



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:00 PM UTC

PDB ID : 1HLG / pdb_00001hlg
Title : CRYSTAL STRUCTURE OF HUMAN GASTRIC LIPASE
Authors : Roussel, A.; Canaan, S.; Verger, R.; Cambillau, C.
Deposited on : 1999-03-12
Resolution : 3.00 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

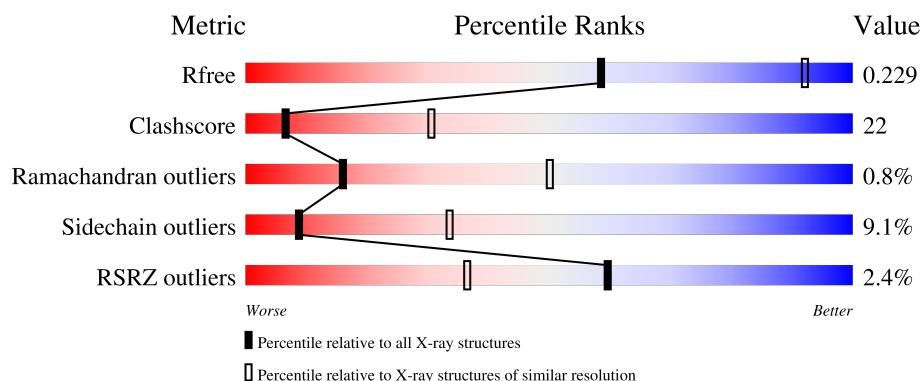
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	371	2% 53% 40% 6% ..
1	B	371	2% 54% 37% 8% .
2	C	2	50% 50%
2	D	2	50% 50%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6192 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 LIPASE, GASTRIC.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	368	Total	C	N	O	S	0	0	0
			2980	1936	489	543	12			
1	B	368	Total	C	N	O	S	0	0	0
			2980	1936	489	543	12			

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	D	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

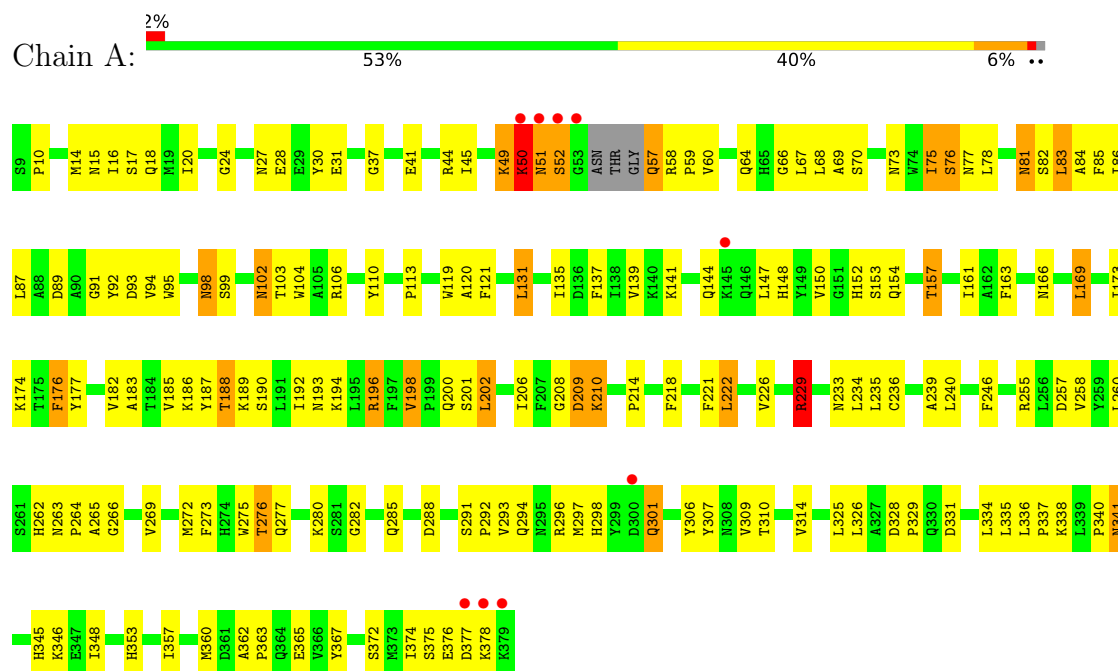
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	46	Total	O	0	0
			46	46		
4	B	46	Total	O	0	0
			46	46		

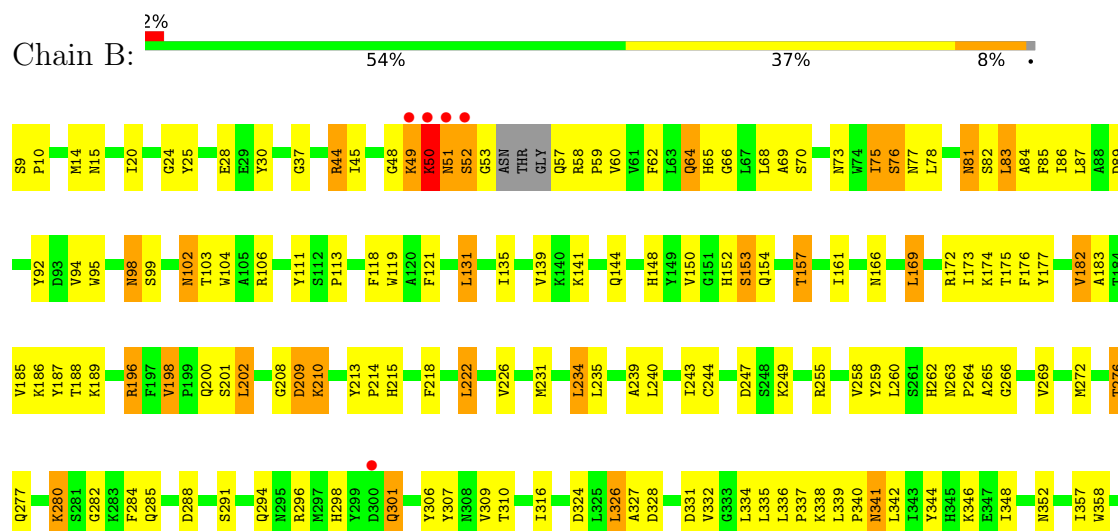
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: LIPASE, GASTRIC



• Molecule 1: LIPASE, GASTRIC





● Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



● Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	I 21 3	Depositor
Cell constants a, b, c, α , β , γ	244.33Å 244.33Å 244.33Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 3.00 15.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.4 (15.00-3.00) 98.6 (15.00-3.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.40 (at 3.01Å)	Xtriage
Refinement program	X-PLOR 3843	Depositor
R, R_{free}	0.210 , 0.228 0.211 , 0.229	Depositor DCC
R_{free} test set	2381 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	58.7	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 53.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.023 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6192	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.68% 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: NAG

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	0.93	13/3071 (0.4%)	1.15	17/4179 (0.4%)
1	B	1.01	17/3071 (0.6%)	1.18	16/4179 (0.4%)
All	All	0.97	30/6142 (0.5%)	1.16	33/8358 (0.4%)

