



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 01:16 AM UTC

PDB ID : 4O7P / pdb_00004o7p
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 types of validation reports are described at

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

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

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

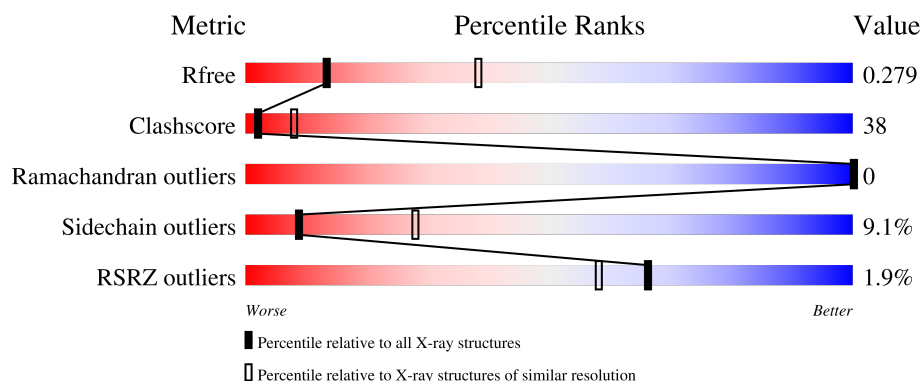
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (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 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	455	<div> <div>2%</div> <div> <div></div> <div>55%</div> <div>33%</div> <div>6%</div> <div>5%</div> </div> </div>
1	B	455	<div> <div>2%</div> <div> <div></div> <div>54%</div> <div>38%</div> <div>7%</div> <div>.</div> </div> </div>
2	C	2	<div> <div></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.

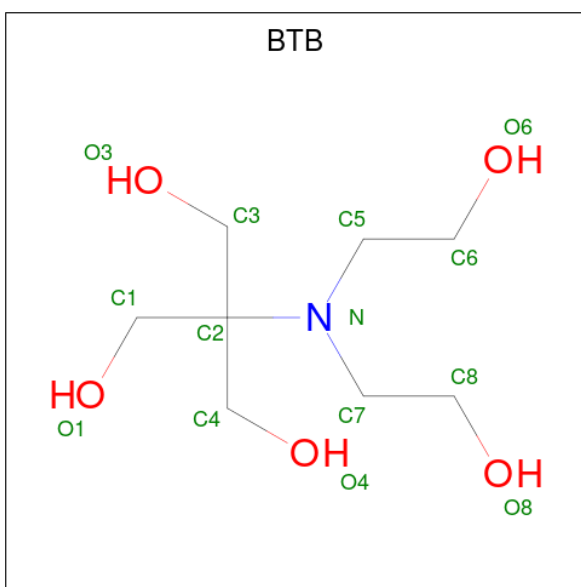
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	433	Total	C	N	O	S	0	0	0
			3349	2114	593	636	6			
1	B	451	Total	C	N	O	S	0	0	0
			3490	2199	618	667	6			

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



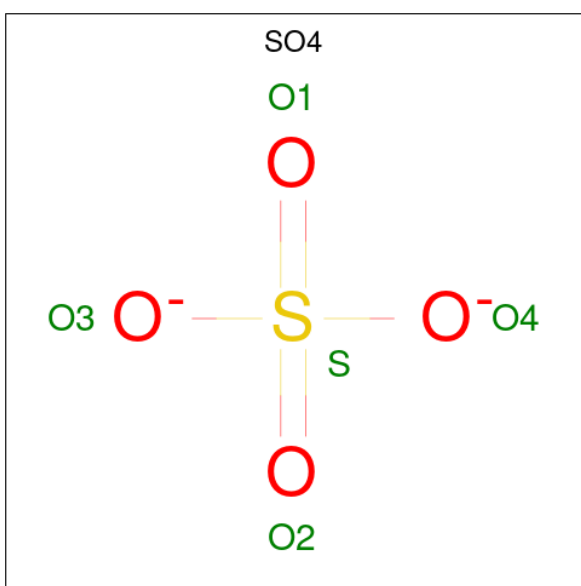
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (CCD ID: 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 (CCD ID: SO₄) (formula: O₄S).

