



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 08:32 AM BST

PDB ID : 4O7P
Title : Crystal structure of Mycobacterium tuberculosis maltose kinase MaK complexed with maltose
Authors : Li, J.; Guan, X.T.; Rao, Z.H.
Deposited on : 2013-12-26
Resolution : 2.90 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

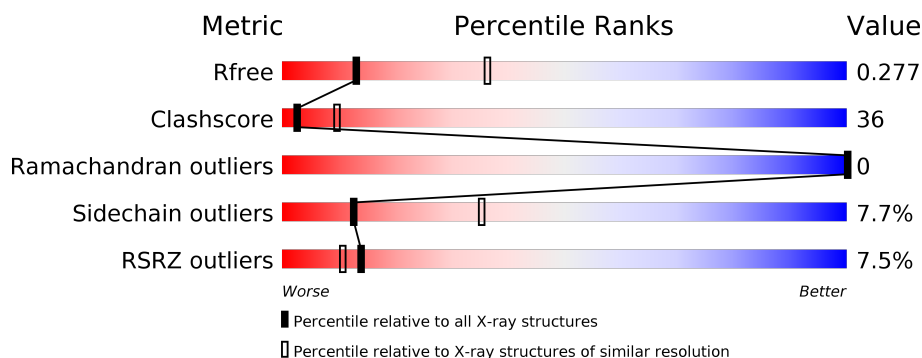
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	455	<div> <div>7%</div> <div> <div></div> <div>57%</div> <div>33%</div> <div>5%</div> <div>5%</div> </div> </div>
1	B	455	<div> <div>7%</div> <div> <div></div> <div>57%</div> <div>38%</div> <div>.</div> <div>.</div> </div> </div>
2	C	2	<div> <div>100%</div> </div>

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6895 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 Maltokinase.

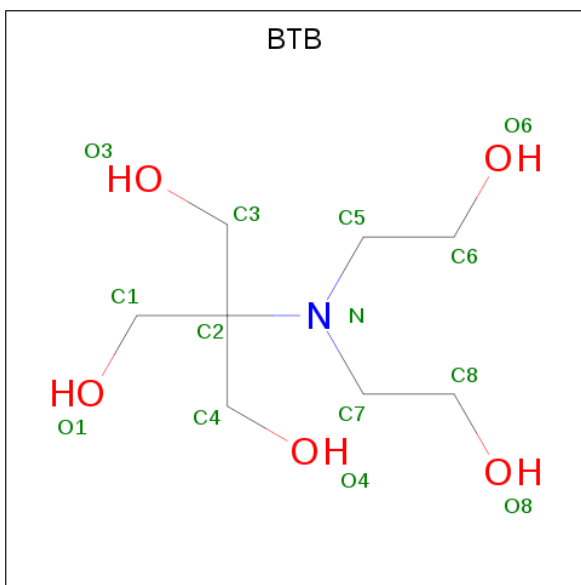
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	433	Total	C	N	O	S	0	0	0
			3349	2114	593	636	6			
1	B	451	Total	C	N	O	S	0	0	0
			3490	2199	618	667	6			

- 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			

- Molecule 3 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (three-letter code: BTB) (formula: C₈H₁₉NO₅).

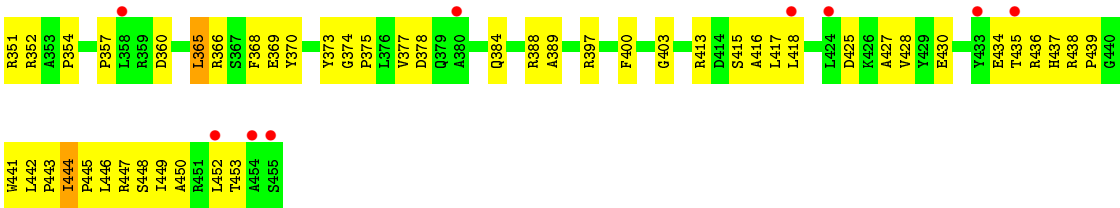


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		

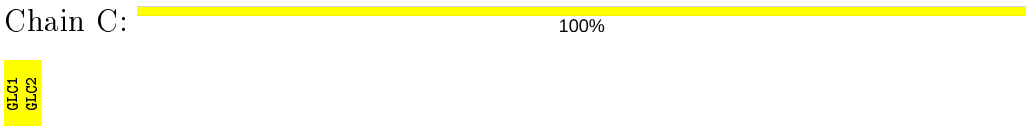
- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	96.73 Å 96.73 Å 461.30 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.60 – 2.90 48.36 – 2.90	Depositor EDS
% Data completeness (in resolution range)	97.2 (44.60-2.90) 99.7 (48.36-2.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.91 Å)	Xtriage
Refinement program	PHENIX 1.7 _650	Depositor
R, R_{free}	0.239 , 0.281 0.235 , 0.277	Depositor DCC
R_{free} test set	1494 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	98.7	Xtriage
Anisotropy	0.219	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 77.5	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	6895	wwPDB-VP
Average B, all atoms (Å ²)	116.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.89% 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: GLC, SO4, BTB

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	0.66	0/3420	0.86	1/4661 (0.0%)
1	B	0.66	0/3568	0.84	2/4867 (0.0%)
All	All	0.66	0/6988	0.85	3/9528 (0.0%)

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	3
1	B	0	1
All	All	0	4

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	265	LEU	CB-CG-CD1	-8.29	96.90	111.00
1	A	138	VAL	N-CA-C	-5.93	95.00	111.00
1	B	329	LEU	CB-CG-CD1	-5.33	101.94	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	142	GLU	Peptide
1	A	293	PRO	Peptide
1	A	344	PRO	Peptide

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Mol	Chain	Res	Type	Group
1	B	115	THR	Peptide

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	3349	0	3283	241	0
1	B	3490	0	3409	253	0
2	C	23	0	19	0	0
3	A	28	0	38	7	0
4	A	5	0	0	0	0
All	All	6895	0	6749	494	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 36.

All (494) 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:198:ARG:HG2	1:B:204:LEU:HD13	1.25	1.17
1:A:10:LYS:NZ	1:A:85:ALA:HB1	1.56	1.17
1:B:303:GLU:HA	1:B:306:PHE:CE2	1.87	1.10
1:A:302:ILE:HG22	1:A:306:PHE:HE2	1.10	1.10
1:A:7:LEU:HG	1:A:10:LYS:HE3	1.26	1.09
1:A:295:LEU:HD23	1:A:298:TYR:HE2	0.96	1.09
1:A:295:LEU:HD23	1:A:298:TYR:CE2	1.87	1.07
1:A:446:LEU:HA	1:A:449:ILE:HD12	1.36	1.05
1:B:289:VAL:HG23	1:B:295:LEU:HD21	1.07	1.05
1:B:289:VAL:HA	1:B:295:LEU:HD11	1.29	1.04
1:B:306:PHE:CE1	1:B:307:GLN:HG2	1.92	1.04
1:A:302:ILE:CG2	1:A:306:PHE:HE2	1.72	1.03
1:A:431:THR:O	1:A:435:THR:HB	1.57	1.01
1:A:442:LEU:O	1:A:446:LEU:HD13	1.61	1.01
1:B:289:VAL:HG21	1:B:296:ARG:HH12	1.24	1.00
1:A:96:ASP:O	1:A:100:PRO:HD3	1.62	0.99
1:A:18:SER:HA	1:A:23:TYR:CD2	1.99	0.98