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	378	LYS	C-O	11.53	1.37	1.24
1	B	50	LYS	C-O	11.36	1.38	1.24
1	B	377	ASP	CG-OD1	8.68	1.41	1.25
1	B	377	ASP	CG-OD2	8.45	1.41	1.25
1	A	229	ARG	CZ-NH1	8.30	1.44	1.32
1	B	378	LYS	CA-C	8.18	1.63	1.52
1	A	120	ALA	CA-CB	-7.84	1.43	1.52
1	A	229	ARG	CZ-NH2	7.67	1.43	1.33
1	B	50	LYS	CG-CD	7.59	1.75	1.52
1	A	52	SER	CA-C	7.52	1.62	1.53
1	A	229	ARG	NE-CZ	7.22	1.41	1.33
1	B	51	ASN	CB-CG	7.06	1.69	1.52
1	A	51	ASN	CA-C	7.04	1.62	1.52
1	B	51	ASN	CG-OD1	6.76	1.36	1.23
1	A	51	ASN	CG-OD1	6.72	1.36	1.23
1	B	52	SER	CA-C	6.51	1.61	1.53
1	B	231	MET	CG-SD	6.33	1.96	1.80
1	B	50	LYS	CB-CG	6.28	1.71	1.52
1	B	51	ASN	C-O	6.28	1.31	1.23
1	B	231	MET	SD-CE	6.17	1.95	1.79
1	B	50	LYS	CE-NZ	6.01	1.67	1.49
1	B	50	LYS	CD-CE	6.01	1.70	1.52
1	A	360	MET	SD-CE	5.86	1.94	1.79
1	A	52	SER	CB-OG	5.59	1.53	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	50	LYS	C-N	5.48	1.41	1.33
1	A	51	ASN	C-O	5.27	1.30	1.24
1	B	52	SER	C-O	5.16	1.30	1.23
1	A	297	MET	SD-CE	5.10	1.92	1.79
1	A	52	SER	N-CA	5.10	1.53	1.46
1	A	51	ASN	CG-ND2	5.07	1.44	1.33

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	198	VAL	N-CA-C	9.88	119.37	107.61
1	A	198	VAL	N-CA-C	8.74	118.02	107.61
1	A	210	LYS	N-CA-C	8.51	120.59	108.00
1	B	210	LYS	N-CA-C	8.26	120.23	108.00
1	B	76	SER	N-CA-C	8.18	122.92	113.19
1	A	37	GLY	N-CA-C	7.92	125.21	114.92
1	A	76	SER	N-CA-C	7.88	122.57	113.19
1	B	37	GLY	N-CA-C	7.73	124.97	114.92
1	B	45	ILE	CA-C-N	7.15	127.14	119.28
1	B	45	ILE	C-N-CA	7.15	127.14	119.28
1	B	183	ALA	N-CA-C	-6.91	104.84	113.55
1	A	45	ILE	CA-C-N	6.64	126.33	119.56
1	A	45	ILE	C-N-CA	6.64	126.33	119.56
1	A	209	ASP	N-CA-C	-6.45	100.02	109.63
1	B	301	GLN	N-CA-C	-6.23	100.07	108.86
1	B	52	SER	N-CA-C	-6.13	102.79	111.24
1	A	183	ALA	N-CA-C	-6.05	104.88	112.93
1	B	153	SER	CB-CA-C	-6.04	109.59	116.54
1	A	110	TYR	N-CA-C	6.04	121.57	113.72
1	A	301	GLN	N-CA-C	-6.02	100.63	109.18
1	A	353	HIS	N-CA-C	5.97	118.27	111.11
1	A	218	PHE	N-CA-C	5.97	120.56	113.28
1	B	189	LYS	N-CA-C	-5.89	105.40	112.59
1	A	94	VAL	N-CA-C	5.85	117.17	108.23
1	B	50	LYS	N-CA-C	-5.81	98.43	110.80
1	B	209	ASP	N-CA-C	-5.71	101.12	109.63
1	B	218	PHE	N-CA-C	5.61	120.12	113.28
1	B	94	VAL	N-CA-C	5.49	116.63	108.23
1	A	210	LYS	CB-CA-C	-5.33	108.88	116.34
1	A	51	ASN	CA-C-N	-5.29	115.01	122.63
1	A	51	ASN	C-N-CA	-5.29	115.01	122.63
1	A	52	SER	N-CA-C	-5.08	104.42	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	182	VAL	CB-CA-C	-5.00	103.09	111.29

There are no chirality outliers.

There are no planarity outliers.

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	2980	0	2879	132	0
1	B	2980	0	2879	124	0
2	C	28	0	25	3	0
2	D	28	0	25	0	0
3	A	42	0	39	0	0
3	B	42	0	39	1	0
4	A	46	0	0	0	0
4	B	46	0	0	1	0
All	All	6192	0	5886	260	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 22.