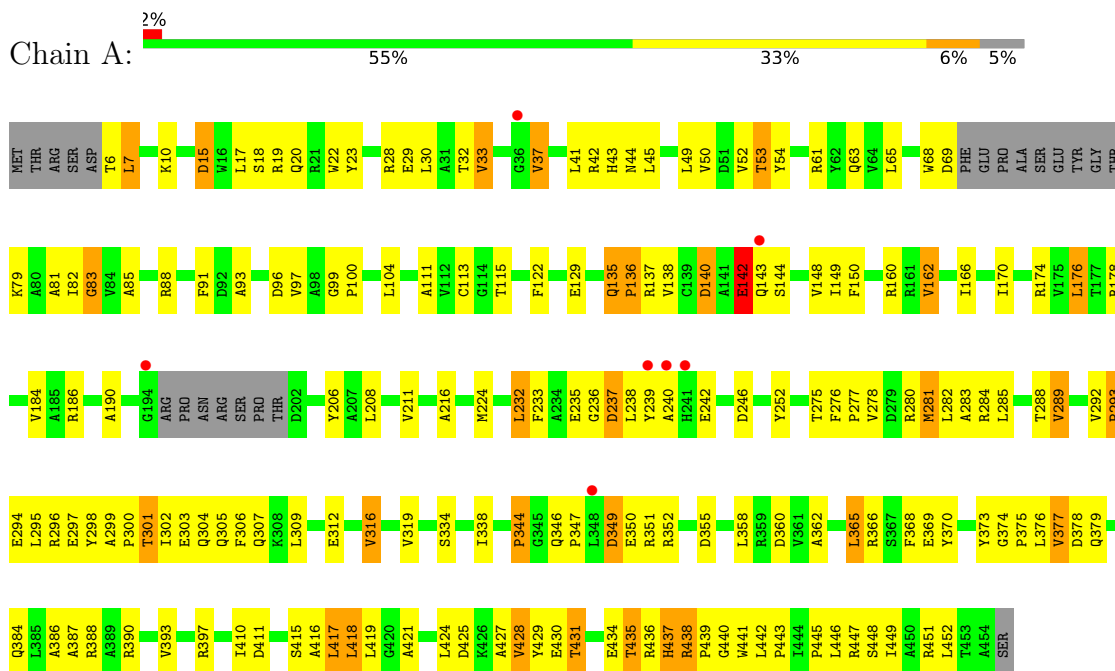


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

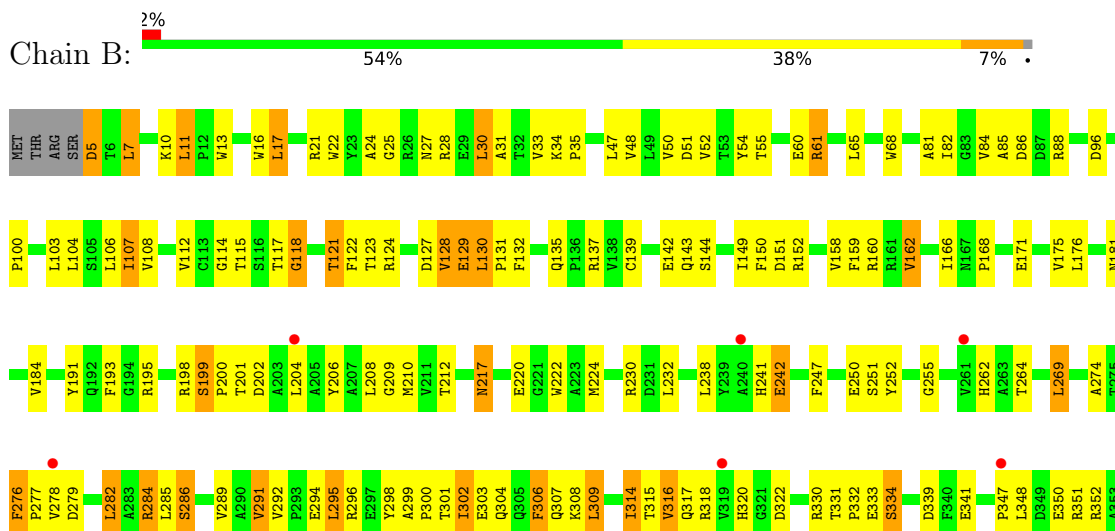
3 Residue-property plots

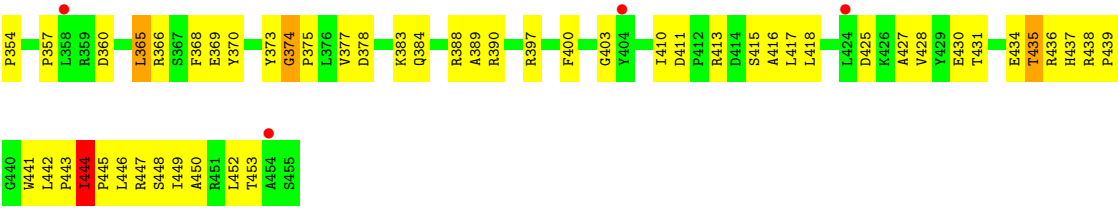
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Maltokinase



• Molecule 1: Maltokinase





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

Chain C: 100%

GLC1
GLC2

4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	96.73Å 96.73Å 461.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.60 – 2.90 44.60 – 2.90	Depositor EDS
% Data completeness (in resolution range)	97.2 (44.60-2.90) 99.6 (44.60-2.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 2.91Å)	Xtriage
Refinement program	PHENIX 1.7_650	Depositor
R, R_{free}	0.239 , 0.281 0.235 , 0.279	Depositor DCC
R_{free} test set	1498 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)	0.33 , 92.2	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	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, BTB, SO4

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	0.83	1/3420 (0.0%)	1.18	14/4661 (0.3%)
1	B	0.84	4/3568 (0.1%)	1.18	25/4867 (0.5%)
All	All	0.83	5/6988 (0.1%)	1.18	39/9528 (0.4%)