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:65:LEU:HD11	1:B:103:LEU:HD11	1.44	0.98
1:A:238:LEU:HD12	1:A:239:TYR:N	1.79	0.97
1:A:285:LEU:O	1:A:288:THR:HG22	1.64	0.97
1:B:96:ASP:O	1:B:100:PRO:HD2	1.63	0.97
1:B:198:ARG:HG2	1:B:204:LEU:CD1	1.94	0.97
1:B:289:VAL:CG2	1:B:295:LEU:HD21	1.95	0.96
1:A:302:ILE:HG22	1:A:306:PHE:CE2	2.00	0.96
1:B:306:PHE:HD1	1:B:307:GLN:N	1.63	0.96
1:A:295:LEU:CD2	1:A:298:TYR:HE2	1.79	0.95
1:B:306:PHE:CD1	1:B:307:GLN:N	2.35	0.95
1:A:424:LEU:HD12	1:A:425:ASP:H	1.30	0.95
1:B:303:GLU:HA	1:B:306:PHE:CD2	2.04	0.93
1:B:10:LYS:NZ	1:B:86:ASP:H	1.67	0.91
1:B:306:PHE:HE1	1:B:307:GLN:HG2	1.24	0.91
1:A:10:LYS:HZ3	1:A:85:ALA:HB1	1.34	0.91
1:A:135:GLN:NE2	1:A:136:PRO:HD3	1.84	0.91
1:A:424:LEU:HD12	1:A:425:ASP:N	1.85	0.91
1:A:135:GLN:HG3	1:A:136:PRO:CD	2.00	0.90
1:A:365:LEU:HD12	1:A:397:ARG:HG2	1.51	0.90
1:A:135:GLN:HG3	1:A:136:PRO:HD2	1.52	0.90
1:A:417:LEU:HD13	1:A:418:LEU:HD12	1.54	0.90
1:A:7:LEU:HG	1:A:10:LYS:CE	2.01	0.90
1:A:10:LYS:HZ2	1:A:85:ALA:HB1	1.35	0.90
1:A:69:ASP:C	1:A:88:ARG:HG2	1.90	0.90
1:B:195:ARG:HG3	1:B:198:ARG:NH1	1.87	0.90
1:A:238:LEU:HD12	1:A:239:TYR:H	1.34	0.89
1:B:10:LYS:HZ2	1:B:86:ASP:H	0.89	0.89
1:A:346:GLN:HE21	1:A:350:GLU:HG2	1.39	0.87
1:B:289:VAL:HG23	1:B:295:LEU:CD2	1.99	0.87
1:A:373:TYR:OH	1:A:390:ARG:HG2	1.74	0.87
1:B:301:THR:HG21	1:B:453:THR:HG21	1.55	0.86
1:B:106:LEU:HB3	1:B:122:PHE:CD2	2.11	0.86
1:A:347:PRO:HG2	1:A:350:GLU:HB2	1.57	0.85
1:A:7:LEU:HD21	1:A:68:TRP:CH2	2.12	0.85
1:B:11:LEU:HD12	1:B:85:ALA:HB2	1.58	0.85
1:A:7:LEU:CD2	1:A:68:TRP:CH2	2.60	0.84
1:B:289:VAL:HG21	1:B:296:ARG:NH1	1.91	0.84
1:B:30:LEU:HD11	1:B:33:VAL:HG23	1.57	0.84
1:B:306:PHE:HE1	1:B:307:GLN:CG	1.90	0.84
1:A:451:ARG:NH1	1:A:451:ARG:HB2	1.92	0.84
1:B:436:ARG:O	1:B:436:ARG:HD2	1.78	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:11:LEU:HD11	1:B:68:TRP:HH2	1.42	0.84
1:A:301:THR:O	1:A:305:GLN:HG3	1.79	0.83
1:A:438:ARG:HH22	3:A:501:BTB:H12	1.44	0.83
1:A:10:LYS:NZ	1:A:85:ALA:CB	2.41	0.83
1:B:296:ARG:HH21	1:B:299:ALA:CB	1.90	0.82
1:B:308:LYS:HD2	1:B:308:LYS:O	1.79	0.82
1:A:417:LEU:HD13	1:A:418:LEU:CD1	2.09	0.82
1:A:280:ARG:NH1	1:A:284:ARG:HH12	1.77	0.82
1:B:299:ALA:O	1:B:302:ILE:HG12	1.79	0.82
1:B:370:TYR:CE1	1:B:444:ILE:HD11	2.14	0.81
1:B:445:PRO:O	1:B:449:ILE:HG23	1.81	0.81
1:B:10:LYS:HZ2	1:B:86:ASP:N	1.75	0.81
1:B:276:PHE:HD2	1:B:278:VAL:HG13	1.46	0.81
1:A:302:ILE:CG2	1:A:306:PHE:CE2	2.63	0.80
1:B:296:ARG:NH2	1:B:299:ALA:CB	2.44	0.80
1:B:238:LEU:O	1:B:388:ARG:NH1	2.13	0.80
1:B:286:SER:O	1:B:289:VAL:HG12	1.82	0.80
1:A:451:ARG:HB2	1:A:451:ARG:CZ	2.13	0.79
1:A:346:GLN:NE2	1:A:350:GLU:HG2	1.98	0.79
1:B:198:ARG:CG	1:B:204:LEU:HD13	2.09	0.79
1:B:300:PRO:HA	1:B:303:GLU:OE1	1.82	0.79
1:A:390:ARG:HH22	1:A:447:ARG:HD3	1.46	0.79
1:B:303:GLU:CA	1:B:306:PHE:CE2	2.67	0.78
1:A:7:LEU:CG	1:A:10:LYS:HE3	2.13	0.77
1:A:142:GLU:HA	1:A:143:GLN:C	2.05	0.77
1:B:444:ILE:HG13	1:B:445:PRO:CD	2.14	0.77
1:B:298:TYR:O	1:B:301:THR:HG22	1.84	0.77
1:B:252:TYR:CE1	1:B:403:GLY:HA2	2.19	0.77
1:A:232:LEU:C	1:A:232:LEU:HD13	2.05	0.76
1:B:28:ARG:HG3	1:B:54:TYR:HE2	1.51	0.76
1:A:79:LYS:HE3	1:A:81:ALA:CB	2.14	0.76
1:A:135:GLN:NE2	1:A:136:PRO:CD	2.48	0.76
1:A:417:LEU:N	1:A:417:LEU:HD12	2.01	0.75
1:B:299:ALA:O	1:B:303:GLU:OE1	2.04	0.75
1:B:441:TRP:O	1:B:444:ILE:HG13	1.86	0.75
1:A:96:ASP:O	1:A:100:PRO:CD	2.34	0.75
1:B:427:ALA:HA	1:B:430:GLU:OE1	1.86	0.75
1:B:108:VAL:HG13	1:B:124:ARG:NH1	2.02	0.75
1:B:278:VAL:O	1:B:282:LEU:HD13	1.87	0.75
1:B:384:GLN:OE1	1:B:388:ARG:HG2	1.87	0.74
1:A:79:LYS:HE3	1:A:81:ALA:HB2	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:296:ARG:HH21	1:B:299:ALA:HB3	1.52	0.74
1:A:421:ALA:O	1:A:424:LEU:HD12	1.87	0.74
1:A:304:GLN:HA	1:A:307:GLN:HG2	1.68	0.74
1:A:82:ILE:HG22	1:A:91:PHE:HA	1.68	0.74
1:A:295:LEU:CD2	1:A:298:TYR:CE2	2.61	0.74
1:B:299:ALA:O	1:B:302:ILE:CG1	2.36	0.74
1:A:277:PRO:HG2	1:A:281:MET:CE	2.19	0.73
1:A:370:TYR:OH	1:A:430:GLU:HB3	1.88	0.73
1:A:135:GLN:HE21	1:A:136:PRO:CD	2.02	0.73
1:A:7:LEU:O	1:A:7:LEU:HD23	1.89	0.72
1:A:45:LEU:HD21	1:A:99:GLY:HA2	1.71	0.72
1:A:439:PRO:O	1:A:442:LEU:HD13	1.89	0.72
1:A:135:GLN:CD	1:A:136:PRO:HD3	2.10	0.72
1:A:285:LEU:O	1:A:289:VAL:HG13	1.90	0.72
1:A:280:ARG:HH12	1:A:284:ARG:HH12	1.35	0.71
1:A:428:VAL:CG2	1:A:429:TYR:N	2.53	0.71
1:B:303:GLU:HG3	1:B:306:PHE:HE2	1.56	0.71
1:B:439:PRO:HA	1:B:442:LEU:HD13	1.72	0.71
1:B:106:LEU:HD13	1:B:122:PHE:CE2	2.26	0.70
1:B:307:GLN:OE1	1:B:307:GLN:HA	1.91	0.70
1:A:174:ARG:HD3	1:A:178:ARG:HH21	1.57	0.70
1:B:446:LEU:O	1:B:449:ILE:HG12	1.92	0.70
1:B:13:TRP:O	1:B:17:LEU:HB2	1.92	0.70
1:B:128:VAL:O	1:B:130:LEU:HD13	1.89	0.70
1:A:7:LEU:HD23	1:A:7:LEU:C	2.13	0.69
1:B:369:GLU:HG2	1:B:448:SER:OG	1.93	0.69
1:A:232:LEU:O	1:A:232:LEU:HD13	1.91	0.69
1:B:415:SER:HB3	1:B:418:LEU:HD13	1.72	0.69
1:A:7:LEU:O	1:A:10:LYS:HG2	1.93	0.69
1:A:135:GLN:CG	1:A:136:PRO:CD	2.71	0.69
1:B:122:PHE:HD1	1:B:191:TYR:CB	2.05	0.69
1:B:295:LEU:HD23	1:B:296:ARG:N	2.06	0.69
1:B:65:LEU:HD11	1:B:103:LEU:CD1	2.22	0.69
1:A:97:VAL:C	1:A:100:PRO:HD2	2.12	0.69
1:B:52:VAL:HG22	1:B:54:TYR:CE1	2.28	0.69
1:B:52:VAL:HG22	1:B:54:TYR:HE1	1.58	0.69
1:A:115:THR:HG22	1:A:122:PHE:HB2	1.75	0.69
1:B:122:PHE:HD1	1:B:191:TYR:HB2	1.58	0.68
1:B:106:LEU:HD13	1:B:122:PHE:HE2	1.57	0.68