All (260) 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:50:LYS:CD	1:B:50:LYS:CG	1.75	1.56
1:A:69:ALA:HB1	1:A:73:ASN:HD21	1.18	1.06
1:B:69:ALA:HB1	1:B:73:ASN:HD21	1.35	0.92
1:A:291:SER:OG	1:A:294:GLN:HG3	1.71	0.91
1:A:229:ARG:HA	1:A:229:ARG:HE	1.35	0.90
1:A:73:ASN:HB3	1:A:357:ILE:HG21	1.55	0.87
1:A:78:LEU:H	1:A:81:ASN:HD21	1.27	0.81
1:A:66:GLY:HA3	1:A:153:SER:HB3	1.61	0.80
1:B:78:LEU:H	1:B:81:ASN:HD21	1.29	0.79
1:B:73:ASN:HB3	1:B:357:ILE:HG21	1.64	0.79
1:A:69:ALA:HB1	1:A:73:ASN:ND2	1.97	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:66:GLY:HA3	1:B:153:SER:HB3	1.62	0.79
1:B:336:LEU:HB2	1:B:337:PRO:HD3	1.64	0.77
1:B:186:LYS:HD2	1:B:187:TYR:CZ	2.20	0.77
1:A:186:LYS:HB3	1:A:282:GLY:HA2	1.66	0.77
1:A:154:GLN:O	1:A:157:THR:HG23	1.85	0.76
1:B:291:SER:OG	1:B:294:GLN:HG3	1.86	0.75
1:B:363:PRO:HA	1:B:367:TYR:HB2	1.69	0.75
1:A:75:ILE:HD12	1:A:84:ALA:HB3	1.69	0.73
1:A:185:VAL:HG22	1:A:188:THR:HG21	1.69	0.72
1:B:69:ALA:HB1	1:B:73:ASN:ND2	2.04	0.72
1:B:157:THR:HB	1:B:182:VAL:O	1.89	0.72
1:B:69:ALA:HA	1:B:263:ASN:HD22	1.54	0.71
1:A:154:GLN:HA	1:A:157:THR:CG2	2.20	0.70
1:B:99:SER:O	1:B:102:ASN:ND2	2.25	0.69
1:B:154:GLN:HA	1:B:157:THR:CG2	2.22	0.69
1:A:60:VAL:H	1:A:378:LYS:HZ3	1.40	0.69
1:B:44:ARG:HG3	1:B:95:TRP:CE2	2.28	0.69
1:A:99:SER:O	1:A:102:ASN:ND2	2.26	0.69
1:B:78:LEU:N	1:B:81:ASN:HD21	1.91	0.68
1:A:272:MET:O	1:A:276:THR:HG23	1.93	0.68
1:B:59:PRO:HG2	1:B:92:TYR:CD2	2.29	0.68
1:B:60:VAL:H	1:B:378:LYS:HZ3	1.43	0.67
1:B:60:VAL:H	1:B:378:LYS:NZ	1.94	0.65
1:A:157:THR:HB	1:A:182:VAL:O	1.97	0.65
1:B:185:VAL:HG22	1:B:188:THR:HG21	1.79	0.65
1:B:81:ASN:HD22	1:B:81:ASN:C	2.03	0.64
1:B:15:ASN:HB2	1:B:104:TRP:CH2	2.32	0.64
1:B:10:PRO:O	1:B:14:MET:HE3	1.97	0.64
1:A:44:ARG:HG3	1:A:95:TRP:CE2	2.33	0.64
1:B:119:TRP:CG	1:B:210:LYS:HA	2.33	0.64
1:B:310:THR:HG22	1:B:338:LYS:HB3	1.80	0.64
1:A:363:PRO:HA	1:A:367:TYR:HB2	1.80	0.63
1:B:76:SER:O	1:B:255:ARG:NH1	2.31	0.63
1:B:14:MET:O	1:B:262:HIS:HE1	1.81	0.63
1:B:154:GLN:O	1:B:157:THR:HG23	1.98	0.63
1:A:60:VAL:H	1:A:378:LYS:NZ	1.97	0.62
1:A:78:LEU:N	1:A:81:ASN:HD21	1.97	0.62
1:B:57:GLN:HG3	1:B:58:ARG:H	1.66	0.61
1:B:154:GLN:HA	1:B:157:THR:HG23	1.82	0.61
1:A:273:PHE:O	1:A:277:GLN:HG3	2.00	0.61
1:A:229:ARG:HE	1:A:229:ARG:CA	2.05	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:174:LYS:HD3	1:A:378:LYS:HB2	1.82	0.60
1:A:258:VAL:O	1:A:262:HIS:HD2	1.84	0.60
1:A:185:VAL:HG13	1:A:185:VAL:O	2.00	0.60
1:B:340:PRO:O	1:B:341:ASN:CB	2.49	0.60
1:B:152:HIS:ND1	1:B:357:ILE:HD11	2.16	0.60
1:B:51:ASN:HB3	1:B:89:ASP:HB3	1.82	0.60
1:B:196:ARG:HA	1:B:280:LYS:HE2	1.83	0.60
1:A:59:PRO:HG2	1:A:92:TYR:CD2	2.37	0.60
1:B:102:ASN:HA	1:B:266:GLY:HA3	1.84	0.59
1:A:186:LYS:HB3	1:A:282:GLY:CA	2.32	0.59
1:B:70:SER:HB2	1:B:265:ALA:HB2	1.83	0.59
1:B:106:ARG:O	1:B:113:PRO:HG3	2.02	0.59
1:A:15:ASN:OD1	1:A:18:GLN:HG3	2.03	0.59
1:A:277:GLN:HE22	1:A:288:ASP:H	1.49	0.59
1:A:310:THR:HG22	1:A:338:LYS:O	2.02	0.58
1:A:103:THR:HG22	1:A:104:TRP:CD2	2.37	0.58
1:B:198:VAL:HG13	1:B:202:LEU:HB3	1.85	0.58
1:A:309:VAL:HG21	1:A:335:LEU:CD1	2.33	0.58
1:A:348:ILE:HD13	1:A:365:GLU:HB3	1.84	0.58
1:A:15:ASN:HB2	1:A:104:TRP:CH2	2.38	0.58
1:B:235:LEU:HD12	1:B:260:LEU:HD21	1.85	0.58
1:B:234:LEU:HD13	1:B:260:LEU:HD13	1.86	0.58
1:B:103:THR:HG22	1:B:104:TRP:CD2	2.39	0.57
1:A:214:PRO:HG3	1:A:239:ALA:HB2	1.85	0.57
1:A:14:MET:O	1:A:262:HIS:HE1	1.88	0.57
1:A:336:LEU:HD11	1:A:345:HIS:CE1	2.39	0.57
1:B:285:GLN:HB3	1:B:306:TYR:CD1	2.40	0.56
1:A:76:SER:O	1:A:255:ARG:NH1	2.38	0.56
1:A:119:TRP:CG	1:A:210:LYS:HA	2.40	0.56
1:A:340:PRO:O	1:A:341:ASN:CB	2.52	0.56
1:A:152:HIS:ND1	1:A:357:ILE:HD11	2.22	0.56
1:A:102:ASN:HA	1:A:266:GLY:HA3	1.86	0.55
1:B:51:ASN:HB3	1:B:89:ASP:CB	2.36	0.55
1:B:214:PRO:HG3	1:B:239:ALA:HB2	1.88	0.55
1:A:198:VAL:HG13	1:A:202:LEU:HB3	1.88	0.55