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

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	444	ILE	CA-CB	6.23	1.62	1.54
1	B	107	ILE	CA-CB	-5.38	1.48	1.54
1	B	81	ALA	CA-CB	-5.36	1.46	1.53
1	A	186	ARG	CD-NE	5.33	1.53	1.46
1	B	158	VAL	CA-CB	-5.11	1.47	1.53

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	138	VAL	N-CA-C	-9.02	95.00	108.17
1	B	118	GLY	N-CA-C	7.97	122.79	110.91
1	B	274	ALA	N-CA-C	7.46	119.03	108.74
1	A	334	SER	N-CA-C	7.03	118.86	108.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	314	ILE	N-CA-C	6.80	119.33	108.85
1	A	431	THR	N-CA-C	-6.67	104.04	111.71
1	B	199	SER	CA-C-N	6.52	128.24	121.65
1	B	199	SER	C-N-CA	6.52	128.24	121.65
1	B	411	ASP	CA-C-N	6.50	126.73	119.32
1	B	411	ASP	C-N-CA	6.50	126.73	119.32
1	B	294	GLU	N-CA-C	6.48	118.00	111.07
1	A	438	ARG	CA-C-N	6.19	125.92	119.05
1	A	438	ARG	C-N-CA	6.19	125.92	119.05
1	A	437	HIS	N-CA-C	5.86	120.58	113.38
1	B	302	ILE	N-CA-C	-5.83	104.24	110.36
1	B	181	ASN	CA-C-N	5.77	125.87	119.87
1	B	181	ASN	C-N-CA	5.77	125.87	119.87
1	A	135	GLN	N-CA-C	5.58	117.51	109.48
1	B	390	ARG	N-CA-C	-5.54	105.24	111.28
1	B	114	GLY	N-CA-C	5.54	117.47	110.38
1	B	334	SER	N-CA-C	5.43	117.45	109.24
1	B	217	ASN	N-CA-C	5.38	119.47	113.02
1	B	276	PHE	CA-C-N	5.34	129.12	121.91
1	B	276	PHE	C-N-CA	5.34	129.12	121.91
1	B	374	GLY	CA-C-N	5.23	124.84	119.56
1	B	374	GLY	C-N-CA	5.23	124.84	119.56
1	B	264	THR	CA-C-N	-5.20	113.31	120.28
1	B	264	THR	C-N-CA	-5.20	113.31	120.28
1	A	135	GLN	CA-C-O	-5.19	114.87	119.75
1	B	431	THR	N-CA-C	-5.18	105.52	111.07
1	A	83	GLY	CA-C-N	-5.17	115.92	123.06
1	A	83	GLY	C-N-CA	-5.17	115.92	123.06
1	A	136	PRO	N-CA-C	5.15	120.54	113.84
1	B	410	ILE	N-CA-C	5.15	115.53	108.12
1	A	216	ALA	N-CA-C	5.09	117.57	111.71
1	B	383	LYS	N-CA-C	-5.09	105.86	111.71
1	A	411	ASP	CA-C-N	5.05	124.84	119.28
1	A	411	ASP	C-N-CA	5.05	124.84	119.28
1	B	435	THR	N-CA-C	-5.01	105.81	111.28

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

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Mol	Chain	Res	Type	Group
1	A	344	PRO	Peptide
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	252	0
1	B	3490	0	3409	264	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	516	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 38.

