



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 09:10 PM BST

PDB ID : 5ANM
Title : Crystal structure of IgE Fc in complex with a neutralizing antibody
Authors : Cohen, E.S.; Dobson, C.L.; Kack, H.; Wang, B.; Sims, D.A.; Lloyd, C.O.; England, E.; Rees, D.G.; Guo, H.; Karagiannis, S.N.; O'Brien, S.; Persdotter, S.; Ekdahl, H.; Butler, R.; Keyes, F.; Oakley, S.; Carlsson, M.; Briend, E.; Wilkinson, T.; Anderson, I.K.; Monk, P.D.; vonWachenfeldt, K.; Eriksson, P.O.; Gould, H.J.; Vaughan, T.J.; May, R.D.
Deposited on : 2015-09-07
Resolution : 2.85 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

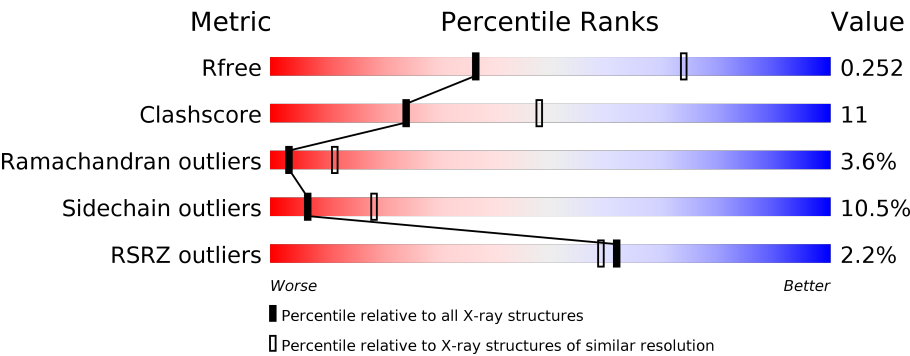
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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	217	<div><div></div><div>75%18%...</div></div>
1	C	217	<div><div></div><div>75%21%...</div></div>
1	L	217	<div><div>4%</div><div>58%24%5%13%</div></div>
2	B	229	<div><div>%</div><div>76%15%...</div></div>
2	D	229	<div><div></div><div>77%15%...</div></div>
2	H	229	<div><div>4%</div><div>54%29%7%9%</div></div>

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Mol	Chain	Length	Quality of chain
3	E	242	<div><div><div></div><div></div><div></div><div></div><div></div></div><div>4%47%31%7%15%</div></div>
3	F	242	<div><div><div></div><div></div><div></div><div></div><div></div></div><div>2%62%20%14%</div></div>
3	G	242	<div><div><div></div><div></div><div></div><div></div><div></div></div><div>2%63%18%6%13%</div></div>
4	I	6	<div><div><div></div><div></div><div></div><div></div><div></div></div><div>17%83%</div></div>
5	J	5	<div><div><div></div><div></div><div></div><div></div><div></div></div><div>80%20%</div></div>
6	K	6	<div><div><div></div><div></div><div></div><div></div><div></div></div><div>83%17%</div></div>

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 14788 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 IMMUNOGLOBULIN G.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	213	Total	C	N	O	S	0	0	0
			1582	991	262	325	4			
1	C	213	Total	C	N	O	S	0	0	0
			1582	991	262	325	4			
1	L	189	Total	C	N	O	S	0	0	0
			1383	865	228	286	4			

- Molecule 2 is a protein called IMMUNOGLOBULIN G.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	220	Total	C	N	O	S	0	0	0
			1673	1068	272	325	8			
2	D	219	Total	C	N	O	S	0	0	0
			1664	1062	270	324	8			
2	H	208	Total	C	N	O	S	0	0	0
			1580	1008	255	309	8			

- Molecule 3 is a protein called IG EPSILON CHAIN C REGION.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	206	Total	C	N	O	S	0	0	0
			1636	1026	301	303	6			
3	F	208	Total	C	N	O	S	0	0	0
			1628	1016	301	305	6			
3	G	211	Total	C	N	O	S	0	0	0
			1665	1042	307	310	6			

There are 72 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	326	ALA	-	expression tag	UNP P01854
E	327	ASP	-	expression tag	UNP P01854