1:B:65:HIS:HD2	1:B:66:GLY:O	1.90	0.55
1:A:51:ASN:HB3	1:A:89:ASP:HB3	1.88	0.55
1:B:28:GLU:HB3	1:B:30:TYR:CE2	2.42	0.55
1:A:75:ILE:HG13	1:A:85:PHE:CZ	2.40	0.55
1:A:154:GLN:HA	1:A:157:THR:HG23	1.88	0.54
1:B:75:ILE:HG13	1:B:85:PHE:CZ	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:LEU:CD2	1:A:87:LEU:HG	2.37	0.54
1:B:161:ILE:HG23	1:B:307:TYR:CE1	2.42	0.54
1:B:15:ASN:HB2	1:B:104:TRP:CZ2	2.42	0.54
1:A:77:ASN:HA	1:A:255:ARG:HH11	1.73	0.53
1:A:185:VAL:HG22	1:A:188:THR:CG2	2.36	0.53
1:A:292:PRO:O	1:A:296:ARG:HG3	2.09	0.53
1:A:20:ILE:HG12	1:A:75:ILE:HG12	1.90	0.53
1:A:309:VAL:HG21	1:A:335:LEU:HD11	1.91	0.53
1:B:187:TYR:HE2	1:B:334:LEU:HD11	1.74	0.53
1:A:75:ILE:CD1	1:A:84:ALA:HB3	2.37	0.53
1:B:10:PRO:C	1:B:14:MET:HE3	2.34	0.53
1:B:139:VAL:HG13	1:B:144:GLN:O	2.08	0.53
1:A:336:LEU:CD1	1:A:345:HIS:CE1	2.92	0.53
1:A:152:HIS:CG	1:A:357:ILE:HD11	2.43	0.52
1:A:196:ARG:HA	1:A:280:LYS:HE2	1.91	0.52
1:A:57:GLN:HG3	1:A:58:ARG:H	1.74	0.52
1:B:173:ILE:N	1:B:173:ILE:HD12	2.25	0.52
1:B:328:ASP:OD1	1:B:331:ASP:OD2	2.28	0.52
1:B:186:LYS:HB3	1:B:282:GLY:HA2	1.91	0.52
1:B:272:MET:O	1:B:276:THR:HG23	2.09	0.52
1:A:68:LEU:O	1:A:264:PRO:HB3	2.10	0.52
1:A:28:GLU:HB3	1:A:30:TYR:CE1	2.44	0.52
1:B:50:LYS:HB2	1:B:50:LYS:NZ	2.25	0.52
1:A:310:THR:HG22	1:A:338:LYS:HB3	1.91	0.52
1:B:83:LEU:HD22	1:B:87:LEU:HG	1.92	0.52
1:B:258:VAL:O	1:B:262:HIS:HD2	1.93	0.52
1:B:75:ILE:HD12	1:B:84:ALA:HB3	1.90	0.52
1:A:131:LEU:HD22	1:A:135:ILE:HD12	1.92	0.52
1:A:102:ASN:HB2	1:A:265:ALA:CB	2.40	0.51
1:A:189:LYS:HB3	1:A:325:LEU:HD22	1.92	0.51
1:B:62:PHE:HE2	1:B:64:GLN:HG2	1.75	0.51
1:A:222:LEU:O	1:A:226:VAL:HG23	2.11	0.51
1:A:291:SER:HG	1:A:294:GLN:HG3	1.74	0.51
1:B:362:ALA:N	1:B:363:PRO:CD	2.74	0.51
1:A:154:GLN:C	1:A:157:THR:HG23	2.36	0.51
3:B:1:NAG:H61	4:B:388:HOH:O	2.10	0.51
1:B:309:VAL:HG21	1:B:335:LEU:CD1	2.40	0.51
1:B:277:GLN:HE22	1:B:288:ASP:H	1.58	0.51
1:B:348:ILE:HD13	1:B:365:GLU:HB3	1.94	0.51
1:A:372:SER:O	1:A:376:GLU:HG3	2.11	0.50
1:B:20:ILE:HG12	1:B:75:ILE:HG12	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:2:NAG:O7	2:C:2:NAG:H3	2.11	0.50
1:A:50:LYS:HB2	1:A:50:LYS:NZ	2.26	0.50
1:B:24:GLY:O	1:B:49:LYS:HB2	2.11	0.50
1:A:77:ASN:HA	1:A:255:ARG:NH1	2.27	0.49
1:A:81:ASN:C	1:A:81:ASN:HD22	2.20	0.49
1:B:25:TYR:CZ	1:B:49:LYS:HD2	2.47	0.49
1:A:148:HIS:CE1	1:A:374:ILE:HG23	2.48	0.49
1:B:77:ASN:HA	1:B:255:ARG:HH11	1.78	0.49
1:A:10:PRO:C	1:A:14:MET:HE3	2.37	0.49
1:A:336:LEU:HB2	1:A:337:PRO:HD3	1.95	0.49
1:B:68:LEU:O	1:B:264:PRO:HB3	2.12	0.49
1:A:69:ALA:HA	1:A:263:ASN:HD22	1.78	0.49
1:A:235:LEU:HD12	1:A:260:LEU:HD21	1.95	0.49
1:A:362:ALA:N	1:A:363:PRO:CD	2.76	0.49
1:B:57:GLN:HG3	1:B:58:ARG:N	2.28	0.49
1:B:222:LEU:O	1:B:226:VAL:HG23	2.13	0.48
1:B:372:SER:O	1:B:376:GLU:HG3	2.13	0.48
1:A:52:SER:HB3	1:A:91:GLY:HA2	1.95	0.48
1:B:121:PHE:CD1	1:B:121:PHE:C	2.90	0.48
1:B:208:GLY:O	1:B:269:VAL:HG11	2.12	0.48
1:A:44:ARG:HH11	1:A:93:ASP:CG	2.20	0.48
1:A:51:ASN:O	1:A:52:SER:HB2	2.13	0.47
1:A:293:VAL:HA	1:A:296:ARG:NH1	2.29	0.47
1:A:348:ILE:CD1	1:A:365:GLU:HB3	2.44	0.47
1:A:150:VAL:HG22	1:A:177:TYR:HB2	1.97	0.47
1:B:81:ASN:C	1:B:81:ASN:ND2	2.72	0.47
1:B:152:HIS:ND1	1:B:357:ILE:CD1	2.77	0.47
1:B:174:LYS:HD3	1:B:378:LYS:HB2	1.96	0.47
1:A:49:LYS:O	1:A:50:LYS:HB2	2.14	0.47
1:A:70:SER:HB2	1:A:265:ALA:HB2	1.96	0.47
1:B:259:TYR:HB3	1:B:358:TRP:CH2	2.50	0.47
1:B:119:TRP:CB	1:B:210:LYS:HA	2.44	0.47
1:B:131:LEU:HD22	1:B:135:ILE:HD12	1.97	0.47
1:A:296:ARG:HA	1:A:301:GLN:O	2.16	0.46
1:B:186:LYS:HD2	1:B:187:TYR:CE2	2.50	0.46
1:A:187:TYR:HE2	1:A:334:LEU:HD11	1.80	0.46
1:A:119:TRP:CB	1:A:210:LYS:HA	2.45	0.46
1:B:166:ASN:CG	1:B:169:LEU:HB2	2.41	0.46
1:A:82:SER:O	1:A:83:LEU:C	2.58	0.46
1:B:185:VAL:CG2	1:B:188:THR:HG21	2.45	0.46
1:B:247:ASP:OD1	1:B:249:LYS:HG2	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:ASP:OD1	1:A:331:ASP:OD2	2.34	0.46