All (516) 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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:18:SER:HA	1:A:23:TYR:CD2	1.99	0.98
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:A:69:ASP:C	1:A:88:ARG:HG2	1.90	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: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:30:LEU:HD11	1:B:33:VAL:HG23	1.57	0.84
1:B:289:VAL:HG21	1:B:296:ARG:NH1	1.91	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:436:ARG:HD2	1:B:436:ARG:O	1.78	0.84
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:A:142:GLU:HA	1:A:143:GLN:C	2.05	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:A:232:LEU:C	1:A:232:LEU:HD13	2.05	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: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:135:GLN:CD	1:A:136:PRO:HD3	2.10	0.77
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:7:LEU:C	1:A:7:LEU:HD23	2.13	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:97:VAL:C	1:A:100:PRO:HD2	2.12	0.73
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:HD23	1:A:7:LEU:O	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: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: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:B:307:GLN:OE1	1:B:307:GLN:HA	1.91	0.69
1:B:224:MET:HE2	1:B:247:PHE:HE1	1.57	0.69
1:B:369:GLU:HG2	1:B:448:SER:OG	1.93	0.69
1:A:17:LEU:HD23	1:A:17:LEU:C	2.17	0.69
1:A:232:LEU:HD13	1:A:232:LEU:O	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: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:106:LEU:HD13	1:B:122:PHE:HE2	1.57	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:122:PHE:HD1	1:B:191:TYR:HB2	1.58	0.68
1:A:277:PRO:HG2	1:A:281:MET:HE3	1.75	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
1:B:448:SER:O	1:B:452:LEU:HD13	1.93	0.68
1:A:7:LEU:HD22	1:A:68:TRP:CZ2	2.30	0.67
1:A:28:ARG:NH1	1:A:54:TYR:CD2	2.62	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:H	1:A:417:LEU:CD1	2.07	0.67
1:B:106:LEU:HB3	1:B:122:PHE:CE2	2.29	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:B:449:ILE:HG13	1:B:450:ALA:N	2.09	0.66
1:A:424:LEU:O	1:A:428:VAL:HG13	1.96	0.65
1:A:232:LEU:C	1:A:232:LEU:CD1	2.70	0.65
1:B:436:ARG:HG3	1:B:437:HIS:CE1	2.32	0.65
1:A:18:SER:HA	1:A:23:TYR:HD2	1.60	0.64
1:A:97:VAL:O	1:A:100:PRO:HD2	1.97	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:162:VAL:HG13	1:B:206:TYR:HD1	1.65	0.62
1:B:10:LYS:HB3	1:B:85:ALA:HB1	1.81	0.62
1:B:27:ASN:OD1	1:B:28:ARG:HD2	1.98	0.62
1:B:299:ALA:HB3	1:B:300:PRO:CD	2.30	0.62
1:B:299:ALA:C	1:B:303:GLU:OE1	2.43	0.62
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:B:47:LEU:C	1:B:47:LEU:HD23	2.26	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:82:ILE:HG23	1:A:83:GLY:H	1.65	0.60
1:A:430:GLU:O	1:A:434:GLU:HG2	2.01	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:10:LYS:CE	1:A:85:ALA:HB1	2.30	0.60
1:A:424:LEU:CD1	1:A:425:ASP:N	2.62	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:306:PHE:HE1	1:B:307:GLN:CD	2.09	0.60
1:B:444:ILE:HG13	1:B:445:PRO:HD3	1.82	0.60
1:A:417:LEU:HD12	1:A:417:LEU:H	1.62	0.60
1:B:210:MET:HE2	1:B:212:THR:HG23	1.82	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:291:VAL:CG2	1:B:435:THR:HG21	2.32	0.59
1:B:299:ALA:HB3	1:B:300:PRO:HD3	1.84	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:415:SER:C	1:A:417:LEU:HD12	2.28	0.58
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: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
1:B:159:PHE:HD2	1:B:208:LEU:C	2.11	0.58
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:28:ARG:NH1	1:A:54:TYR:CG	2.73	0.57
1:A:275:THR:HG22	1:A:276:PHE:O	2.04	0.57
1:A:37:VAL:HG13	1:A:49:LEU:HB2	1.85	0.57
1:A:276:PHE:CD2	1:A:309:LEU:HB2	2.39	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:A:111:ALA:HB2	3:A:502:BTB:H62	1.87	0.57
1:A:292:VAL:HG13	1:A:292:VAL:O	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:A:174:ARG:HD3	1:A:178:ARG:NH2	2.19	0.56
1:B:16:TRP:CD1	1:B:82:ILE:HA	2.40	0.56
1:B:47:LEU:HD23	1:B:48:VAL:N	2.20	0.56
1:A:7:LEU:CD2	1:A:7:LEU:C	2.78	0.56
1:B:65:LEU:CD1	1:B:103:LEU:CD1	2.83	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:A:277:PRO:O	1:A:281:MET:HG2	2.05	0.56
1:B:295:LEU:HD23	1:B:295:LEU:C	2.30	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:A:434:GLU:HB2	1:A:438:ARG:HB3	1.88	0.55
1:A:15:ASP:N	1:A:15:ASP:OD1	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:393:VAL:O	1:A:397:ARG:HG3	2.07	0.55
1:B:30:LEU:HD21	1:B:52:VAL:HG23	1.89	0.55
1:B:303:GLU:CG	1:B:306:PHE:CE2	2.90	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:299:ALA:C	1:B:302:ILE:HG12	2.33	0.54
1:B:303:GLU:O	1:B:306:PHE:CD1	2.61	0.54
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:N	1:B:54:TYR:HD1	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: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:B:333:GLU:HG2	1:B:334:SER:H	1.72	0.54
1:B:436:ARG:HD2	1:B:436:ARG:C	2.32	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: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:5:ASP:OD1	1:B:5:ASP:N	2.41	0.53
1:B:282:LEU:HD13	1:B:282:LEU:N	2.23	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:129:GLU:CD	1:B:129:GLU:H	2.17	0.53
1:B:282:LEU:N	1:B:282:LEU:CD1	2.71	0.53
1:B:54:TYR:N	1:B:54:TYR:CD1	2.73	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:A:43:HIS:O	1:A:44:ASN:HB2	2.09	0.53
1:A:232:LEU:HD11	1:A:388:ARG:HB3	1.91	0.53
1:A:135:GLN:CG	1:A:136:PRO:HD3	2.39	0.52
1:A:349:ASP:OD1	1:A:349:ASP:N	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201:THR:O	1:B:202:ASP:HB2	2.08	0.52
1:A:427:ALA:HA	1:A:430:GLU:OE2	2.10	0.52
1:B:441:TRP:C	1:B:443:PRO:HD2	2.36	0.51
1:B:306:PHE:CD1	1:B:306:PHE:C	2.88	0.51
1:B:370:TYR:OH	1:B:430:GLU:HB3	2.11	0.51
1:A:439:PRO:C	1:A:442:LEU:HD13	2.34	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
