB2	2.06	0.55
1:B:336:LEU:HD22	1:B:366:PHE:CE1	2.41	0.55
1:B:313:ILE:HD13	1:A:33:VAL:HG11	1.88	0.55
1:B:88:HIS:HB3	1:B:92:ILE:HD11	1.88	0.55
1:B:221:THR:CG2	1:B:222:ARG:N	2.69	0.55
1:B:300:LYS:HZ3	1:A:133:SER:HA	1.68	0.55
1:B:303:GLU:OE2	1:B:305:ASP:HB2	2.06	0.55
1:A:151:LYS:HG3	1:A:158:LEU:HD21	1.88	0.55
1:B:246:MET:HG2	1:B:247:SER:N	2.21	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:232:LEU:HD11	1:B:248:LEU:HD22	1.88	0.55
1:B:365:PRO:HB3	1:A:40:TRP:CZ2	2.41	0.55
1:B:59:TYR:HB3	1:B:66:TYR:HD2	1.72	0.55
1:A:215:ILE:CG2	1:A:222:ARG:HH12	2.18	0.55
1:A:246:MET:HG2	1:A:247:SER:N	2.22	0.54
1:A:198:ASP:HA	1:A:201:MET:HG3	1.89	0.54
1:A:172:LEU:HD23	1:A:185:LEU:HD23	1.88	0.54
1:A:324:VAL:HG13	1:A:325:PRO:CD	2.36	0.54
1:B:345:GLU:O	1:B:349:THR:HG22	2.07	0.54
1:B:59:TYR:HB3	1:B:66:TYR:CD2	2.42	0.54
1:A:42:ARG:HB3	1:A:61:GLN:HB3	1.88	0.54
1:B:141:THR:O	1:B:142:ASP:HB2	2.08	0.54
1:A:221:THR:CG2	1:A:222:ARG:N	2.71	0.53
1:A:305:ASP:O	1:A:306:ASN:C	2.47	0.53
1:B:61:GLN:HG3	1:B:61:GLN:O	2.09	0.53
1:B:198:ASP:HA	1:B:201:MET:HG3	1.90	0.53
1:A:308:GLU:C	1:A:309:LEU:HD23	2.29	0.53
1:A:69:TYR:HA	1:A:73:ASN:ND2	2.23	0.53
1:B:355:THR:HG23	1:B:358:GLU:HG3	1.91	0.53
1:B:237:ALA:HB3	1:B:245:SER:HB2	1.91	0.53
1:A:213:SER:CB	1:A:214:PRO:HD3	2.25	0.52
1:A:345:GLU:O	1:A:349:THR:HG22	2.09	0.52
1:A:77:GLN:HB2	1:A:80:TRP:CD1	2.45	0.52
1:A:22:LEU:C	1:A:23:GLN:HG2	2.30	0.52
1:A:356:PHE:CD2	1:A:356:PHE:C	2.83	0.52
5:B:2502:CIT:O4	5:B:2502:CIT:H21	2.09	0.52
1:A:119:TYR:HA	1:A:126:LEU:HD23	1.91	0.52
1:A:69:TYR:OH	1:A:94:PRO:HG2	2.10	0.52
1:B:216:SER:N	1:B:219:HIS:CD2	2.75	0.52
1:B:300:LYS:O	1:B:302:ASP:N	2.43	0.52
1:A:213:SER:HB2	1:A:322:GLY:HA3	1.92	0.52
1:A:76:LEU:HB3	1:A:80:TRP:CE3	2.45	0.52
1:A:7:VAL:HG23	1:A:8:GLU:N	2.24	0.51
1:B:208:HIS:O	1:B:225:PRO:HD3	2.10	0.51
1:A:361:PHE:CE1	1:A:370:LYS:HE2	2.46	0.51
1:A:65:GLU:CD	5:A:2504:CIT:O1	2.49	0.51
1:A:352:ASP:O	1:A:354:LYS:HG2	2.11	0.51
1:B:213:SER:CB	1:B:214:PRO:HD3	2.25	0.51
1:B:28:CYS:SG	1:B:50:HIS:ND1	2.84	0.51
1:B:130:GLU:O	1:B:131:ALA:HB3	2.10	0.51
1:B:213:SER:HB2	1:B:322:GLY:HA3	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:301:LEU:HD21	1:A:112:TYR:HE1	1.76	0.51
1:B:43:LEU:CD1	1:B:60:VAL:HG22	2.41	0.51
1:A:70:THR:CG2	1:A:71:SER:N	2.74	0.50
1:B:215:ILE:CG2	1:B:222:ARG:HH12	2.24	0.50
1:B:88:HIS:O	1:B:92:ILE:HG13	2.11	0.50
1:A:105:LEU:HD12	1:A:178:ASN:HD22	1.77	0.50
1:A:301:LEU:O	1:A:302:ASP:C	2.50	0.50
1:A:70:THR:HG22	1:A:71:SER:N	2.25	0.50
1:B:319:LEU:HD12	1:B:323:GLU:HB2	1.94	0.50
1:A:6:ASP:HA	1:A:9:ARG:HG2	1.93	0.50
1:A:141:THR:O	1:A:142:ASP:HB2	2.12	0.50
1:A:283:ARG:NH1	6:A:393:HOH:O	2.44	0.50
1:A:6:ASP:O	1:A:9:ARG:CG	2.59	0.50
1:B:304:GLU:O	1:B:305:ASP:C	2.50	0.50
1:B:309:LEU:O	1:B:325:PRO:HG2	2.10	0.50
1:B:357:LEU:HD23	1:B:357:LEU:N	2.26	0.49
1:A:33:VAL:HA	1:A:41:ARG:O	2.11	0.49
1:B:364:CYS:C	1:B:366:PHE:H	2.15	0.49
1:B:44:LYS:HE2	1:B:46:TYR:HE1	1.77	0.49
1:A:130:GLU:O	1:A:131:ALA:HB3	2.13	0.49