1:B:301:THR:CG2	1:B:453:THR:HG21	2.23	0.68
1:B:220:GLU:OE2	1:B:222:TRP:HB3	1.93	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:448:SER:O	1:B:452:LEU:HD13	1.93	0.68
1:A:28:ARG:NH1	1:A:54:TYR:CD2	2.62	0.67
1:A:7:LEU:HD22	1:A:68:TRP:CZ2	2.30	0.67
1:A:10:LYS:HZ2	1:A:85:ALA:CB	2.03	0.67
1:B:374:GLY:N	1:B:375:PRO:HD2	2.10	0.67
1:A:417:LEU:CD1	1:A:417:LEU:H	2.07	0.67
1:B:106:LEU:HB3	1:B:122:PHE:CE2	2.29	0.67
1:B:449:ILE:HG13	1:B:450:ALA:N	2.09	0.67
1:A:63:GLN:OE1	1:A:208:LEU:HD22	1.95	0.67
1:B:52:VAL:CG2	1:B:54:TYR:HE1	2.08	0.66
1:B:65:LEU:CD1	1:B:103:LEU:HD11	2.21	0.66
1:B:292:VAL:O	1:B:295:LEU:HD13	1.95	0.66
1:B:347:PRO:HG2	1:B:350:GLU:OE1	1.95	0.66
1:B:373:TYR:HE2	1:B:447:ARG:HD3	1.61	0.66
1:A:424:LEU:O	1:A:428:VAL:HG13	1.96	0.65
1:B:436:ARG:HG3	1:B:437:HIS:CE1	2.32	0.65
1:A:17:LEU:C	1:A:17:LEU:HD23	2.17	0.64
1:A:97:VAL:O	1:A:100:PRO:HD2	1.97	0.64
1:A:18:SER:HA	1:A:23:TYR:HD2	1.60	0.64
1:B:444:ILE:HG13	1:B:445:PRO:HD2	1.78	0.64
1:A:166:ILE:O	1:A:352:ARG:HD2	1.97	0.64
1:A:276:PHE:CD1	1:A:277:PRO:HD2	2.32	0.64
1:B:171:GLU:OE1	1:B:352:ARG:NH1	2.31	0.64
1:B:373:TYR:CE2	1:B:447:ARG:HD3	2.33	0.64
1:B:299:ALA:HA	1:B:302:ILE:HG12	1.80	0.64
1:B:442:LEU:O	1:B:446:LEU:HD13	1.99	0.63
1:A:282:LEU:O	1:A:285:LEU:HB3	1.99	0.63
1:B:195:ARG:HG3	1:B:198:ARG:HH12	1.63	0.63
1:B:295:LEU:CD2	1:B:296:ARG:N	2.62	0.63
1:A:7:LEU:HD22	1:A:68:TRP:CH2	2.33	0.63
1:B:276:PHE:CD2	1:B:278:VAL:HG13	2.31	0.63
1:A:390:ARG:NH2	1:A:447:ARG:HD3	2.14	0.62
1:B:30:LEU:HD13	1:B:31:ALA:N	2.14	0.62
1:A:96:ASP:HB2	1:B:132:PHE:CE1	2.34	0.62
1:A:434:GLU:O	1:A:442:LEU:HD11	2.00	0.62
1:B:300:PRO:CA	1:B:303:GLU:OE1	2.47	0.62
1:B:122:PHE:CD1	1:B:191:TYR:HB2	2.34	0.62
1:B:162:VAL:HG13	1:B:206:TYR:HD1	1.65	0.62
1:B:27:ASN:OD1	1:B:28:ARG:HD2	1.98	0.62
1:B:10:LYS:HB3	1:B:85:ALA:HB1	1.81	0.62
1:B:299:ALA:HB3	1:B:300:PRO:CD	2.30	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:252:TYR:CE1	1:B:403:GLY:CA	2.83	0.61
1:A:309:LEU:HD23	1:A:312:GLU:OE1	2.00	0.61
1:A:428:VAL:HG23	1:A:429:TYR:N	2.16	0.61
1:A:421:ALA:O	1:A:424:LEU:CD1	2.48	0.61
1:A:140:ASP:HB3	1:B:21:ARG:HD2	1.80	0.61
1:B:306:PHE:CE1	1:B:307:GLN:CG	2.71	0.61
1:B:316:VAL:HG13	1:B:354:PRO:HB2	1.83	0.61
1:A:280:ARG:NH1	1:A:284:ARG:NH1	2.46	0.61
1:A:238:LEU:HD12	1:A:239:TYR:HB2	1.83	0.61
1:A:430:GLU:O	1:A:434:GLU:HG2	2.01	0.60
1:A:82:ILE:HG23	1:A:83:GLY:H	1.65	0.60
1:B:298:TYR:O	1:B:301:THR:CG2	2.49	0.60
1:B:96:ASP:O	1:B:100:PRO:CD	2.44	0.60
1:A:424:LEU:CD1	1:A:425:ASP:N	2.62	0.60
1:A:10:LYS:CE	1:A:85:ALA:HB1	2.30	0.60
1:A:10:LYS:HZ3	1:A:85:ALA:CB	2.07	0.60
1:A:282:LEU:CD2	1:A:303:GLU:HG3	2.31	0.60
1:B:444:ILE:HG13	1:B:445:PRO:HD3	1.82	0.60
1:A:232:LEU:CD1	1:A:232:LEU:C	2.70	0.60
1:A:417:LEU:HD12	1:A:417:LEU:H	1.62	0.60
3:A:501:BTB:O1	3:A:501:BTB:H52	2.00	0.60
1:A:425:ASP:O	1:A:428:VAL:HG22	2.01	0.60
1:A:346:GLN:HG3	1:A:347:PRO:HD2	1.84	0.59
1:B:299:ALA:HB3	1:B:300:PRO:HD3	1.84	0.59
1:B:291:VAL:CG2	1:B:435:THR:HG21	2.32	0.59
1:A:282:LEU:HD22	1:A:303:GLU:HG3	1.84	0.59
1:A:278:VAL:HA	1:A:281:MET:HG3	1.83	0.59
1:B:269:LEU:HD23	1:B:318:ARG:HG3	1.84	0.59
1:B:303:GLU:HG3	1:B:306:PHE:CE2	2.36	0.59
1:A:386:ALA:O	1:A:390:ARG:HG3	2.01	0.59
1:A:104:LEU:HD22	1:B:150:PHE:CE2	2.37	0.59
1:B:289:VAL:CG2	1:B:296:ARG:NH1	2.65	0.59
1:B:11:LEU:HD11	1:B:68:TRP:CH2	2.30	0.59
1:A:142:GLU:HA	1:A:144:SER:HA	1.85	0.59
3:A:501:BTB:H42	3:A:501:BTB:O8	2.02	0.58
1:A:17:LEU:HD21	1:A:23:TYR:CD1	2.37	0.58
1:A:288:THR:HG23	1:A:289:VAL:N	2.18	0.58
1:A:277:PRO:HG2	1:A:281:MET:HE1	1.86	0.58
1:B:30:LEU:HD11	1:B:33:VAL:CG2	2.32	0.58
1:B:444:ILE:CD1	1:B:445:PRO:HD3	2.34	0.58
1:A:427:ALA:O	1:A:431:THR:HG23	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:370:TYR:CE1	1:B:444:ILE:CD1	2.85	0.58
1:B:28:ARG:HG3	1:B:54:TYR:CE2	2.36	0.58
1:B:441:TRP:O	1:B:444:ILE:CG1	2.51	0.58
1:A:362:ALA:O	1:A:366:ARG:HG2	2.03	0.58
1:A:275:THR:HG22	1:A:276:PHE:O	2.04	0.57
1:A:28:ARG:NH1	1:A:54:TYR:CG	2.73	0.57
1:A:276:PHE:CD2	1:A:309:LEU:HB2	2.39	0.57
1:A:37:VAL:HG13	1:A:49:LEU:HB2	1.85	0.57
1:B:373:TYR:CE1	1:B:389:ALA:HB1	2.40	0.57
1:A:42:ARG:HE	1:A:45:LEU:HD13	1.69	0.57
1:B:299:ALA:C	1:B:303:GLU:OE1	2.43	0.57
1:A:111:ALA:HB2	3:A:502:BTB:H62	1.87	0.57
1:A:292:VAL:O	1:A:292:VAL:HG13	2.04	0.57
1:B:122:PHE:CD1	1:B:191:TYR:CB	2.88	0.57
1:A:232:LEU:HD12	1:A:233:PHE:CD1	2.40	0.56
1:B:339:ASP:OD1	1:B:341:GLU:HG2	2.05	0.56
1:A:104:LEU:HD12	1:A:211:VAL:HG21	1.88	0.56
1:A:309:LEU:CD2	1:A:312:GLU:OE1	2.53	0.56
1:B:47:LEU:HD23	1:B:47:LEU:C	2.26	0.56
1:A:174:ARG:HD3	1:A:178:ARG:NH2	2.19	0.56
1:B:47:LEU:HD23	1:B:48:VAL:N	2.20	0.56
1:B:16:TRP:CD1	1:B:82:ILE:HA	2.40	0.56
1:B:199:SER:OG	1:B:201:THR:HG23	2.06	0.56
1:B:308:LYS:HG3	1:B:309:LEU:HD23	1.86	0.56
1:B:65:LEU:CD1	1:B:103:LEU:CD1	2.83	0.56
1:A:277:PRO:O	1:A:281:MET:HG2	2.05	0.56
1:B:434:GLU:OE1	1:B:438:ARG:HB2	2.05	0.56
1:A:79:LYS:HE3	1:A:81:ALA:HB3	1.86	0.56
1:B:306:PHE:HE1	1:B:307:GLN:CD	2.09	0.56
1:A:434:GLU:HB2	1:A:438:ARG:HB3	1.88	0.55
1:A:15:ASP:OD1	1:A:15:ASP:N	2.39	0.55
1:A:293:PRO:O	1:A:296:ARG:HG3	2.07	0.55
1:B:302:ILE:HG13	1:B:303:GLU:N	2.20	0.55
1:A:440:GLY:O	1:A:443:PRO:HD2	2.06	0.55
1:A:393:VAL:O	1:A:397:ARG:HG3	2.07	0.55
1:B:303:GLU:CG	1:B:306:PHE:CE2	2.90	0.55
1:B:30:LEU:HD21	1:B:52:VAL:HG23	1.89	0.55