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:210:MET:HE2	1:B:212:THR:CG2	2.41	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:122:PHE:HE1	1:B:191:TYR:CD1	2.29	0.50
1:B:142:GLU:HG3	1:B:144:SER:H	1.75	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:276:PHE:CD2	1:B:276:PHE:C	2.90	0.50
1:A:442:LEU:N	1:A:442:LEU:CD1	2.75	0.49
1:A:93:ALA:HB1	1:A:99:GLY:HA3	1.93	0.49
1:A:297:GLU:O	1:A:300:PRO:HD2	2.12	0.49
1:A:418:LEU:HD12	1:A:418:LEU:H	1.76	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:51:ASP:OD2	1:B:61:ARG:HD2	2.13	0.49
1:A:252:TYR:CD1	1:A:252:TYR:C	2.91	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:295:LEU:CD2	1:B:295:LEU:C	2.85	0.49
1:B:84:VAL:HA	1:B:88:ARG:O	2.13	0.48
1:B:108:VAL:CG1	1:B:124:ARG:NH1	2.74	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:444:ILE:CG1	1:B:445:PRO:CD	2.89	0.48
1:A:288:THR:CG2	1:A:289:VAL:N	2.76	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:B:306:PHE:CE1	1:B:307:GLN:CD	2.90	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:118:GLY:HA3	1:B:193:PHE:CZ	2.49	0.47
1:B:162:VAL:HG13	1:B:206:TYR:CD1	2.47	0.47
1:A:30:LEU:HD22	1:A:33:VAL:HG12	1.94	0.47
1:B:52:VAL:CG2	1:B:54:TYR:CE1	2.91	0.47
1:B:299:ALA:CA	1:B:302:ILE:HG12	2.41	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:306:PHE:HD1	1:B:306:PHE:C	2.21	0.47
1:B:384:GLN:OE1	1:B:384:GLN:O	2.32	0.47
1:A:162:VAL:HG12	1:A:206:TYR:HD1	1.80	0.47
1:A:278:VAL:C	1:A:281:MET:HG3	2.40	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:OE1	1:A:242:GLU:HA	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:17:LEU:C	1:A:17:LEU:CD2	2.87	0.47
1:A:33:VAL:O	1:A:33:VAL:CG2	2.63	0.47
1:A:276:PHE:CE2	1:A:309:LEU:HB2	2.50	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:A:281:MET:HG2	1:A:281:MET:H	1.35	0.47
1:B:442:LEU:N	1:B:443:PRO:CD	2.78	0.47
1:B:151:ASP:C	1:B:152:ARG:HG2	2.40	0.47
1:B:159:PHE:CE2	1:B:209:GLY:HA2	2.50	0.46
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:175:VAL:HG21	1:B:269:LEU:HD11	1.97	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:10:LYS:HZ3	1:A:85:ALA:C	2.23	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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: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: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:198:ARG:O	1:B:199:SER:HB3	2.15	0.46
1:B:224:MET:HE2	1:B:224:MET:HB3	1.58	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
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:298:TYR:O	1:B:299:ALA:C	2.58	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:A:294:GLU:C	1:A:294:GLU:OE1	2.59	0.45
1:B:166:ILE:HD12	1:B:352:ARG:NH2	2.31	0.45
1:B:224:MET:HE2	1:B:247:PHE:CE1	2.44	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: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:A:293:PRO:O	1:A:294:GLU:C	2.60	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:CD1	1:B:269:LEU:N	2.80	0.45
1:A:137:ARG:HB3	1:A:149:ILE:HD12	1.99	0.44
1:A:427:ALA:HA	1:A:430:GLU:CD	2.42	0.44
1:A:358:LEU:HD23	1:A:419:LEU:HB2	1.99	0.44
1:B:171:GLU:CD	1:B:352:ARG:NH1	2.75	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:235:GLU:HG3	1:A:236:GLY:O	2.17	0.44
1:A:275:THR:O	1:A:276:PHE:C	2.60	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: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:B:415:SER:O	1:B:416:ALA:C	2.61	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:220:GLU:CD	1:B:222:TRP:HB3	2.43	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:A:373:TYR:HH	1:A:390:ARG:HG2	1.75	0.43
1:B:162:VAL:CG1	1:B:206:TYR:CD1	3.01	0.43
1:A:384:GLN:O	1:A:387:ALA:HB3	2.18	0.43
3:A:502:BTB:O6	3:A:502:BTB:H72	2.18	0.43
1:A:30:LEU:HD22	1:A:33:VAL:CG1	2.48	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:168:PRO:HD3	1:B:348:LEU:CD1	2.48	0.43
1:B:333:GLU:HG2	1:B:334:SER:N	2.33	0.43
1:A:434:GLU:C	1:A:442:LEU:HD11	2.44	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:A:135:GLN:CD	1:A:136:PRO:CD	2.83	0.43
1:B:279:ASP:HA	1:B:282:LEU:HD22	2.00	0.42
1:B:365:LEU:O	1:B:368:PHE:HB2	2.19	0.42
1:A:52:VAL:HG12	1:A:54:TYR:CE2	2.54	0.42
1:A:276:PHE:CD2	1:A:309:LEU:CB	3.01	0.42
1:B:341:GLU:O	1:B:351:ARG:HD2	2.19	0.42
1:A:224:MET:HE2	1:A:224:MET:HB3	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:B:303:GLU:O	1:B:304:GLN:C	2.61	0.42
1:A:373:TYR:HD1	1:A:376:LEU:HD12	1.85	0.42
1:B:255:GLY:HA2	1:B:400:PHE:CE1	2.54	0.42
1:B:285:LEU:HD11	1:B:302:ILE:HD11	2.01	0.42
1:A:69:ASP:C	1:A:88:ARG:CG	2.76	0.42
1:B:30:LEU:CD2	1:B:52:VAL:HG23	2.49	0.42
1:B:306:PHE:HD1	1:B:307:GLN:CA	2.30	0.42
1:A:22:TRP:CG	1:A:160:ARG:HG3	2.54	0.42
3:A:502:BTB:H52	3:A:502:BTB:H31	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: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:B:417:LEU:HD12	1:B:417:LEU:O	2.20	0.41
1:A:29:GLU:CD	1:A:30:LEU:H	2.28	0.41
1:A:17:LEU:CD2	1:A:23:TYR:CD1	3.04	0.41
1:A:142:GLU:CA	1:A:143:GLN:C	2.86	0.41
1:A:442:LEU:O	1:A:446:LEU:CD1	2.51	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: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:377:VAL:O	1:B:378:ASP:C	2.64	0.41
1:A:316:VAL:HG23	1:A:355:ASP:C	2.45	0.41
1:B:28:ARG:HB2	1:B:54:TYR:CD2	2.56	0.41
1:B:370:TYR:CD1	1:B:444:ILE:CD1	3.04	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:HD13	1:A:302:ILE:HG21	1.84	0.41
1:B:278:VAL:C	1:B:282:LEU:HD22	2.46	0.41
1:B:284:ARG:HH11	1:B:425:ASP:CG	2.29	0.41
1:B:7:LEU:HD12	1:B:7:LEU:HA	1.93	0.41
1:A:374:GLY:N	1:A:375:PRO:CD	2.83	0.41
1:A:415:SER:O	1:A:416:ALA:C	2.64	0.41
1:A:417:LEU:CD1	1:A:418:LEU:CD1	2.91	0.41
1:B:210:MET:HE3	1:B:210:MET:HB2	1.90	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:442:LEU:CD1	1:A:442:LEU:H	2.34	0.40
1:A:208:LEU:HA	1:A:208:LEU:HD23	1.83	0.40
1:A:418:LEU:CD1	1:A:418:LEU:H	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:N	1:B:373:TYR:CD1	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 [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	342/362 (94%)	311 (91%)	31 (9%)	9	28