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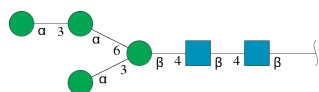
Chain	Residue	Modelled	Actual	Comment	Reference
E	328	PRO	-	expression tag	UNP P01854
E	329	CYS	-	expression tag	UNP P01854
E	548	ASP	-	expression tag	UNP P01854
E	549	TYR	-	expression tag	UNP P01854
E	550	LYS	-	expression tag	UNP P01854
E	551	ASP	-	expression tag	UNP P01854
E	552	ASP	-	expression tag	UNP P01854
E	553	ASP	-	expression tag	UNP P01854
E	554	ASP	-	expression tag	UNP P01854
E	555	LYS	-	expression tag	UNP P01854
E	556	ALA	-	expression tag	UNP P01854
E	557	ALA	-	expression tag	UNP P01854
E	558	HIS	-	expression tag	UNP P01854
E	559	HIS	-	expression tag	UNP P01854
E	560	HIS	-	expression tag	UNP P01854
E	561	HIS	-	expression tag	UNP P01854
E	562	HIS	-	expression tag	UNP P01854
E	563	HIS	-	expression tag	UNP P01854
E	564	HIS	-	expression tag	UNP P01854
E	565	HIS	-	expression tag	UNP P01854
E	566	HIS	-	expression tag	UNP P01854
E	567	HIS	-	expression tag	UNP P01854
F	326	ALA	-	expression tag	UNP P01854
F	327	ASP	-	expression tag	UNP P01854
F	328	PRO	-	expression tag	UNP P01854
F	329	CYS	-	expression tag	UNP P01854
F	548	ASP	-	expression tag	UNP P01854
F	549	TYR	-	expression tag	UNP P01854
F	550	LYS	-	expression tag	UNP P01854
F	551	ASP	-	expression tag	UNP P01854
F	552	ASP	-	expression tag	UNP P01854
F	553	ASP	-	expression tag	UNP P01854
F	554	ASP	-	expression tag	UNP P01854
F	555	LYS	-	expression tag	UNP P01854
F	556	ALA	-	expression tag	UNP P01854
F	557	ALA	-	expression tag	UNP P01854
F	558	HIS	-	expression tag	UNP P01854
F	559	HIS	-	expression tag	UNP P01854
F	560	HIS	-	expression tag	UNP P01854
F	561	HIS	-	expression tag	UNP P01854
F	562	HIS	-	expression tag	UNP P01854
F	563	HIS	-	expression tag	UNP P01854

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Chain	Residue	Modelled	Actual	Comment	Reference
F	564	HIS	-	expression tag	UNP P01854
F	565	HIS	-	expression tag	UNP P01854
F	566	HIS	-	expression tag	UNP P01854
F	567	HIS	-	expression tag	UNP P01854
G	326	ALA	-	expression tag	UNP P01854
G	327	ASP	-	expression tag	UNP P01854
G	328	PRO	-	expression tag	UNP P01854
G	329	CYS	-	expression tag	UNP P01854
G	548	ASP	-	expression tag	UNP P01854
G	549	TYR	-	expression tag	UNP P01854
G	550	LYS	-	expression tag	UNP P01854
G	551	ASP	-	expression tag	UNP P01854
G	552	ASP	-	expression tag	UNP P01854
G	553	ASP	-	expression tag	UNP P01854
G	554	ASP	-	expression tag	UNP P01854
G	555	LYS	-	expression tag	UNP P01854
G	556	ALA	-	expression tag	UNP P01854
G	557	ALA	-	expression tag	UNP P01854
G	558	HIS	-	expression tag	UNP P01854
G	559	HIS	-	expression tag	UNP P01854
G	560	HIS	-	expression tag	UNP P01854
G	561	HIS	-	expression tag	UNP P01854
G	562	HIS	-	expression tag	UNP P01854
G	563	HIS	-	expression tag	UNP P01854
G	564	HIS	-	expression tag	UNP P01854
G	565	HIS	-	expression tag	UNP P01854
G	566	HIS	-	expression tag	UNP P01854
G	567	HIS	-	expression tag	UNP P01854