1:A:347:PRO:HG2	1:A:350:GLU:CB	2.32	0.54
1:A:302:ILE:O	1:A:306:PHE:CD2	2.61	0.54
1:A:438:ARG:HG2	1:A:441:TRP:CD1	2.42	0.54
1:B:303:GLU:O	1:B:306:PHE:CD1	2.61	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:LEU:HG	1:A:297:GLU:HG3	1.89	0.54
1:A:435:THR:CG2	1:A:436:ARG:N	2.70	0.54
1:B:54:TYR:HD1	1:B:54:TYR:N	2.04	0.54
1:A:435:THR:HG22	1:A:436:ARG:N	2.23	0.54
1:A:441:TRP:CD1	1:A:441:TRP:N	2.76	0.54
1:B:159:PHE:HD2	1:B:208:LEU:C	2.11	0.54
1:B:291:VAL:CG2	1:B:291:VAL:O	2.56	0.54
1:B:370:TYR:CZ	1:B:444:ILE:HD11	2.42	0.54
1:B:373:TYR:HE1	1:B:389:ALA:CB	2.20	0.54
1:A:238:LEU:CD1	1:A:239:TYR:H	2.13	0.54
1:A:415:SER:C	1:A:417:LEU:HD12	2.28	0.54
1:B:333:GLU:HG2	1:B:334:SER:H	1.72	0.54
1:A:162:VAL:CG1	1:A:206:TYR:HD1	2.21	0.54
3:A:501:BTB:H42	3:A:501:BTB:C8	2.35	0.53
1:B:201:THR:O	1:B:202:ASP:HB2	2.08	0.53
1:B:30:LEU:HD21	1:B:33:VAL:HG22	1.89	0.53
1:A:150:PHE:HZ	1:B:100:PRO:HB2	1.74	0.53
1:B:282:LEU:N	1:B:282:LEU:HD13	2.23	0.53
1:B:5:ASP:OD1	1:B:5:ASP:N	2.41	0.53
1:B:296:ARG:O	1:B:300:PRO:HD2	2.09	0.53
1:B:278:VAL:O	1:B:282:LEU:HD22	2.08	0.53
1:B:282:LEU:CD1	1:B:282:LEU:N	2.71	0.53
1:B:320:HIS:HE1	1:B:322:ASP:O	1.92	0.53
1:B:373:TYR:CE1	1:B:389:ALA:CB	2.92	0.53
1:B:54:TYR:N	1:B:54:TYR:CD1	2.73	0.53
1:A:232:LEU:HD11	1:A:388:ARG:HB3	1.91	0.53
1:A:43:HIS:O	1:A:44:ASN:HB2	2.09	0.53
1:A:135:GLN:CG	1:A:136:PRO:HD3	2.39	0.52
1:A:349:ASP:N	1:A:349:ASP:OD1	2.38	0.52
1:B:128:VAL:O	1:B:130:LEU:CD1	2.56	0.52
1:B:317:GLN:O	1:B:354:PRO:HA	2.08	0.52
1:A:446:LEU:CA	1:A:449:ILE:HD12	2.24	0.52
1:A:377:VAL:O	1:A:379:GLN:OE1	2.27	0.52
1:A:280:ARG:O	1:A:283:ALA:HB3	2.10	0.52
1:A:427:ALA:HA	1:A:430:GLU:OE2	2.10	0.52
1:A:7:LEU:CD2	1:A:7:LEU:C	2.78	0.52
1:B:370:TYR:OH	1:B:430:GLU:HB3	2.11	0.51
1:B:295:LEU:HD23	1:B:295:LEU:C	2.30	0.51
1:A:295:LEU:HD11	1:A:297:GLU:OE2	2.10	0.51
1:A:451:ARG:NH1	1:A:451:ARG:CB	2.71	0.51
1:B:444:ILE:CG1	1:B:445:PRO:HD3	2.40	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:428:VAL:HG22	1:A:429:TYR:H	1.74	0.51
1:A:445:PRO:O	1:A:449:ILE:HG13	2.11	0.51
1:A:302:ILE:HG23	1:A:306:PHE:CE2	2.45	0.51
1:A:49:LEU:HB3	1:A:61:ARG:HD3	1.92	0.51
1:B:303:GLU:CB	1:B:306:PHE:CE2	2.94	0.51
1:B:425:ASP:O	1:B:428:VAL:HB	2.11	0.51
1:A:140:ASP:CB	1:B:21:ARG:HD2	2.41	0.50
1:B:276:PHE:CD2	1:B:278:VAL:CG1	2.94	0.50
1:A:174:ARG:HG2	1:A:178:ARG:HD2	1.94	0.50
1:B:137:ARG:HB2	1:B:149:ILE:HB	1.93	0.50
1:B:142:GLU:HG3	1:B:144:SER:H	1.75	0.50
1:B:122:PHE:HE1	1:B:191:TYR:CD1	2.29	0.50
1:B:296:ARG:NH2	1:B:299:ALA:HB2	2.26	0.50
1:B:130:LEU:HB3	1:B:131:PRO:HD2	1.93	0.50
1:B:299:ALA:C	1:B:302:ILE:HG12	2.33	0.50
1:A:442:LEU:N	1:A:442:LEU:CD1	2.75	0.49
1:A:297:GLU:O	1:A:300:PRO:HD2	2.12	0.49
1:A:418:LEU:H	1:A:418:LEU:HD12	1.76	0.49
1:A:93:ALA:HB1	1:A:99:GLY:HA3	1.93	0.49
1:B:48:VAL:HG12	1:B:50:VAL:HG23	1.92	0.49
1:B:377:VAL:HG22	1:B:378:ASP:H	1.76	0.49
1:B:436:ARG:C	1:B:436:ARG:HD2	2.32	0.49
1:B:51:ASP:OD2	1:B:61:ARG:HD2	2.13	0.49
1:B:370:TYR:CZ	1:B:444:ILE:CD1	2.96	0.49
1:A:135:GLN:CG	1:A:136:PRO:HD2	2.34	0.49
1:A:142:GLU:HG3	1:A:144:SER:HB3	1.95	0.49
1:A:238:LEU:HD12	1:A:239:TYR:CB	2.42	0.49
1:A:447:ARG:HG2	1:A:451:ARG:HH12	1.77	0.49
1:B:108:VAL:CG1	1:B:124:ARG:NH1	2.74	0.48
1:B:444:ILE:CG1	1:B:445:PRO:CD	2.89	0.48
1:B:84:VAL:HA	1:B:88:ARG:O	2.13	0.48
1:A:288:THR:CG2	1:A:289:VAL:N	2.76	0.48
1:B:129:GLU:CD	1:B:129:GLU:H	2.17	0.48
1:B:299:ALA:O	1:B:302:ILE:HG13	2.11	0.48
1:A:278:VAL:O	1:A:281:MET:HG3	2.13	0.48
1:B:250:GLU:OE1	1:B:330:ARG:NH1	2.43	0.48
1:A:428:VAL:CG2	1:A:429:TYR:H	2.26	0.48
1:A:142:GLU:HA	1:A:144:SER:N	2.28	0.48
1:A:174:ARG:CD	1:A:178:ARG:HH21	2.25	0.47
1:B:162:VAL:HG13	1:B:206:TYR:CD1	2.47	0.47
1:B:118:GLY:HA3	1:B:193:PHE:CZ	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:LEU:HD22	1:A:33:VAL:HG12	1.94	0.47
1:B:299:ALA:CA	1:B:302:ILE:HG12	2.41	0.47
1:B:52:VAL:CG2	1:B:54:TYR:CE1	2.91	0.47
1:B:252:TYR:CD1	1:B:403:GLY:CA	2.97	0.47
1:B:373:TYR:HE1	1:B:389:ALA:HB1	1.77	0.47
1:A:344:PRO:HA	1:A:351:ARG:HH21	1.79	0.47
1:B:384:GLN:O	1:B:384:GLN:OE1	2.32	0.47
1:A:162:VAL:HG12	1:A:206:TYR:HD1	1.80	0.47
1:A:346:GLN:HG3	1:A:347:PRO:CD	2.43	0.47
1:B:309:LEU:HD22	1:B:417:LEU:HG	1.95	0.47
1:A:242:GLU:HA	1:A:242:GLU:OE1	2.14	0.47
1:B:314:ILE:HG22	1:B:315:THR:O	2.14	0.47
1:B:444:ILE:HD12	1:B:445:PRO:HD3	1.97	0.47
1:A:276:PHE:CE2	1:A:309:LEU:HB2	2.50	0.47
1:A:33:VAL:O	1:A:33:VAL:CG2	2.63	0.47
1:A:319:VAL:HB	1:A:360:ASP:OD2	2.14	0.47
1:B:112:VAL:HG22	1:B:121:THR:HG23	1.96	0.47
1:B:441:TRP:C	1:B:443:PRO:HD2	2.36	0.47
1:A:281:MET:HG2	1:A:281:MET:H	1.35	0.47
1:B:306:PHE:C	1:B:306:PHE:CD1	2.88	0.47
1:B:442:LEU:N	1:B:443:PRO:CD	2.78	0.47
1:A:439:PRO:C	1:A:442:LEU:HD13	2.34	0.47
1:B:176:LEU:HD13	1:B:184:VAL:HG11	1.98	0.46
1:B:199:SER:HA	1:B:200:PRO:HD3	1.64	0.46
1:B:159:PHE:CE2	1:B:209:GLY:HA2	2.50	0.46
1:B:175:VAL:HG21	1:B:269:LEU:HD11	1.97	0.46
1:B:198:ARG:O	1:B:199:SER:HB3	2.15	0.46
1:B:230:ARG:HG2	1:B:230:ARG:HH21	1.79	0.46
1:B:350:GLU:OE1	1:B:350:GLU:N	2.48	0.46
1:A:174:ARG:CG	1:A:178:ARG:HD2	2.45	0.46
1:A:20:GLN:HG3	1:A:22:TRP:CZ2	2.51	0.46
1:A:237:ASP:HA	1:A:240:ALA:CB	2.45	0.46
1:A:415:SER:O	1:A:417:LEU:HD12	2.16	0.46
1:A:32:THR:CG2	1:A:53:THR:OG1	2.64	0.46
1:A:82:ILE:HG23	1:A:83:GLY:N	2.30	0.46
1:B:34:LYS:HA	1:B:35:PRO:HD3	1.79	0.46