1:B:119:TYR:HA	1:B:126:LEU:HD23	1.93	0.49
1:A:237:ALA:HB3	1:A:245:SER:HB2	1.95	0.49
1:B:84:ASP:O	1:B:88:HIS:HB2	2.12	0.49
1:A:221:THR:O	1:A:222:ARG:HG3	2.13	0.48
1:A:305:ASP:O	1:A:307:GLU:N	2.47	0.48
1:B:48:PRO:HG2	1:B:55:GLY:O	2.13	0.48
1:A:223:LEU:HD13	1:A:233:LYS:HB2	1.96	0.48
1:A:302:ASP:C	1:A:304:GLU:H	2.17	0.48
1:B:221:THR:O	1:B:222:ARG:HG3	2.14	0.48
1:B:76:LEU:HD21	1:B:85:PHE:CE2	2.49	0.48
1:B:301:LEU:O	1:B:302:ASP:C	2.52	0.48
1:A:11:LEU:HD23	1:A:15:TYR:HD1	1.77	0.47
1:A:361:PHE:CD2	1:A:361:PHE:C	2.87	0.47
1:B:361:PHE:CD2	1:B:361:PHE:C	2.86	0.47
1:A:312:ASN:HB2	1:A:337:LYS:O	2.14	0.47
1:A:6:ASP:O	1:A:9:ARG:HG3	2.15	0.47
1:B:227:LYS:HZ1	1:B:259:ARG:HH22	1.63	0.47
1:B:70:THR:HG23	1:B:72:ALA:H	1.79	0.47
1:B:355:THR:HG22	1:B:358:GLU:HG3	1.96	0.47
1:A:229:ARG:HH22	1:A:332:ALA:HB1	1.80	0.47
1:A:357:LEU:HD23	1:A:357:LEU:N	2.27	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:43:LEU:HD11	1:A:58:LEU:CD1	2.37	0.47
1:B:311:CYS:H	1:B:325:PRO:HG3	1.79	0.47
1:B:312:ASN:HB2	1:B:337:LYS:O	2.14	0.47
1:A:229:ARG:HH22	1:A:332:ALA:CB	2.28	0.47
1:A:256:GLU:OE2	1:A:256:GLU:HA	2.15	0.46
1:A:319:LEU:HD12	1:A:323:GLU:HB2	1.96	0.46
1:A:211:GLN:HA	1:A:212:PRO:C	2.34	0.46
1:A:305:ASP:C	1:A:306:ASN:HD22	2.19	0.46
1:A:361:PHE:CA	1:A:373:THR:HG23	2.45	0.46
1:A:97:LYS:HD3	1:A:97:LYS:N	2.31	0.46
1:B:361:PHE:CA	1:B:373:THR:HG23	2.45	0.46
1:B:223:LEU:HD13	1:B:233:LYS:HB2	1.98	0.46
1:A:59:TYR:CZ	1:A:68:LEU:HD13	2.51	0.46
1:B:300:LYS:O	1:B:301:LEU:HG	2.15	0.46
1:B:335:VAL:O	1:B:335:VAL:HG23	2.16	0.46
1:A:380:ASN:ND2	1:A:381:ASP:N	2.63	0.46
1:A:304:GLU:HG2	1:A:306:ASN:HD22	1.80	0.46
1:B:211:GLN:HA	1:B:212:PRO:C	2.36	0.46
1:B:324:VAL:HG13	1:B:325:PRO:CD	2.43	0.46
1:B:229:ARG:HH22	1:B:332:ALA:CB	2.30	0.45
1:B:351:MET:HE1	1:B:356:PHE:HB2	1.97	0.45
1:A:151:LYS:HG3	1:A:158:LEU:HD11	1.99	0.45
1:B:194:ARG:N	1:B:195:PRO:CD	2.80	0.45
1:B:335:VAL:O	1:B:336:LEU:CB	2.64	0.45
1:B:42:ARG:HB3	1:B:61:GLN:HG2	1.98	0.45
1:B:299:PRO:HB3	1:A:112:TYR:CE1	2.52	0.45
1:B:310:ARG:CZ	1:A:173:THR:HG21	2.47	0.45
1:B:335:VAL:O	1:B:336:LEU:CG	2.64	0.45
1:B:44:LYS:HE2	1:B:46:TYR:CE1	2.52	0.45
1:A:67:LYS:HD3	1:A:69:TYR:CE1	2.51	0.45
1:B:229:ARG:HH22	1:B:332:ALA:HB1	1.82	0.45
1:A:102:PRO:HB2	1:A:177:GLY:HA2	1.99	0.45
1:B:311:CYS:H	1:B:325:PRO:CG	2.30	0.45
1:A:311:CYS:H	1:A:325:PRO:HG3	1.83	0.44
1:B:256:GLU:HA	1:B:256:GLU:OE2	2.16	0.44
2:C:1:GLC:C6	2:C:2:GLC:H5	2.43	0.44
1:B:156:LEU:CD2	1:A:310:ARG:HH11	2.25	0.44
1:B:380:ASN:ND2	1:B:381:ASP:N	2.64	0.44
1:A:7:VAL:HG23	1:A:8:GLU:H	1.82	0.44
1:A:136:ARG:HG3	1:A:149:GLU:HG2	1.99	0.44
1:B:49:HIS:O	1:B:50:HIS:C	2.55	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:TRP:CD1	1:B:84:ASP:HB3	2.53	0.44
1:B:38:ASP:O	1:B:39:THR:HB	2.18	0.43
1:B:8:GLU:HA	1:B:11:LEU:HB2	1.99	0.43
1:B:41:ARG:NH1	1:B:90:PRO:HA	2.33	0.43
1:A:161:SER:O	1:A:162:LEU:HB2	2.18	0.43
1:B:151:LYS:HG3	1:B:158:LEU:HD11	2.00	0.43
1:B:165:CYS:SG	1:B:220:LYS:HD3	2.59	0.43
1:B:311:CYS:HB3	1:B:312:ASN:H	1.52	0.43