1:B:51:ASN:CB	1:B:89:ASP:HB3	2.46	0.46
1:B:150:VAL:HG22	1:B:177:TYR:HB2	1.97	0.46
1:A:20:ILE:HD12	1:A:27:ASN:ND2	2.30	0.46
1:A:83:LEU:HD22	1:A:87:LEU:HG	1.98	0.46
1:A:106:ARG:O	1:A:113:PRO:HG3	2.16	0.46
1:A:310:THR:HA	1:A:340:PRO:HD3	1.98	0.46
1:B:77:ASN:HA	1:B:255:ARG:NH1	2.31	0.46
1:B:185:VAL:HG22	1:B:188:THR:CG2	2.46	0.46
1:B:213:TYR:O	1:B:215:HIS:HD2	1.98	0.46
1:A:173:ILE:HD12	1:A:173:ILE:N	2.31	0.45
1:A:202:LEU:HD22	1:A:206:ILE:CD1	2.47	0.45
1:B:59:PRO:HG2	1:B:92:TYR:HD2	1.77	0.45
1:A:139:VAL:HG13	1:A:144:GLN:O	2.16	0.45
1:B:369:ASP:O	1:B:373:MET:HB2	2.16	0.45
1:B:102:ASN:HB2	1:B:265:ALA:CB	2.47	0.45
1:B:352:ASN:OD1	1:B:352:ASN:C	2.58	0.45
1:B:75:ILE:CD1	1:B:84:ALA:HB3	2.46	0.45
1:A:374:ILE:C	1:A:376:GLU:H	2.26	0.44
1:B:296:ARG:HA	1:B:301:GLN:O	2.17	0.44
1:A:257:ASP:N	1:A:257:ASP:OD1	2.49	0.44
2:C:1:NAG:O7	2:C:1:NAG:H3	2.18	0.44
1:A:194:LYS:HE3	1:A:221:PHE:CD1	2.53	0.44
1:A:161:ILE:HG23	1:A:307:TYR:CE1	2.52	0.44
1:A:189:LYS:CB	1:A:325:LEU:HD22	2.48	0.44
1:B:82:SER:O	1:B:83:LEU:C	2.60	0.44
1:B:346:LYS:NZ	1:B:365:GLU:O	2.45	0.44
1:A:16:ILE:HG23	1:A:17:SER:N	2.33	0.44
1:A:51:ASN:HB3	1:A:89:ASP:CB	2.48	0.44
1:A:208:GLY:O	1:A:269:VAL:HG11	2.17	0.44
1:A:280:LYS:HE3	1:A:280:LYS:HB2	1.80	0.44
1:B:185:VAL:HG13	1:B:185:VAL:O	2.18	0.44
1:A:137:PHE:C	1:A:137:PHE:CD1	2.95	0.43
1:B:294:GLN:O	1:B:298:HIS:HD2	2.01	0.43
1:B:243:ILE:HG22	1:B:243:ILE:O	2.18	0.43
1:B:284:PHE:CD2	1:B:307:TYR:HB2	2.53	0.43
1:A:67:LEU:HD13	1:A:275:TRP:NE1	2.33	0.43
1:A:163:PHE:CZ	1:A:173:ILE:HG12	2.53	0.43
1:B:50:LYS:O	1:B:53:GLY:N	2.51	0.43
1:B:296:ARG:C	1:B:298:HIS:H	2.26	0.43
1:B:51:ASN:O	1:B:52:SER:C	2.61	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:175:THR:HB	1:B:177:TYR:CE1	2.53	0.43
1:B:342:LEU:HD11	1:B:344:TYR:O	2.18	0.43
1:A:285:GLN:HB3	1:A:306:TYR:CD2	2.54	0.43
1:B:98:ASN:HD22	1:B:98:ASN:HA	1.59	0.43
1:B:234:LEU:HD13	1:B:260:LEU:CD1	2.49	0.43
1:B:374:ILE:C	1:B:376:GLU:H	2.27	0.43
1:A:147:LEU:C	1:A:147:LEU:HD12	2.44	0.43
1:A:340:PRO:O	1:A:340:PRO:HG2	2.19	0.43
1:A:44:ARG:HD3	1:A:93:ASP:OD1	2.19	0.43
1:A:176:PHE:CD1	1:A:176:PHE:C	2.96	0.43
1:B:174:LYS:HZ2	1:B:378:LYS:CG	2.32	0.43
1:A:166:ASN:CG	1:A:169:LEU:HB2	2.44	0.42
1:A:173:ILE:CG2	1:A:314:VAL:HG11	2.49	0.42
1:B:310:THR:HA	1:B:340:PRO:HD3	2.01	0.42
1:A:154:GLN:CA	1:A:157:THR:HG23	2.49	0.42
1:B:48:GLY:HA3	1:B:51:ASN:O	2.18	0.42
1:B:148:HIS:NE2	1:B:378:LYS:HB2	2.35	0.42
1:B:106:ARG:HA	1:B:118:PHE:HZ	1.85	0.42
1:A:24:GLY:O	1:A:49:LYS:HB2	2.20	0.42
1:B:174:LYS:NZ	1:B:378:LYS:HG3	2.35	0.42
1:B:69:ALA:CB	1:B:73:ASN:HD21	2.18	0.42
1:A:190:SER:C	1:A:192:ILE:N	2.78	0.42
1:B:174:LYS:HZ2	1:B:378:LYS:HD2	1.85	0.41
1:A:296:ARG:C	1:A:298:HIS:H	2.28	0.41
1:A:60:VAL:N	1:A:378:LYS:HZ3	2.14	0.41
1:A:121:PHE:CD1	1:A:121:PHE:C	2.98	0.41
1:B:324:ASP:OD2	1:B:327:ALA:HB3	2.20	0.41
1:B:340:PRO:O	1:B:341:ASN:HB3	2.19	0.41
1:A:196:ARG:HE	1:A:196:ARG:HB3	1.76	0.41
1:B:316:ILE:HG21	1:B:339:LEU:HD13	2.02	0.41
1:A:345:HIS:CD2	1:A:346:LYS:N	2.88	0.41
1:A:98:ASN:HD22	1:A:98:ASN:HA	1.62	0.41
1:A:152:HIS:ND1	1:A:357:ILE:CD1	2.83	0.41
2:C:1:NAG:H61	2:C:2:NAG:C1	2.51	0.41
1:A:340:PRO:O	1:A:341:ASN:HB3	2.21	0.41
1:B:244:CYS:SG	1:B:326:LEU:HD12	2.61	0.41
1:A:236:CYS:HA	1:A:246:PHE:CE1	2.56	0.41
1:B:328:ASP:O	1:B:332:VAL:HG23	2.21	0.40
1:A:193:ASN:O	1:A:196:ARG:HB3	2.21	0.40
1:A:258:VAL:O	1:A:262:HIS:CD2	2.69	0.40
1:A:15:ASN:HB2	1:A:104:TRP:CZ2	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:ASN:ND2	1:B:102:ASN:H	2.20	0.40
1:A:31:GLU:HG3	1:A:41:GLU:OE2	2.22	0.40
1:A:174:LYS:HE2	1:A:377:ASP:HB3	2.04	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	364/371 (98%)	333 (92%)	28 (8%)	3 (1%)	16	50
1	B	364/371 (98%)	330 (91%)	31 (8%)	3 (1%)	16	50
All	All	728/742 (98%)	663 (91%)	59 (8%)	6 (1%)	16	50