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	358/362 (99%)	325 (91%)	33 (9%)	8	27
All	All	700/724 (97%)	636 (91%)	64 (9%)	9	28

All (64) 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	129	GLU
1	A	140	ASP
1	A	142	GLU
1	A	148	VAL
1	A	162	VAL
1	A	176	LEU
1	A	232	LEU
1	A	237	ASP
1	A	246	ASP
1	A	281	MET
1	A	289	VAL
1	A	301	THR
1	A	316	VAL
1	A	349	ASP
1	A	365	LEU
1	A	369	GLU
1	A	377	VAL
1	A	410	ILE
1	A	417	LEU
1	A	418	LEU
1	A	428	VAL
1	A	435	THR
1	A	448	SER
1	B	5	ASP
1	B	7	LEU
1	B	11	LEU
1	B	17	LEU
1	B	30	LEU

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Mol	Chain	Res	Type
1	B	55	THR
1	B	61	ARG
1	B	104	LEU
1	B	107	ILE
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	130	LEU
1	B	135	GLN
1	B	143	GLN
1	B	162	VAL
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

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	173	ASN
1	A	273	GLN
1	A	346	GLN
1	A	437	HIS
1	B	63	GLN
1	B	437	HIS

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	2.07	4 (33%)	17,17,17	1.27	2 (11%)
2	GLC	C	2	2	11,11,12	2.51	3 (27%)	15,15,17	1.32	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	-6.56	1.42	1.52
2	C	1	GLC	C3-C2	-4.21	1.41	1.52
2	C	1	GLC	O5-C1	3.37	1.51	1.42
2	C	2	GLC	C4-C3	-3.13	1.44	1.52
2	C	2	GLC	O5-C1	-2.75	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	-2.16	1.37	1.43

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	GLC	O3-C3-C4	-3.03	103.25	110.38
2	C	2	GLC	O2-C2-C3	-2.91	104.13	110.15
2	C	2	GLC	O2-C2-C1	2.77	115.57	109.22
2	C	1	GLC	O5-C5-C4	2.18	113.62	109.70

There are no chirality outliers.