1:A:339:HIS:HB2	1:A:342:CYS:SG	2.58	0.43
1:A:97:LYS:N	1:A:97:LYS:CD	2.82	0.43
1:B:41:ARG:HH12	1:B:90:PRO:HA	1.84	0.43
1:B:200:PHE:CZ	1:B:236:ILE:HD12	2.54	0.43
1:B:212:PRO:HG2	1:B:219:HIS:CG	2.54	0.43
1:A:146:HIS:CD2	1:A:161:SER:HB2	2.54	0.43
1:A:216:SER:C	1:A:218:LYS:H	2.22	0.43
1:B:294:ASP:OD1	1:B:295:TRP:N	2.51	0.43
1:A:6:ASP:O	1:A:9:ARG:HG2	2.18	0.43
1:B:310:ARG:HH12	1:A:173:THR:HG21	1.83	0.43
1:A:250:ILE:HD13	1:A:260:LEU:HB3	2.01	0.42
1:A:70:THR:HB	1:A:73:ASN:ND2	2.34	0.42
1:A:14:LYS:HZ2	1:A:83:GLU:CG	2.32	0.42
1:B:364:CYS:C	1:B:366:PHE:N	2.73	0.42
1:B:61:GLN:CG	1:B:61:GLN:O	2.68	0.42
1:A:137:PHE:CD2	1:A:183:LEU:HD13	2.55	0.42
1:A:194:ARG:N	1:A:195:PRO:CD	2.82	0.42
1:A:301:LEU:HG	1:A:302:ASP:N	2.33	0.42
1:B:310:ARG:HH12	1:A:173:THR:CG2	2.32	0.42
1:B:70:THR:N	1:B:73:ASN:HB2	2.29	0.42
1:A:7:VAL:O	1:A:10:LEU:HB3	2.20	0.42
1:A:212:PRO:HG2	1:A:219:HIS:CG	2.54	0.42
1:A:379:LEU:O	1:A:381:ASP:N	2.47	0.42
1:A:97:LYS:CE	1:A:97:LYS:H	2.30	0.42
1:B:355:THR:CG2	1:B:357:LEU:HG	2.48	0.42
1:A:113:TYR:CD2	1:A:113:TYR:C	2.93	0.42
1:B:303:GLU:CD	1:B:305:ASP:H	2.23	0.42
1:B:349:THR:CG2	1:A:16:PRO:O	2.68	0.42
1:B:28:CYS:CB	1:B:50:HIS:ND1	2.83	0.42
1:B:78:ASP:N	1:B:78:ASP:OD1	2.53	0.42
1:A:220:LYS:HE3	1:A:237:ALA:O	2.20	0.42
1:A:354:LYS:H	1:A:354:LYS:HG2	1.56	0.42
1:B:306:ASN:O	1:B:308:GLU:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:361:PHE:HD2	1:B:361:PHE:C	2.23	0.42
1:B:284:MET:HG2	1:B:285:PHE:CE2	2.55	0.41
1:A:196:PHE:HA	1:A:242:CYS:SG	2.60	0.41
1:A:284:MET:HG2	1:A:285:PHE:CE2	2.55	0.41
1:A:203:ILE:O	1:A:207:CYS:HB2	2.20	0.41
1:A:106:CYS:HB3	1:A:107:ARG:NH1	2.35	0.41
1:A:134:MET:HG2	1:A:135:ILE:N	2.36	0.41
1:A:364:CYS:HA	1:A:365:PRO:HD3	1.93	0.41
1:B:300:LYS:HD3	1:A:153:PRO:HD3	2.02	0.41
1:B:216:SER:C	1:B:218:LYS:H	2.24	0.41
1:A:311:CYS:HB3	1:A:312:ASN:H	1.55	0.41
1:A:311:CYS:H	1:A:325:PRO:CG	2.34	0.41
1:A:89:LEU:HD12	1:A:89:LEU:HA	1.59	0.41
1:B:134:MET:HG2	1:B:135:ILE:N	2.36	0.41
1:B:301:LEU:O	1:B:304:GLU:CD	2.58	0.41
1:A:324:VAL:HA	1:A:339:HIS:HE2	1.85	0.41
1:B:136:ARG:HG3	1:B:149:GLU:HG2	2.03	0.41
1:B:146:HIS:CD2	1:B:161:SER:HB2	2.56	0.41
1:B:59:TYR:CZ	1:B:68:LEU:HD13	2.56	0.41
1:A:9:ARG:NH1	1:A:13:GLN:OE1	2.54	0.41
1:A:56:PHE:C	1:A:56:PHE:CD1	2.94	0.41
1:B:134:MET:HB3	1:B:134:MET:HE3	1.93	0.41
1:B:137:PHE:CD2	1:B:183:LEU:HD13	2.55	0.41
1:B:250:ILE:HD13	1:B:260:LEU:HB3	2.03	0.41
1:B:35:GLY:HA3	1:A:366:PHE:CD1	2.56	0.41
1:B:16:PRO:HD3	1:A:365:PRO:HB3	2.03	0.41
1:A:17:GLY:O	1:A:32:GLY:HA3	2.21	0.40
1:A:303:GLU:H	1:A:303:GLU:HG2	1.46	0.40
1:A:70:THR:H	1:A:73:ASN:HD22	1.63	0.40
1:A:331:ASN:HB3	1:A:334:CYS:HB2	2.03	0.40
1:A:361:PHE:HD2	1:A:361:PHE:C	2.24	0.40
1:B:335:VAL:O	1:B:336:LEU:HG	2.22	0.40
1:B:339:HIS:HB2	1:B:342:CYS:SG	2.61	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	372/381 (98%)	329 (88%)	34 (9%)	9 (2%)	6	34
1	B	360/381 (94%)	315 (88%)	33 (9%)	12 (3%)	4	25
All	All	732/762 (96%)	644 (88%)	67 (9%)	21 (3%)	4	28