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	50	LYS
1	B	50	LYS
1	B	375	SER
1	A	341	ASN
1	A	375	SER
1	B	341	ASN

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	324/326 (99%)	295 (91%)	29 (9%)	9	34
1	B	324/326 (99%)	294 (91%)	30 (9%)	8	32
All	All	648/652 (99%)	589 (91%)	59 (9%)	9	33

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	LYS
1	A	50	LYS
1	A	57	GLN
1	A	64	GLN
1	A	75	ILE
1	A	81	ASN
1	A	83	LEU
1	A	86	ILE
1	A	98	ASN
1	A	102	ASN
1	A	131	LEU
1	A	141	LYS
1	A	157	THR
1	A	169	LEU
1	A	176	PHE
1	A	188	THR
1	A	196	ARG
1	A	200	GLN
1	A	201	SER
1	A	202	LEU
1	A	209	ASP
1	A	222	LEU
1	A	229	ARG
1	A	233	ASN
1	A	234	LEU
1	A	240	LEU
1	A	276	THR
1	A	326	LEU
1	A	329	PRO
1	B	9	SER
1	B	44	ARG
1	B	49	LYS
1	B	50	LYS
1	B	64	GLN
1	B	75	ILE

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Mol	Chain	Res	Type
1	B	81	ASN
1	B	83	LEU
1	B	86	ILE
1	B	98	ASN
1	B	102	ASN
1	B	111	TYR
1	B	131	LEU
1	B	141	LYS
1	B	157	THR
1	B	169	LEU
1	B	172	ARG
1	B	176	PHE
1	B	196	ARG
1	B	200	GLN
1	B	201	SER
1	B	202	LEU
1	B	209	ASP
1	B	222	LEU
1	B	234	LEU
1	B	240	LEU
1	B	276	THR
1	B	280	LYS
1	B	326	LEU
1	B	364	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	27	ASN
1	A	43	ASN
1	A	57	GLN
1	A	65	HIS
1	A	73	ASN
1	A	81	ASN
1	A	98	ASN
1	A	102	ASN
1	A	144	GLN
1	A	193	ASN
1	A	215	HIS
1	A	262	HIS
1	A	263	ASN
1	A	277	GLN

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Mol	Chain	Res	Type
1	B	27	ASN
1	B	43	ASN
1	B	65	HIS
1	B	73	ASN
1	B	81	ASN
1	B	98	ASN
1	B	102	ASN
1	B	144	GLN
1	B	193	ASN
1	B	215	HIS
1	B	262	HIS
1	B	263	ASN
1	B	277	GLN
1	B	285	GLN
1	B	298	HIS
1	B	301	GLN
1	B	313	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	NAG	C	1	1,2	14,14,15	0.70	0	17,19,21	0.97	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	C	2	2	14,14,15	0.74	0	17,19,21	0.73	0
2	NAG	D	1	1,2	14,14,15	0.74	0	17,19,21	0.98	1 (5%)
2	NAG	D	2	2	14,14,15	0.67	0	17,19,21	0.78	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
2	NAG	C	1	1,2	-	3/6/23/26	0/1/1/1
2	NAG	C	2	2	-	3/6/23/26	0/1/1/1
2	NAG	D	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	D	2	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C4-C3-C2	-3.21	106.32	111.02
2	D	1	NAG	C4-C3-C2	-2.34	107.59	111.02

There are no chirality outliers.

All (9) torsion outliers are listed below:

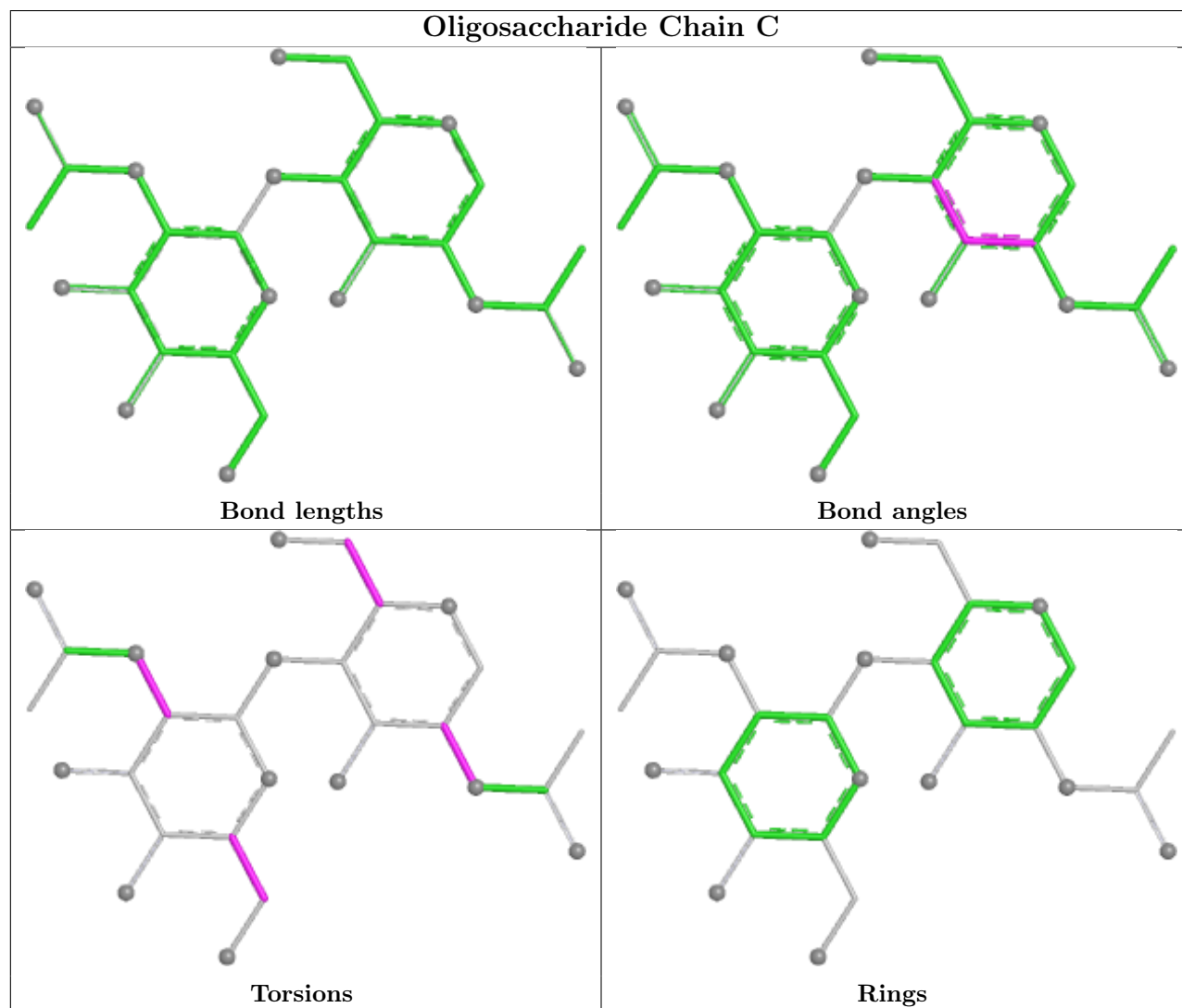
Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C3-C2-N2-C7
2	D	2	NAG	C1-C2-N2-C7
2	C	2	NAG	O5-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
2	C	2	NAG	C3-C2-N2-C7
2	D	1	NAG	C3-C2-N2-C7
2	D	2	NAG	C3-C2-N2-C7
2	C	1	NAG	C4-C5-C6-O6
2	C	1	NAG	O5-C5-C6-O6