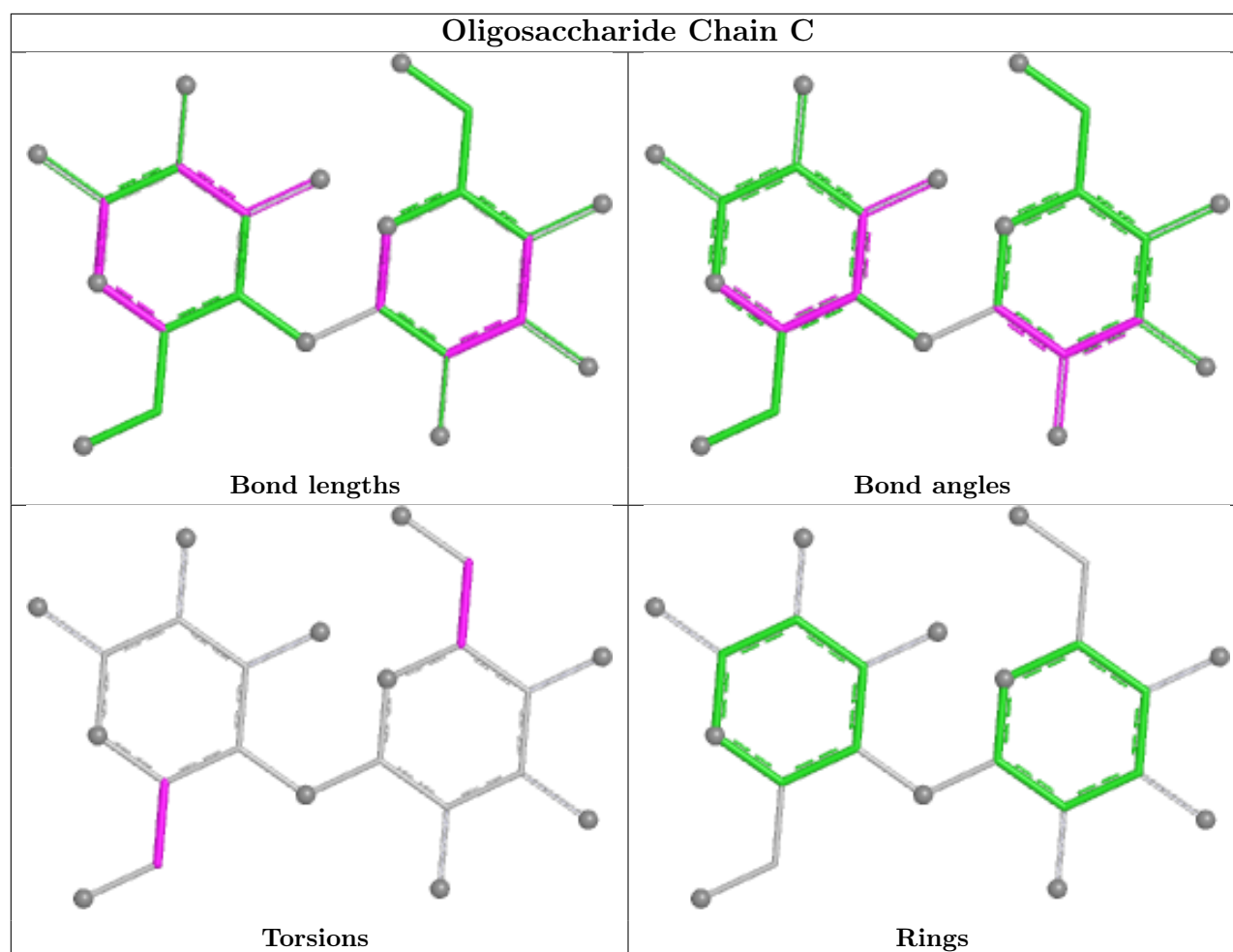
All (4) torsion outliers are listed below:

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	1.60	4 (30%)	7,16,16	2.50	2 (28%)
4	SO4	A	503	-	4,4,4	0.32	0	6,6,6	0.31	0
3	BTB	A	501	-	13,13,13	0.61	0	7,16,16	3.48	5 (71%)

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	-2.72	1.50	1.53
3	A	502	BTB	C1-C2	-2.59	1.50	1.53
3	A	502	BTB	C4-C2	-2.49	1.50	1.53
3	A	502	BTB	C5-N	-2.10	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	5.13	123.45	111.40
3	A	501	BTB	O4-C4-C2	5.01	123.17	111.40
3	A	501	BTB	O1-C1-C2	4.00	120.82	111.40
3	A	502	BTB	O4-C4-C2	3.85	120.46	111.40
3	A	502	BTB	O1-C1-C2	3.81	120.36	111.40
3	A	501	BTB	O6-C6-C5	2.31	120.75	111.22
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	501	BTB	N-C2-C3-O3
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	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	502	BTB	N-C7-C8-O8