1:B:159:PHE:CD2	1:B:209:GLY:HA2	2.51	0.46
1:B:434:GLU:CB	1:B:442:LEU:HD12	2.46	0.45
1:A:390:ARG:HH12	1:A:447:ARG:HD3	1.80	0.45
1:B:139:CYS:SG	1:B:149:ILE:HD11	2.57	0.45
1:A:113:CYS:O	1:A:113:CYS:SG	2.74	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:377:VAL:O	1:A:378:ASP:HB2	2.15	0.45
1:A:436:ARG:HG3	1:A:437:HIS:CE1	2.51	0.45
1:B:10:LYS:O	1:B:10:LYS:HG2	2.16	0.45
1:B:241:HIS:CE1	1:B:242:GLU:HG2	2.51	0.45
1:B:252:TYR:CD1	1:B:403:GLY:N	2.84	0.45
1:A:170:ILE:HD13	1:A:190:ALA:HB1	1.97	0.45
1:B:166:ILE:HD12	1:B:352:ARG:NH2	2.31	0.45
1:B:286:SER:HA	1:B:289:VAL:HG12	1.99	0.45
1:A:142:GLU:HA	1:A:144:SER:CA	2.44	0.45
1:A:276:PHE:CE1	1:A:281:MET:SD	3.10	0.45
1:B:276:PHE:CD2	1:B:276:PHE:C	2.90	0.45
1:B:374:GLY:N	1:B:375:PRO:CD	2.77	0.45
1:B:436:ARG:HG3	1:B:437:HIS:ND1	2.31	0.45
1:B:276:PHE:HA	1:B:277:PRO:HD3	1.61	0.45
1:A:365:LEU:CD1	1:A:397:ARG:HG2	2.37	0.45
1:B:269:LEU:N	1:B:269:LEU:CD1	2.80	0.45
1:A:137:ARG:HB3	1:A:149:ILE:HD12	1.99	0.44
1:A:358:LEU:HD23	1:A:419:LEU:HB2	1.99	0.44
1:A:252:TYR:CD1	1:A:252:TYR:C	2.91	0.44
1:A:434:GLU:HG3	1:A:442:LEU:HD12	2.00	0.44
1:A:446:LEU:HA	1:A:449:ILE:CD1	2.26	0.44
1:B:262:HIS:ND1	1:B:357:PRO:HB3	2.32	0.44
1:A:15:ASP:O	1:A:19:ARG:HG3	2.17	0.44
1:B:295:LEU:C	1:B:295:LEU:CD2	2.85	0.44
1:A:235:GLU:HG3	1:A:236:GLY:O	2.17	0.44
1:A:278:VAL:CA	1:A:281:MET:HG3	2.48	0.44
1:A:304:GLN:CA	1:A:307:GLN:HG2	2.45	0.44
1:B:292:VAL:O	1:B:292:VAL:HG12	2.17	0.44
1:B:306:PHE:CE1	1:B:307:GLN:CD	2.90	0.44
1:A:135:GLN:HE21	1:A:136:PRO:HD2	1.78	0.44
1:A:176:LEU:HG	1:A:184:VAL:HG11	1.99	0.44
1:B:276:PHE:HB2	1:B:314:ILE:HD11	2.00	0.44
1:B:317:GLN:O	1:B:354:PRO:CA	2.66	0.44
1:A:338:ILE:O	1:A:338:ILE:HG13	2.18	0.44
1:B:122:PHE:HD1	1:B:191:TYR:HB3	1.79	0.44
1:B:24:ALA:HA	1:B:25:GLY:HA2	1.73	0.44
1:B:320:HIS:CE1	1:B:322:ASP:O	2.71	0.44
1:B:300:PRO:O	1:B:303:GLU:HB2	2.18	0.44
1:B:331:THR:HB	1:B:332:PRO:CD	2.48	0.44
1:B:446:LEU:O	1:B:449:ILE:CG1	2.65	0.44
1:B:162:VAL:CG1	1:B:206:TYR:CD1	3.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:384:GLN:O	1:A:387:ALA:HB3	2.18	0.43
3:A:502:BTB:H72	3:A:502:BTB:O6	2.18	0.43
1:B:187:LEU:HD13	1:B:210:MET:HE3	2.00	0.43
1:A:30:LEU:HD22	1:A:33:VAL:CG1	2.48	0.43
1:B:224:MET:HE2	1:B:247:PHE:HE1	1.83	0.43
1:B:262:HIS:CE1	1:B:360:ASP:HB2	2.53	0.43
1:B:397:ARG:HH21	1:B:397:ARG:HD2	1.69	0.43
1:B:434:GLU:HB2	1:B:442:LEU:HD12	2.00	0.43
1:B:447:ARG:O	1:B:450:ALA:HB3	2.19	0.43
1:A:365:LEU:HD11	1:A:397:ARG:HA	2.00	0.43
1:B:333:GLU:HG2	1:B:334:SER:N	2.33	0.43
1:B:168:PRO:HD3	1:B:348:LEU:CD1	2.48	0.43
1:A:65:LEU:HD23	1:A:93:ALA:HB3	2.00	0.43
1:A:136:PRO:HG2	1:A:149:ILE:O	2.18	0.43
1:B:166:ILE:O	1:B:348:LEU:HD11	2.19	0.43
1:B:279:ASP:HA	1:B:282:LEU:HD22	2.00	0.42
1:B:306:PHE:C	1:B:306:PHE:HD1	2.21	0.42
1:B:365:LEU:O	1:B:368:PHE:HB2	2.19	0.42
1:A:276:PHE:CD2	1:A:309:LEU:CB	3.01	0.42
1:A:52:VAL:HG12	1:A:54:TYR:CE2	2.54	0.42
1:B:341:GLU:O	1:B:351:ARG:HD2	2.19	0.42
1:B:10:LYS:NZ	1:B:86:ASP:N	2.49	0.42
1:B:22:TRP:CG	1:B:160:ARG:HG3	2.55	0.42
1:A:373:TYR:HD1	1:A:376:LEU:HD12	1.85	0.42
1:B:285:LEU:HD11	1:B:302:ILE:HD11	2.01	0.42
1:B:255:GLY:HA2	1:B:400:PHE:CE1	2.54	0.42
1:A:17:LEU:C	1:A:17:LEU:CD2	2.87	0.42
1:B:306:PHE:HD1	1:B:307:GLN:CA	2.30	0.42
1:B:30:LEU:CD2	1:B:52:VAL:HG23	2.49	0.42
1:A:22:TRP:CG	1:A:160:ARG:HG3	2.54	0.42
3:A:502:BTB:H31	3:A:502:BTB:H52	1.95	0.42
1:A:277:PRO:O	1:A:281:MET:CG	2.68	0.42
1:A:299:ALA:N	1:A:300:PRO:HD2	2.35	0.42
1:A:278:VAL:C	1:A:281:MET:HG3	2.40	0.42
1:A:427:ALA:O	1:A:430:GLU:HB2	2.19	0.42
1:A:365:LEU:O	1:A:368:PHE:HB2	2.20	0.42
1:A:442:LEU:N	1:A:442:LEU:HD12	2.35	0.42
1:B:366:ARG:HH12	1:B:448:SER:CB	2.33	0.42
1:A:291:VAL:O	1:A:291:VAL:HG12	2.20	0.41
1:B:417:LEU:O	1:B:417:LEU:HD12	2.20	0.41
1:A:17:LEU:CD2	1:A:23:TYR:CD1	3.04	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:442:LEU:O	1:A:446:LEU:CD1	2.51	0.41
1:A:10:LYS:HZ3	1:A:85:ALA:C	2.23	0.41
1:A:285:LEU:O	1:A:289:VAL:CG1	2.65	0.41
1:A:435:THR:HG22	1:A:436:ARG:H	1.85	0.41
1:B:298:TYR:O	1:B:299:ALA:C	2.58	0.41
1:B:413:ARG:O	1:B:416:ALA:HB2	2.20	0.41
1:A:233:PHE:CZ	1:A:376:LEU:HD21	2.56	0.41
1:B:151:ASP:C	1:B:152:ARG:HG2	2.40	0.41
1:B:370:TYR:CD1	1:B:444:ILE:CD1	3.04	0.41
1:B:28:ARG:HB2	1:B:54:TYR:CD2	2.56	0.41
1:A:294:GLU:OE1	1:A:295:LEU:N	2.54	0.41
1:A:424:LEU:HB3	1:A:452:LEU:HD13	2.03	0.41
1:B:285:LEU:CD1	1:B:302:ILE:HD11	2.51	0.41
1:A:302:ILE:HG21	1:A:302:ILE:HD13	1.84	0.41
1:B:7:LEU:HA	1:B:7:LEU:HD12	1.93	0.41
1:A:374:GLY:N	1:A:375:PRO:CD	2.83	0.41
1:A:417:LEU:CD1	1:A:418:LEU:CD1	2.91	0.41
1:A:442:LEU:H	1:A:442:LEU:CD1	2.34	0.40
1:A:208:LEU:HD23	1:A:208:LEU:HA	1.83	0.40
1:A:294:GLU:OE1	1:A:294:GLU:C	2.59	0.40
1:A:418:LEU:H	1:A:418:LEU:CD1	2.34	0.40
1:B:52:VAL:CG1	1:B:60:GLU:HB2	2.52	0.40
1:B:251:SER:O	1:B:400:PHE:HA	2.21	0.40
1:B:331:THR:HB	1:B:332:PRO:HD2	2.03	0.40
1:A:346:GLN:HG3	1:A:350:GLU:HB3	2.03	0.40
1:B:373:TYR:CD1	1:B:373:TYR:N	2.89	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	427/455 (94%)	409 (96%)	18 (4%)	0	100	100
1	B	449/455 (99%)	433 (96%)	16 (4%)	0	100	100
All	All	876/910 (96%)	842 (96%)	34 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	342/362 (94%)	314 (92%)	28 (8%)	11	32
1	B	358/362 (99%)	332 (93%)	26 (7%)	14	38
All	All	700/724 (97%)	646 (92%)	54 (8%)	13	35