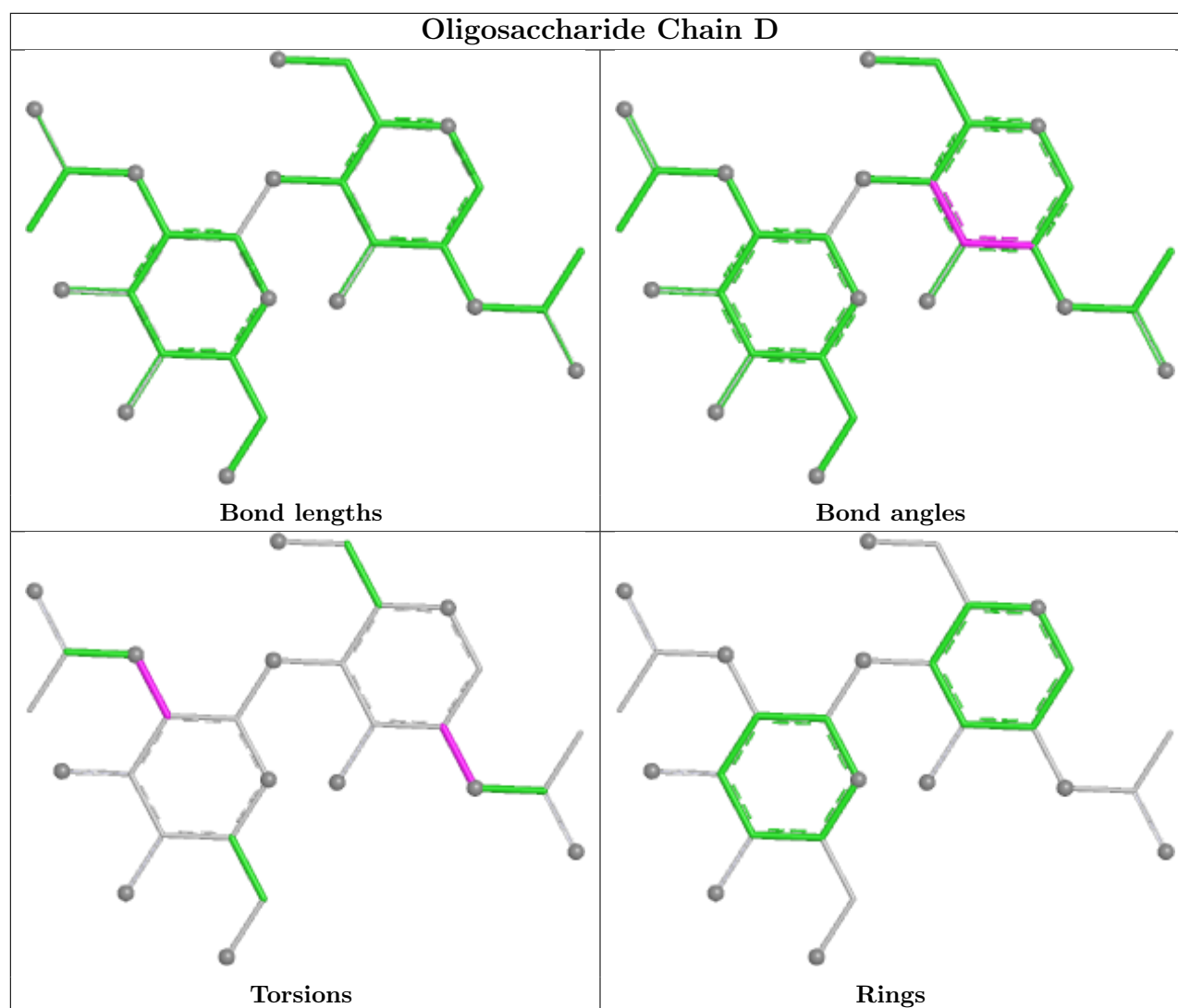
There are no ring outliers.

2 monomers are involved in 3 short contacts:

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

6 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	NAG	B	1	1	14,14,15	0.73	0	17,19,21	0.75	1 (5%)
3	NAG	B	2	1	14,14,15	1.12	2 (14%)	17,19,21	0.82	0
3	NAG	B	5	1	14,14,15	1.01	1 (7%)	17,19,21	1.24	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	A	1	1	14,14,15	0.64	0	17,19,21	0.72	0
3	NAG	A	2	1	14,14,15	0.87	1 (7%)	17,19,21	1.03	1 (5%)
3	NAG	A	5	1	14,14,15	1.15	1 (7%)	17,19,21	1.33	2 (11%)

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	NAG	B	1	1	-	0/6/23/26	0/1/1/1
3	NAG	B	2	1	-	2/6/23/26	0/1/1/1
3	NAG	B	5	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1	1	-	2/6/23/26	0/1/1/1
3	NAG	A	2	1	-	0/6/23/26	0/1/1/1
3	NAG	A	5	1	-	1/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	5	NAG	C1-C2	3.03	1.56	1.52
3	B	2	NAG	C1-C2	2.52	1.55	1.52
3	B	5	NAG	C4-C5	2.11	1.57	1.53
3	B	2	NAG	O5-C5	2.08	1.47	1.43
3	A	2	NAG	C1-C2	2.06	1.55	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	5	NAG	C4-C3-C2	-3.43	105.99	111.02
3	A	5	NAG	C4-C3-C2	-3.24	106.27	111.02
3	A	5	NAG	C1-C2-N2	2.91	115.02	110.43
3	A	2	NAG	C4-C3-C2	-2.43	107.46	111.02
3	B	5	NAG	C1-C2-N2	2.18	113.86	110.43
3	B	1	NAG	C2-N2-C7	-2.13	120.05	122.90

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	5	NAG	C1-C2-N2-C7
3	B	5	NAG	C1-C2-N2-C7
3	B	2	NAG	O5-C5-C6-O6
3	B	2	NAG	C4-C5-C6-O6
3	A	1	NAG	O5-C5-C6-O6
3	A	1	NAG	C4-C5-C6-O6
3	B	5	NAG	C3-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	368/371 (99%)	-0.50	9 (2%) 59 36	27, 47, 70, 95	0
1	B	368/371 (99%)	-0.49	9 (2%) 59 36	31, 48, 71, 93	0
All	All	736/742 (99%)	-0.49	18 (2%) 59 36	27, 48, 71, 95	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	51	ASN	10.0
1	A	51	ASN	7.4
1	A	379	LYS	5.0
1	A	378	LYS	4.8
1	B	378	LYS	4.7
1	A	52	SER	4.6
1	B	52	SER	4.1
1	A	50	LYS	4.1
1	B	379	LYS	4.0
1	B	300	ASP	2.8
1	B	377	ASP	2.6
1	A	145	LYS	2.6
1	B	49	LYS	2.5
1	B	50	LYS	2.4
1	A	300	ASP	2.3
1	A	377	ASP	2.2
1	A	53	GLY	2.1
1	B	364	GLN	2.0