All (21) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	300	LYS
1	B	312	ASN
1	B	336	LEU
1	B	354	LYS
1	A	106	CYS
1	A	312	ASN
1	B	63	SER
1	B	301	LEU
1	B	303	GLU
1	A	306	ASN
1	B	307	GLU
1	A	302	ASP
1	A	311	CYS
1	B	109	GLY
1	B	161	SER
1	B	311	CYS
1	A	109	GLY
1	A	305	ASP
1	B	162	LEU
1	A	162	LEU
1	A	99	PRO

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	340/345 (99%)	297 (87%)	43 (13%)	4	21
1	B	330/345 (96%)	291 (88%)	39 (12%)	5	23
All	All	670/690 (97%)	588 (88%)	82 (12%)	5	22

All (82) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	7	VAL
1	B	9	ARG
1	B	12	CYS
1	B	31	ARG
1	B	64	LEU
1	B	65	GLU
1	B	78	ASP
1	B	87	ASP
1	B	107	ARG
1	B	108	GLU
1	B	125	CYS
1	B	154	SER
1	B	161	SER
1	B	167	SER
1	B	173	THR
1	B	174	LYS
1	B	193	LEU
1	B	201	MET
1	B	217	SER
1	B	221	THR
1	B	222	ARG
1	B	227	LYS
1	B	235	THR
1	B	247	SER
1	B	265	SER
1	B	280	ASN
1	B	284	MET

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Mol	Chain	Res	Type
1	B	301	LEU
1	B	302	ASP
1	B	303	GLU
1	B	310	ARG
1	B	314	CYS
1	B	336	LEU
1	B	338	CYS
1	B	341	VAL
1	B	356	PHE
1	B	357	LEU
1	B	361	PHE
1	B	380	ASN
1	A	6	ASP
1	A	9	ARG
1	A	28	CYS
1	A	39	THR
1	A	42	ARG
1	A	64	LEU
1	A	76	LEU
1	A	87	ASP
1	A	97	LYS
1	A	101	VAL
1	A	104	GLU
1	A	106	CYS
1	A	107	ARG
1	A	125	CYS
1	A	154	SER
1	A	161	SER
1	A	167	SER
1	A	173	THR
1	A	174	LYS
1	A	193	LEU
1	A	201	MET
1	A	217	SER
1	A	221	THR
1	A	222	ARG
1	A	227	LYS
1	A	235	THR
1	A	247	SER
1	A	265	SER
1	A	280	ASN
1	A	284	MET

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Mol	Chain	Res	Type
1	A	301	LEU
1	A	303	GLU
1	A	306	ASN
1	A	308	GLU
1	A	309	LEU
1	A	310	ARG
1	A	338	CYS
1	A	341	VAL
1	A	355	THR
1	A	356	PHE
1	A	357	LEU
1	A	361	PHE
1	A	380	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	73	ASN
1	B	219	HIS
1	B	312	ASN
1	B	380	ASN
1	A	73	ASN
1	A	219	HIS
1	A	306	ASN
1	A	380	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 ⓘ