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	7	LEU
1	A	15	ASP
1	A	33	VAL
1	A	37	VAL
1	A	41	LEU
1	A	50	VAL
1	A	53	THR
1	A	56	ASP
1	A	129	GLU
1	A	140	ASP
1	A	144	SER
1	A	176	LEU
1	A	232	LEU
1	A	239	TYR
1	A	246	ASP
1	A	281	MET
1	A	301	THR

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Mol	Chain	Res	Type
1	A	316	VAL
1	A	349	ASP
1	A	365	LEU
1	A	369	GLU
1	A	417	LEU
1	A	424	LEU
1	A	428	VAL
1	A	435	THR
1	A	436	ARG
1	A	448	SER
1	B	5	ASP
1	B	17	LEU
1	B	30	LEU
1	B	55	THR
1	B	61	ARG
1	B	117	THR
1	B	121	THR
1	B	123	THR
1	B	127	ASP
1	B	128	VAL
1	B	129	GLU
1	B	135	GLN
1	B	217	ASN
1	B	232	LEU
1	B	242	GLU
1	B	269	LEU
1	B	282	LEU
1	B	284	ARG
1	B	286	SER
1	B	291	VAL
1	B	295	LEU
1	B	306	PHE
1	B	309	LEU
1	B	316	VAL
1	B	365	LEU
1	B	444	ILE