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

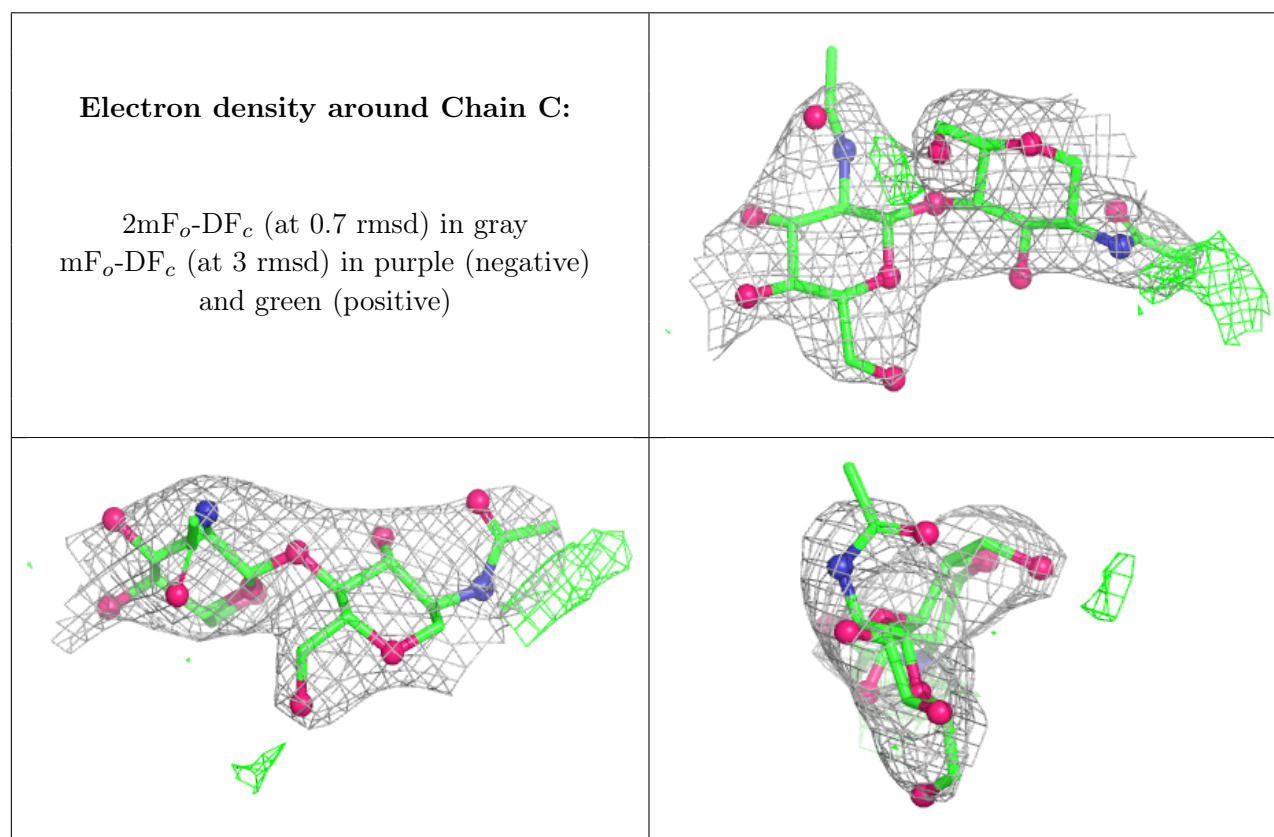
There are no non-standard protein/DNA/RNA residues in this entry.

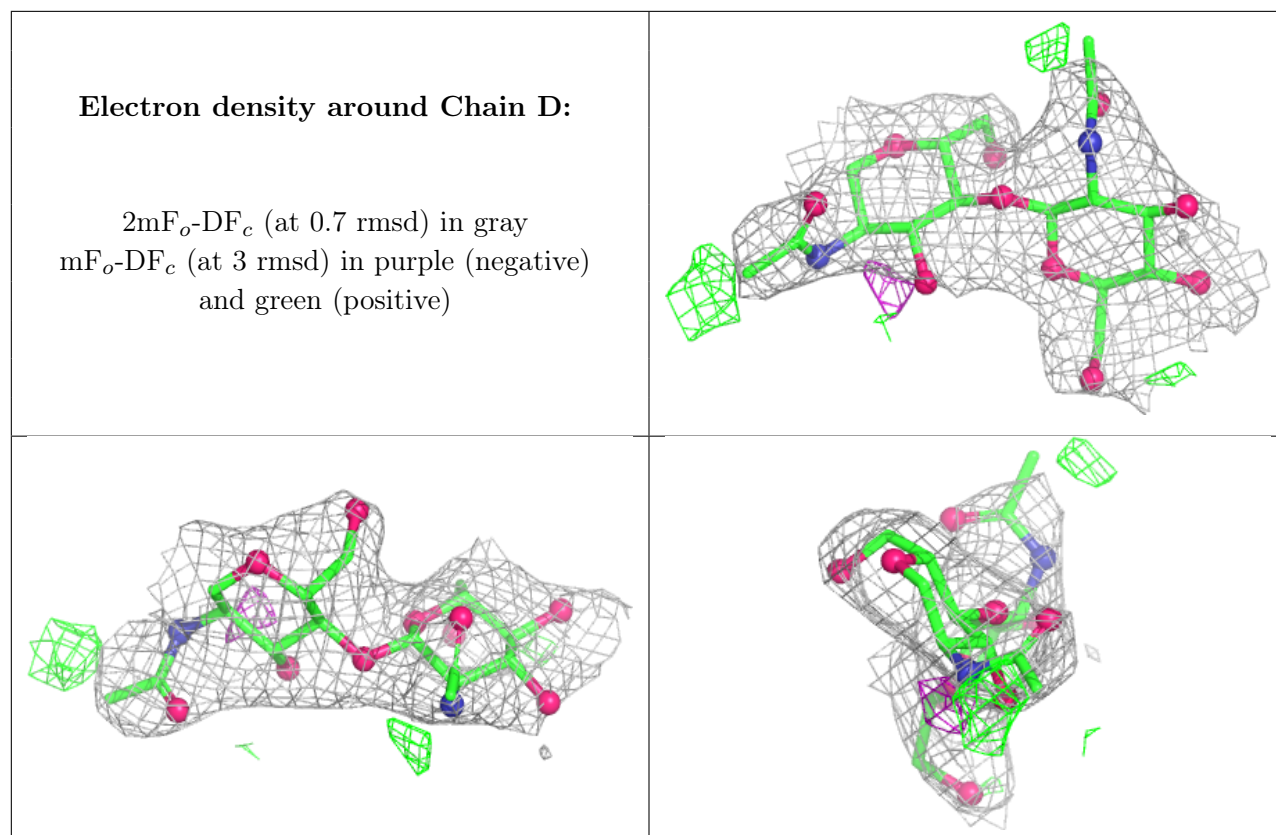
6.3 Carbohydrates [i](#)

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($\text{\AA}^2$)	Q<0.9
2	NAG	C	2	14/15	0.84	0.13	90,95,96,97	0
2	NAG	C	1	14/15	0.85	0.16	67,75,78,83	0
2	NAG	D	2	14/15	0.86	0.12	81,92,95,99	0
2	NAG	D	1	14/15	0.90	0.13	63,71,75,80	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($\text{\AA}^2$)	Q<0.9
3	NAG	B	5	14/15	0.69	0.20	86,91,96,96	0
3	NAG	A	5	14/15	0.71	0.19	83,86,89,90	0
3	NAG	B	1	14/15	0.78	0.17	80,85,92,93	0
3	NAG	A	1	14/15	0.78	0.16	82,88,91,92	0
3	NAG	B	2	14/15	0.88	0.13	51,58,63,66	0
3	NAG	A	2	14/15	0.88	0.14	59,61,75,75	0

6.5 Other polymers [i](#)

There are no such residues in this entry.