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	GLC	C	1	2	12,12,12	0.60	0	17,17,17	1.22	1 (5%)
2	GLC	C	2	2	11,11,12	0.67	0	15,15,17	1.34	2 (13%)
2	GLC	D	1	2	12,12,12	0.63	0	17,17,17	1.07	1 (5%)
2	GLC	D	2	2	11,11,12	0.55	0	15,15,17	1.17	1 (6%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	C	1	2	-	0/2/22/22	0/1/1/1
2	GLC	C	2	2	-	2/2/19/22	0/1/1/1
2	GLC	D	1	2	-	0/2/22/22	0/1/1/1
2	GLC	D	2	2	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	GLC	O5-C1-C2	3.33	116.23	110.28
2	C	2	GLC	O5-C5-C6	3.32	112.41	107.20
2	D	2	GLC	O5-C5-C6	2.49	111.11	107.20
2	C	2	GLC	C1-C2-C3	-2.08	107.10	109.67
2	D	1	GLC	O5-C1-C2	2.08	113.99	110.28

There are no chirality outliers.

All (2) torsion outliers are listed below:

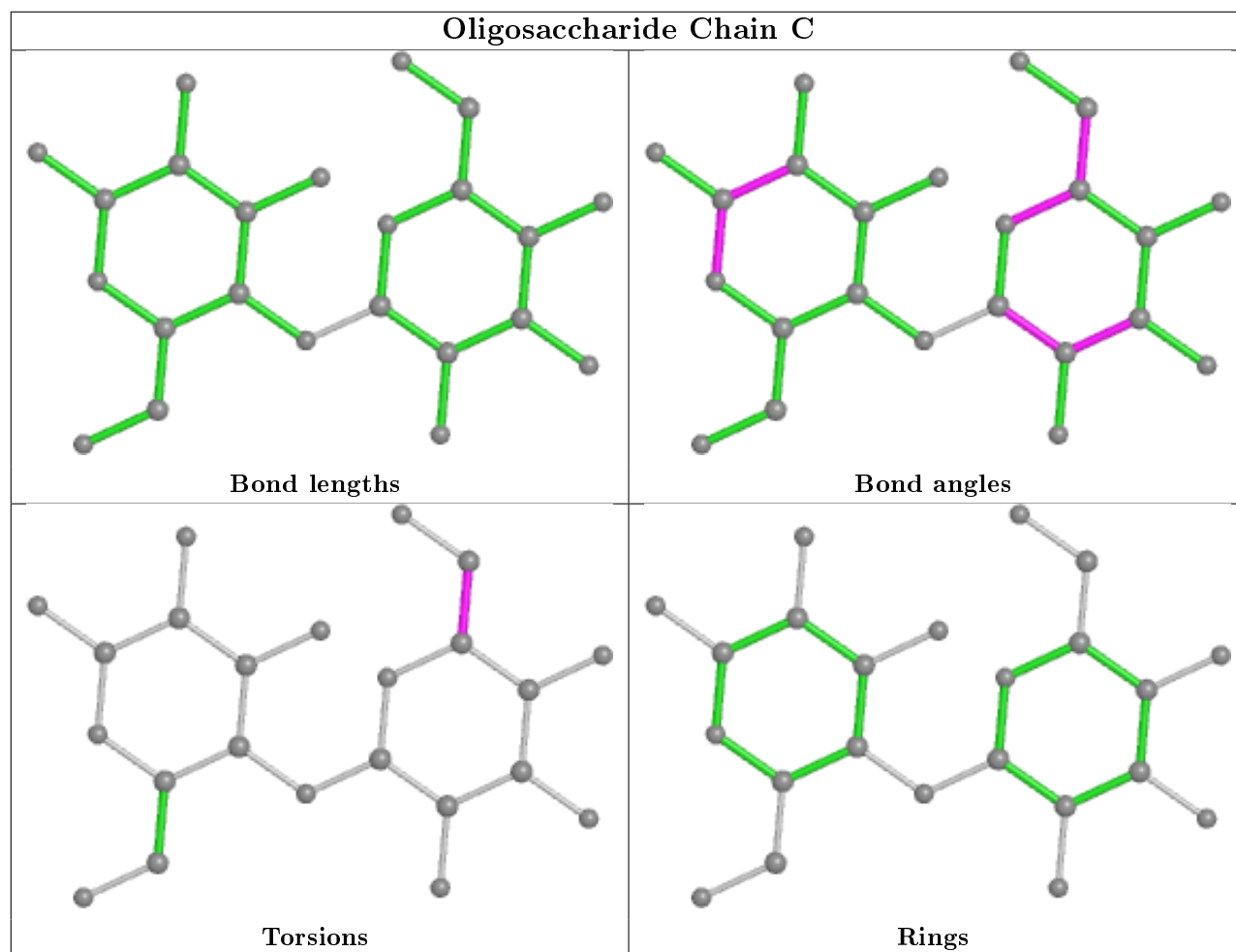
Mol	Chain	Res	Type	Atoms
2	C	2	GLC	O5-C5-C6-O6
2	C	2	GLC	C4-C5-C6-O6