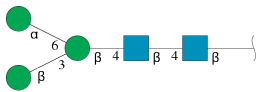
- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	I	6	Total	C	N	O	0	0	0
			72	40	2	30			

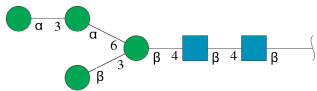
- Molecule 5 is an oligosaccharide called beta-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

se-(1-6)][beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	J	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[beta-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	K	6	Total	C	N	O	0	0	0
			72	40	2	30			

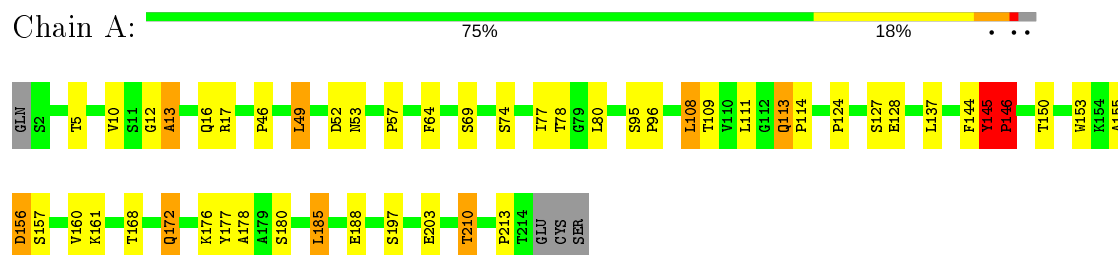
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	14	Total	O	0	0
			14	14		
7	B	60	Total	O	0	0
			60	60		
7	C	16	Total	O	0	0
			16	16		
7	D	60	Total	O	0	0
			60	60		
7	E	6	Total	O	0	0
			6	6		
7	F	10	Total	O	0	0
			10	10		
7	G	22	Total	O	0	0
			22	22		
7	H	2	Total	O	0	0
			2	2		

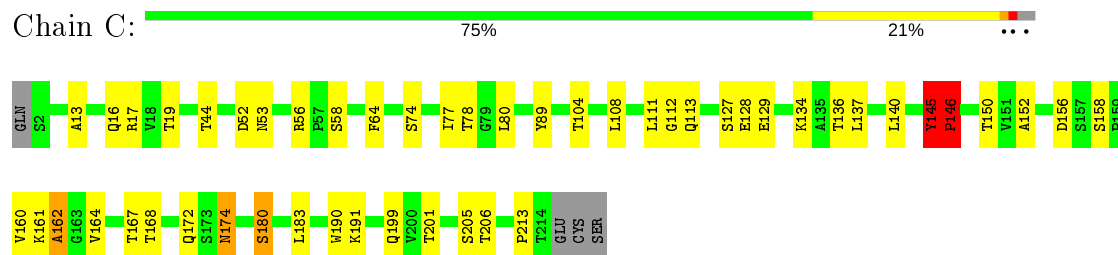
3 Residue-property plots [i](#)

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.