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	346	GLN

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$
2	GLC	C	1	2	12,12,12	2.09	4 (33%)	17,17,17	1.28	2 (11%)
2	GLC	C	2	2	11,11,12	2.60	3 (27%)	15,15,17	1.40	2 (13%)

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	-	2/2/22/22	0/1/1/1
2	GLC	C	2	2	-	2/2/19/22	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	2	GLC	C2-C3	-6.84	1.42	1.52
2	C	1	GLC	C3-C2	-4.29	1.41	1.52
2	C	1	GLC	O5-C1	3.29	1.51	1.42
2	C	2	GLC	C4-C3	-3.19	1.44	1.52
2	C	2	GLC	O5-C1	-2.90	1.39	1.43
2	C	1	GLC	O5-C5	-2.30	1.38	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	GLC	O3-C3	-2.28	1.37	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2	GLC	O2-C2-C1	3.14	115.57	109.15
2	C	1	GLC	O3-C3-C4	-3.07	103.25	110.35
2	C	2	GLC	O2-C2-C3	-3.00	104.13	110.14
2	C	1	GLC	O5-C5-C4	2.16	113.62	109.69

There are no chirality outliers.

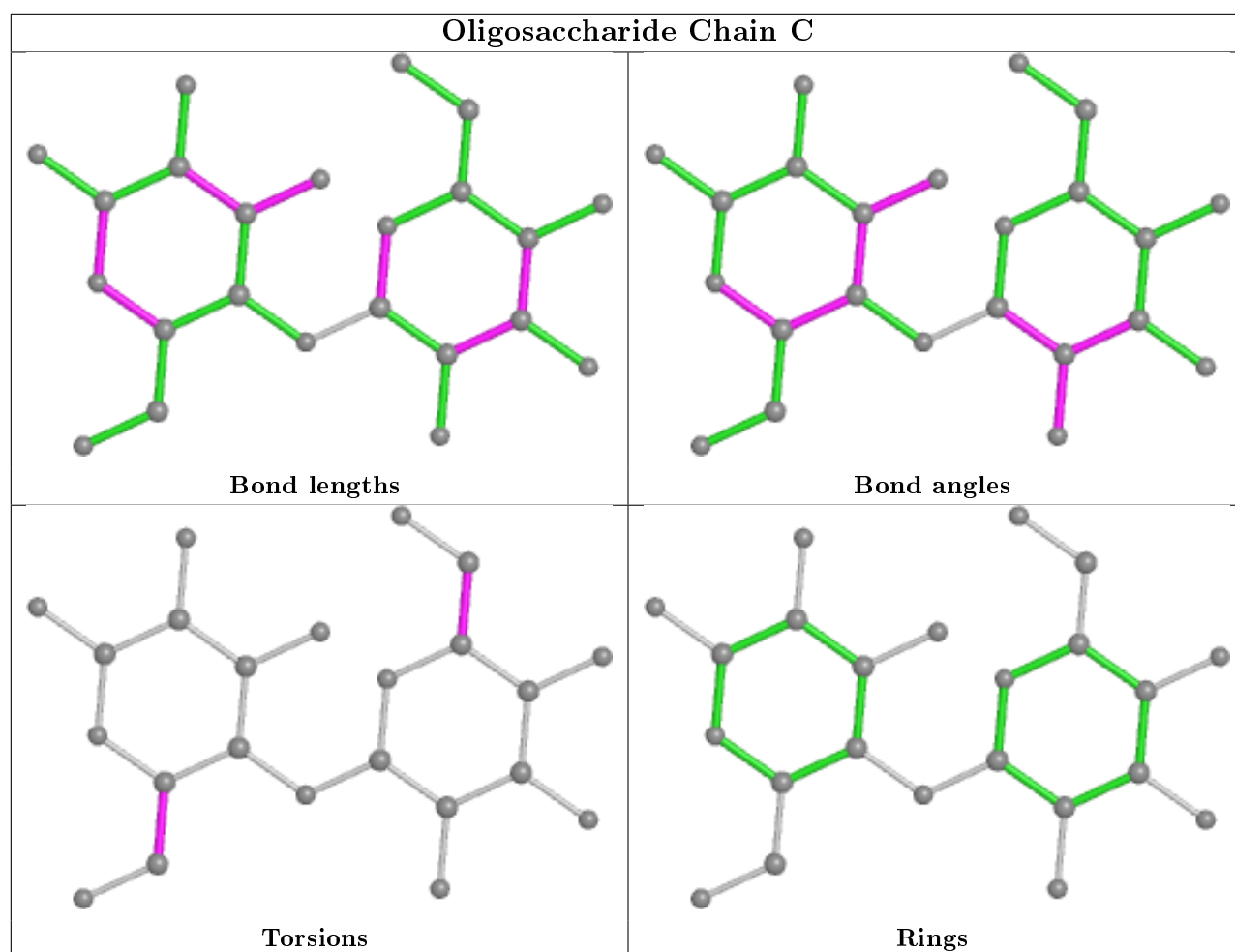
All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	2	GLC	O5-C5-C6-O6
2	C	2	GLC	C4-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	C	1	GLC	O5-C5-C6-O6

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](#)

3 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	BTB	A	502	-	13,13,13	1.50	4 (30%)	7,16,16	2.23	2 (28%)
3	BTB	A	501	-	13,13,13	0.60	0	7,16,16	3.09	5 (71%)
4	SO4	A	503	-	4,4,4	0.23	0	6,6,6	0.36	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	BTB	A	502	-	-	9/21/21/21	-
3	BTB	A	501	-	-	6/21/21/21	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	BTB	C3-C2	-2.48	1.50	1.53
3	A	502	BTB	C1-C2	-2.37	1.50	1.53
3	A	502	BTB	C4-C2	-2.28	1.50	1.53
3	A	502	BTB	C5-N	-2.06	1.45	1.48

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	501	BTB	O3-C3-C2	4.39	123.45	111.44
3	A	501	BTB	O4-C4-C2	4.28	123.17	111.44
3	A	501	BTB	O1-C1-C2	3.43	120.82	111.44
3	A	502	BTB	O4-C4-C2	3.30	120.46	111.44
3	A	502	BTB	O1-C1-C2	3.26	120.36	111.44
3	A	501	BTB	O6-C6-C5	2.31	120.75	111.19
3	A	501	BTB	C6-C5-N	2.23	120.31	111.59