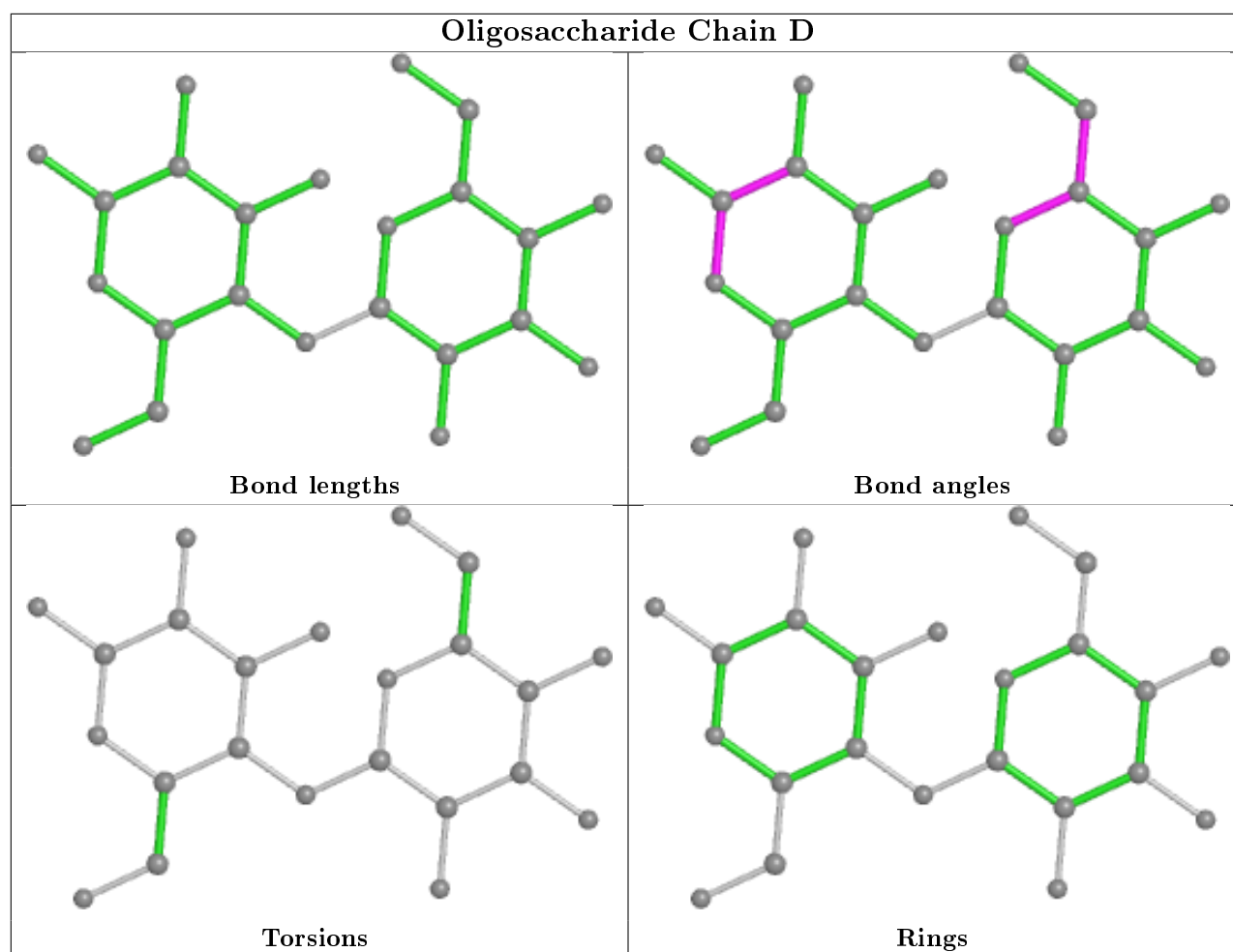
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	GLC	2	0
2	C	2	GLC	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](#)

Of 20 ligands modelled in this entry, 17 are monoatomic - leaving 3 for Mogul analysis.

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	CIT	A	2503	-	3,12,12	1.25	0	3,17,17	1.79	1 (33%)
5	CIT	A	2504	-	3,12,12	1.61	0	3,17,17	3.05	2 (66%)
5	CIT	B	2502	-	3,12,12	1.23	0	3,17,17	1.74	1 (33%)

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	CIT	A	2503	-	-	3/6/16/16	-
5	CIT	A	2504	-	-	6/6/16/16	-
5	CIT	B	2502	-	-	3/6/16/16	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	2504	CIT	C3-C2-C1	-3.85	108.83	114.98
5	A	2504	CIT	C3-C4-C5	-3.55	109.30	114.98
5	A	2503	CIT	C3-C4-C5	-3.08	110.06	114.98
5	B	2502	CIT	C3-C2-C1	-2.30	111.30	114.98