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Mol	Chain	Res	Type	Atoms
3	A	501	BTB	C1-C2-C3-O3
3	A	501	BTB	N-C5-C6-O6
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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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%)	0.11	7 (1%) 70 62	68, 113, 168, 196	0
1	B	451/455 (99%)	0.14	10 (2%) 62 53	60, 107, 177, 205	0
All	All	884/910 (97%)	0.13	17 (1%) 66 58	60, 110, 174, 205	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	319	VAL	3.5
1	A	241	HIS	3.1
1	B	347	PRO	3.0
1	A	143	GLN	2.8
1	B	240	ALA	2.7
1	A	194	GLY	2.5
1	B	358	LEU	2.5
1	B	278	VAL	2.4
1	A	239	TYR	2.4
1	B	261	VAL	2.3
1	A	348	LEU	2.3
1	A	240	ALA	2.3
1	B	404	TYR	2.2
1	A	36	GLY	2.2
1	B	204	LEU	2.2
1	B	454	ALA	2.2
1	B	424	LEU	2.1

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

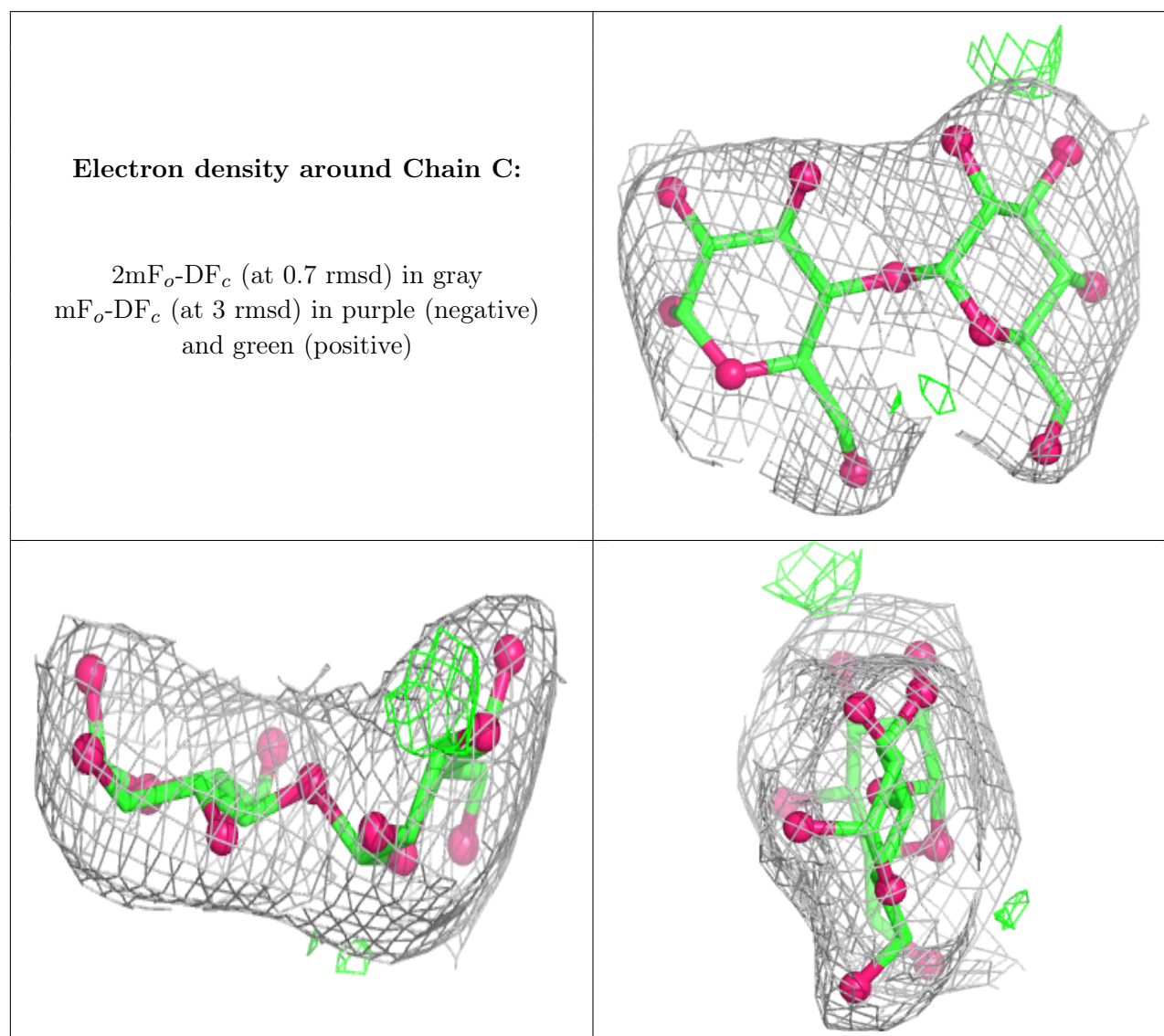
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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	C	2	11/12	0.94	0.07	88,93,107,115	0
2	GLC	C	1	12/12	0.95	0.06	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	502	14/14	0.76	0.11	128,134,141,143	0
3	BTB	A	501	14/14	0.83	0.18	94,111,131,137	0
4	SO4	A	503	5/5	0.92	0.08	96,98,113,123	0

6.5 Other polymers [i](#)

There are no such residues in this entry.