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	BTB	C1-C2-C3-O3
3	A	502	BTB	C4-C2-C3-O3
3	A	502	BTB	N-C2-C3-O3
3	A	502	BTB	C1-C2-C4-O4
3	A	502	BTB	C3-C2-C4-O4
3	A	502	BTB	N-C2-C4-O4
3	A	502	BTB	C6-C5-N-C7
3	A	501	BTB	C1-C2-C4-O4
3	A	501	BTB	C3-C2-C4-O4
3	A	501	BTB	N-C2-C4-O4
3	A	502	BTB	N-C7-C8-O8
3	A	501	BTB	N-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	A	501	BTB	C1-C2-C3-O3
3	A	501	BTB	N-C2-C3-O3
3	A	502	BTB	N-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	502	BTB	3	0
3	A	501	BTB	4	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	433/455 (95%)	0.57	33 (7%) 13 10	68, 113, 168, 196	0
1	B	451/455 (99%)	0.54	33 (7%) 15 11	60, 107, 177, 205	0
All	All	884/910 (97%)	0.55	66 (7%) 14 11	60, 110, 174, 205	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	197	ASN	9.5
1	A	241	HIS	9.0
1	B	380	ALA	8.0
1	B	235	GLU	8.0
1	A	240	ALA	7.4
1	B	196	PRO	6.1
1	B	455	SER	5.6
1	A	30	LEU	5.5
1	A	57	GLY	4.8
1	A	203	ALA	4.7
1	B	238	LEU	4.5
1	B	237	ASP	4.4
1	B	234	ALA	4.4
1	B	454	ALA	4.3
1	A	54	TYR	4.2
1	B	232	LEU	4.1
1	A	204	LEU	3.9
1	B	236	GLY	3.8
1	A	112	VAL	3.7
1	A	237	ASP	3.6
1	B	288	THR	3.5
1	A	376	LEU	3.4
1	B	280	ARG	3.3
1	A	53	THR	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	284	ARG	3.3
1	A	378	ASP	3.3
1	A	26	ARG	3.3
1	A	13	TRP	3.3
1	A	36	GLY	3.3
1	A	113	CYS	3.2
1	B	240	ALA	3.1
1	A	143	GLN	3.1
1	A	243	VAL	3.1
1	B	319	VAL	3.1
1	B	358	LEU	3.0
1	B	133	ALA	3.0
1	A	14	SER	3.0
1	A	120	VAL	2.9
1	A	291	VAL	2.8
1	B	424	LEU	2.8
1	B	418	LEU	2.8
1	B	435	THR	2.8
1	B	132	PHE	2.7
1	B	243	VAL	2.7
1	B	239	TYR	2.5
1	A	377	VAL	2.5
1	A	239	TYR	2.5
1	A	17	LEU	2.5
1	B	26	ARG	2.4
1	A	62	TYR	2.4
1	B	304	GLN	2.3
1	A	114	GLY	2.3
1	A	392	TRP	2.2
1	B	19	ARG	2.2
1	B	452	LEU	2.2
1	B	297	GLU	2.2
1	B	298	TYR	2.2
1	A	236	GLY	2.1
1	B	295	LEU	2.1
1	A	83	GLY	2.1
1	B	244	GLY	2.1
1	A	87	ASP	2.0
1	A	25	GLY	2.0
1	A	380	ALA	2.0
1	B	433	TYR	2.0
1	A	52	VAL	2.0

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

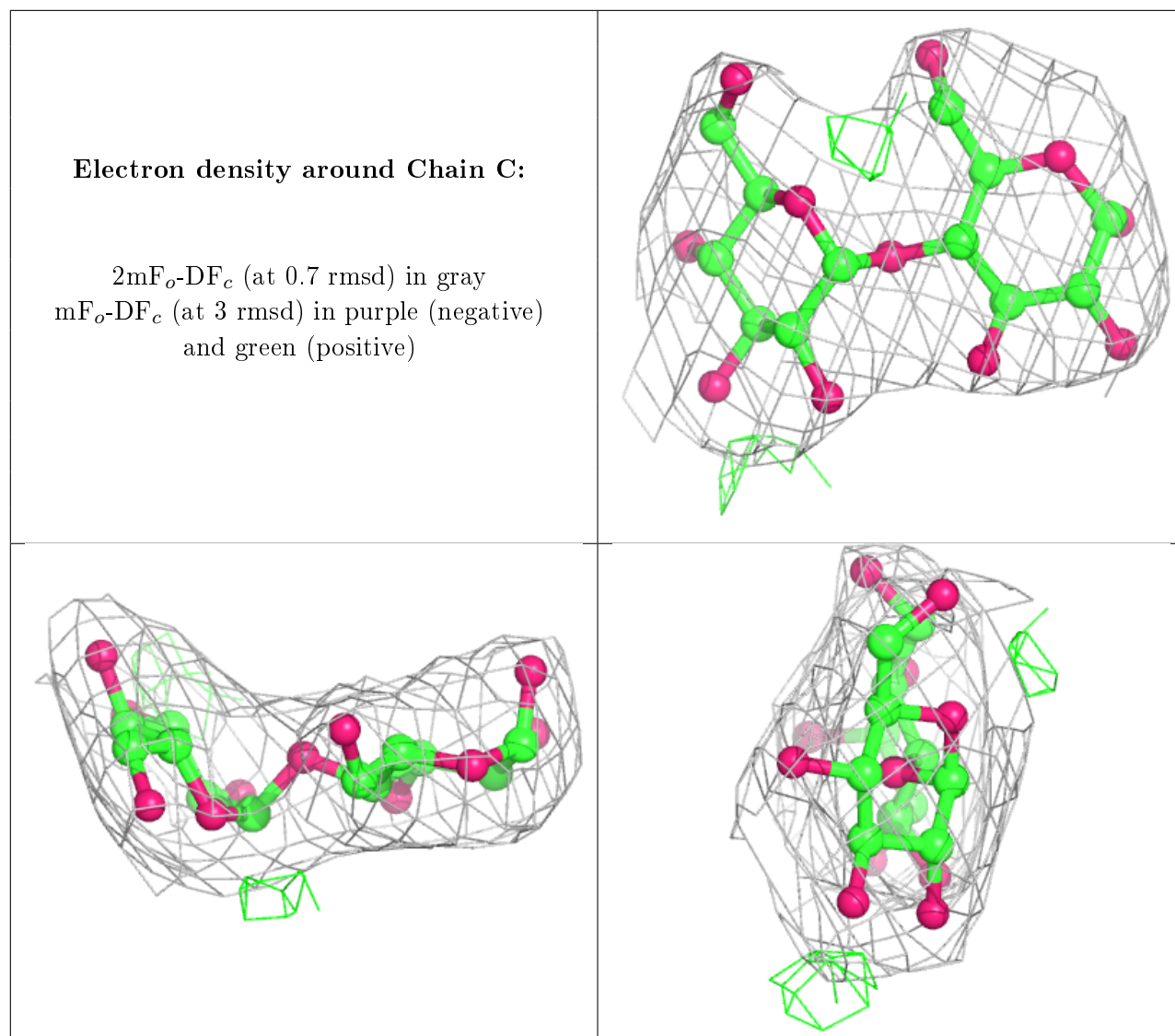
There are no non-standard protein/DNA/RNA residues in this entry.

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(\AA^2)	Q<0.9
2	GLC	C	2	11/12	0.97	0.17	88,93,107,115	0
2	GLC	C	1	12/12	0.97	0.18	87,96,113,113	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 [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
3	BTB	A	501	14/14	0.65	0.29	94,111,131,137	0
3	BTB	A	502	14/14	0.90	0.15	128,134,141,143	0
4	SO4	A	503	5/5	0.94	0.13	96,98,113,123	0

6.5 Other polymers [i](#)

There are no such residues in this entry.