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	2504	CIT	C2-C3-C4-C5
5	A	2504	CIT	O7-C3-C4-C5
5	A	2504	CIT	C6-C3-C4-C5
5	B	2502	CIT	C1-C2-C3-C6
5	A	2504	CIT	C1-C2-C3-C4
5	B	2502	CIT	C1-C2-C3-O7
5	B	2502	CIT	C1-C2-C3-C4
5	A	2503	CIT	C2-C3-C4-C5
5	A	2504	CIT	C1-C2-C3-O7
5	A	2504	CIT	C1-C2-C3-C6
5	A	2503	CIT	C1-C2-C3-C6
5	A	2503	CIT	C1-C2-C3-O7

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	2504	CIT	2	0
5	B	2502	CIT	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	376/381 (98%)	0.01	3 (0%) 86 78	58, 82, 160, 240	0
1	B	364/381 (95%)	0.15	3 (0%) 86 78	60, 91, 184, 279	0
All	All	740/762 (97%)	0.08	6 (0%) 86 78	58, 86, 175, 279	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	289	TYR	3.0
1	A	381	ASP	2.9
1	A	102	PRO	2.8
1	A	100	THR	2.7
1	B	300	LYS	2.3
1	B	64	LEU	2.2

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

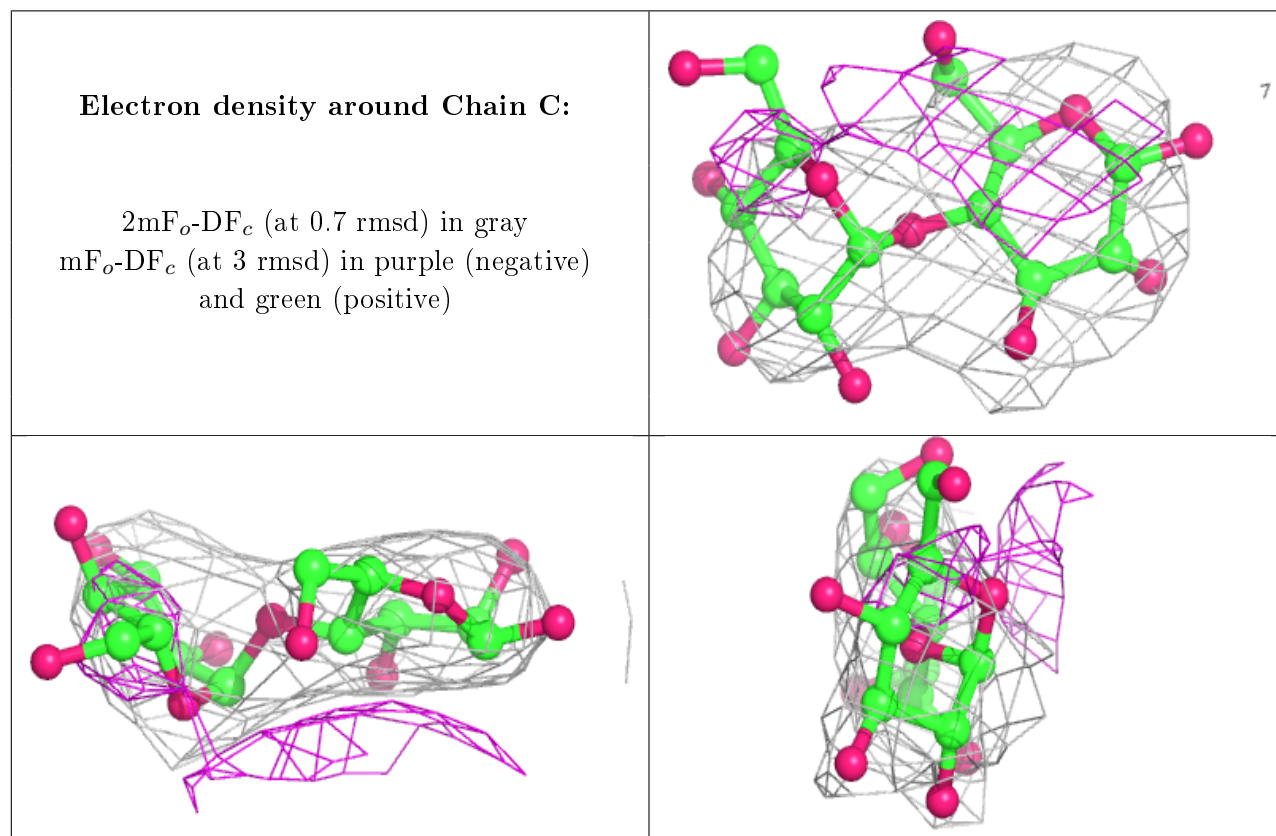
There are no non-standard protein/DNA/RNA residues in this entry.

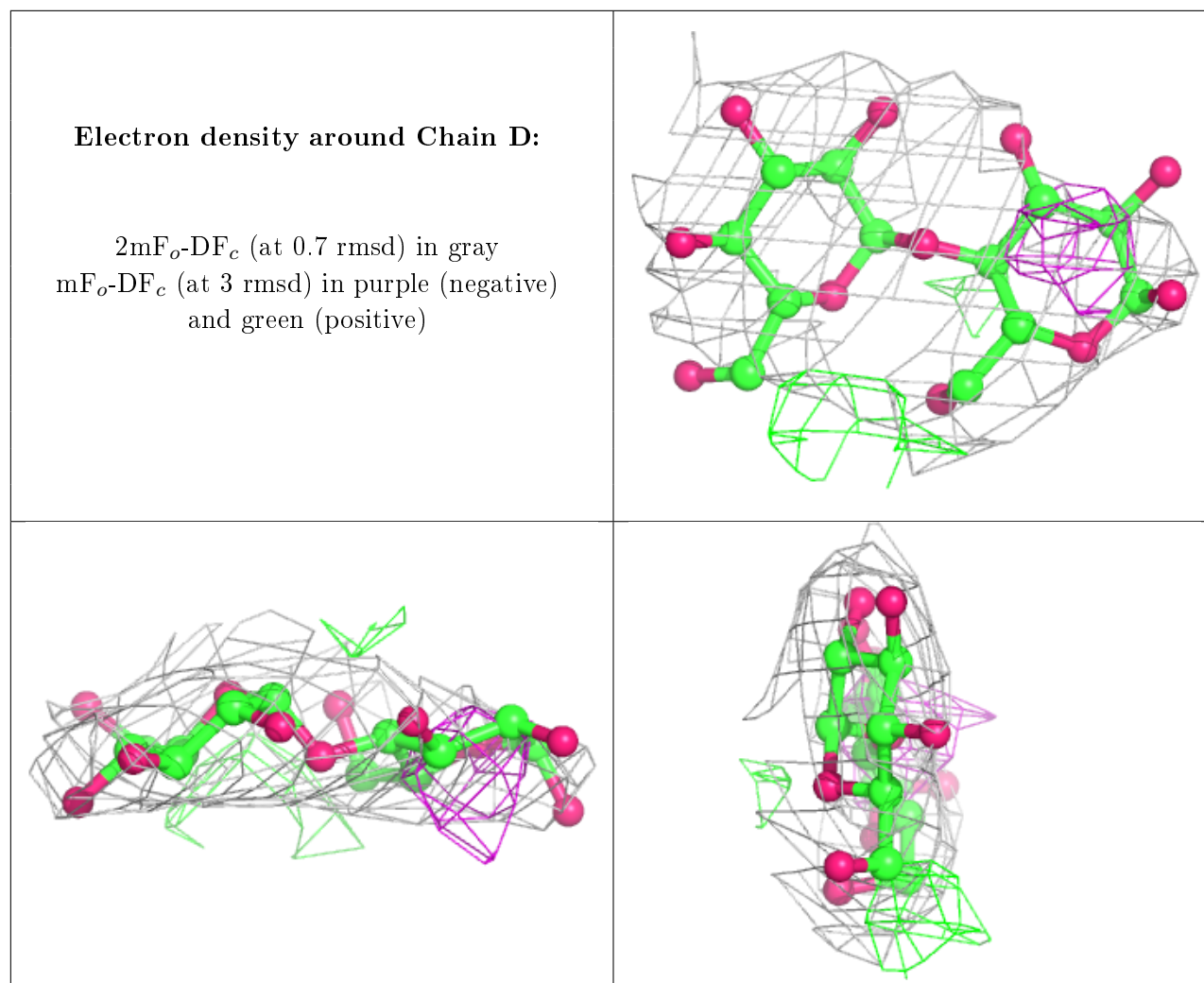
6.3 Carbohydrates ⓘ

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	GLC	D	1	12/12	0.71	0.31	115,115,115,115	12
2	GLC	C	2	11/12	0.80	0.49	115,115,115,115	11
2	GLC	D	2	11/12	0.84	0.20	115,115,115,115	11
2	GLC	C	1	12/12	0.86	0.45	115,115,115,115	12

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(\AA^2)	Q<0.9
5	CIT	A	2503	13/13	0.77	0.25	115,115,115,115	13
5	CIT	A	2504	13/13	0.86	0.32	115,115,115,115	13
4	AU	A	386	1/1	0.89	0.40	78,78,78,78	1
4	AU	A	387	1/1	0.89	0.25	78,78,78,78	1
4	AU	B	387	1/1	0.89	0.36	78,78,78,78	1
4	AU	B	386	1/1	0.90	0.34	78,78,78,78	1
4	AU	B	388	1/1	0.90	0.20	78,78,78,78	1
5	CIT	B	2502	13/13	0.90	0.28	115,115,115,115	13
4	AU	A	384	1/1	0.94	0.40	78,78,78,78	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	AU	B	391	1/1	0.95	0.17	78,78,78,78	1
4	AU	B	389	1/1	0.95	0.17	78,78,78,78	1
4	AU	B	384	1/1	0.95	0.39	78,78,78,78	1
4	AU	B	390	1/1	0.96	0.18	78,78,78,78	1
3	ZN	B	382	1/1	0.97	0.20	213,213,213,213	0
3	ZN	B	383	1/1	0.97	0.23	213,213,213,213	0
3	ZN	A	383	1/1	0.98	0.24	213,213,213,213	0
4	AU	A	388	1/1	0.98	0.20	78,78,78,78	1
3	ZN	A	382	1/1	0.99	0.19	213,213,213,213	0
4	AU	B	385	1/1	0.99	0.31	78,78,78,78	1
4	AU	A	385	1/1	1.00	0.30	78,78,78,78	1

6.5 Other polymers [i](#)

There are no such residues in this entry.