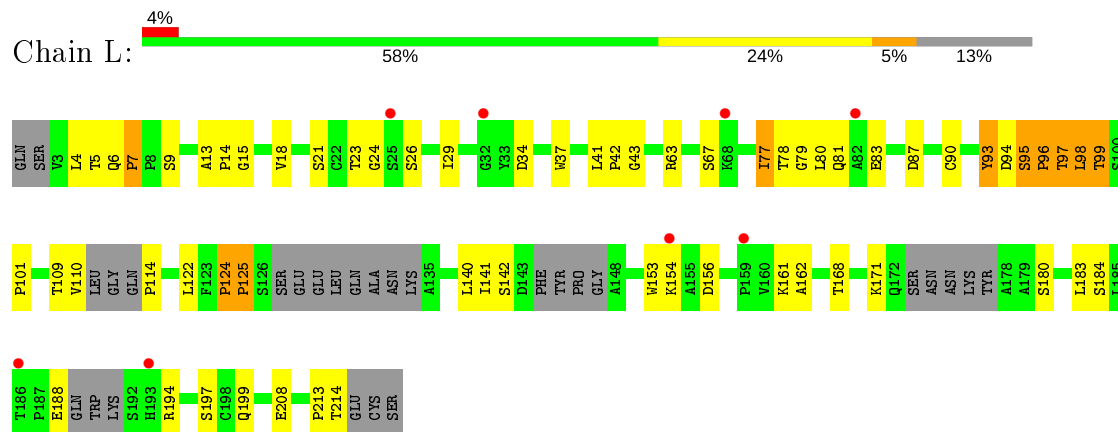
• Molecule 1: IMMUNOGLOBULIN G



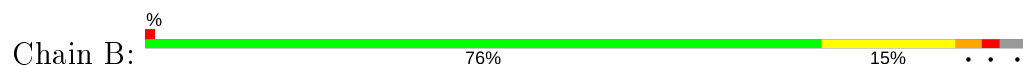
• Molecule 1: IMMUNOGLOBULIN G

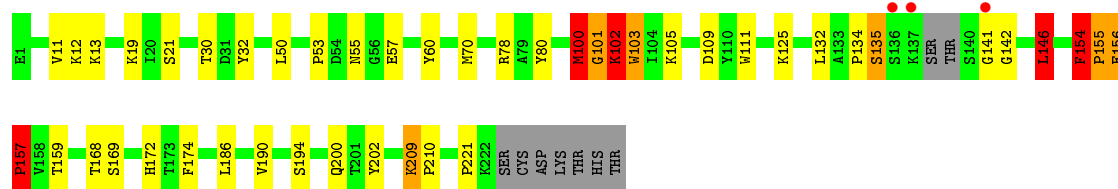


• Molecule 1: IMMUNOGLOBULIN G

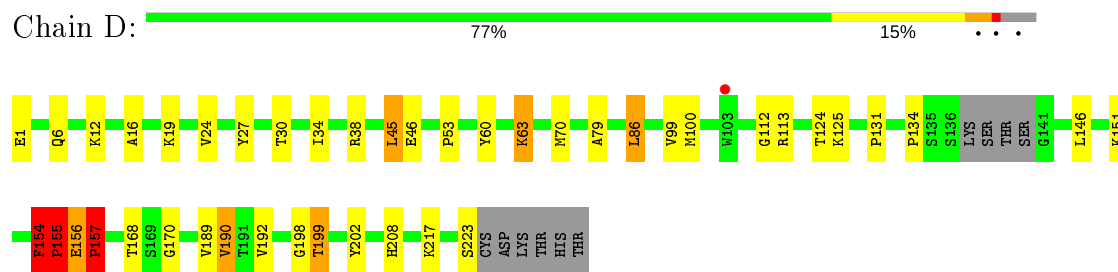


• Molecule 2: IMMUNOGLOBULIN G

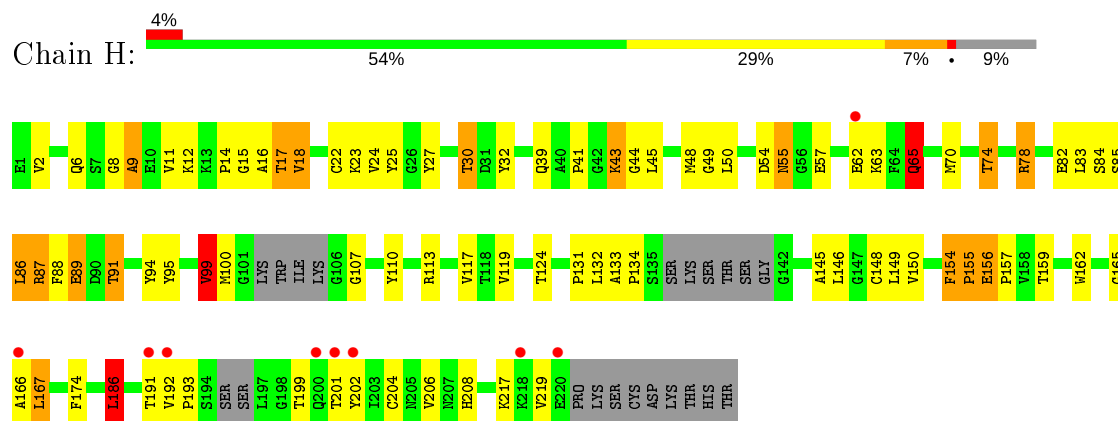




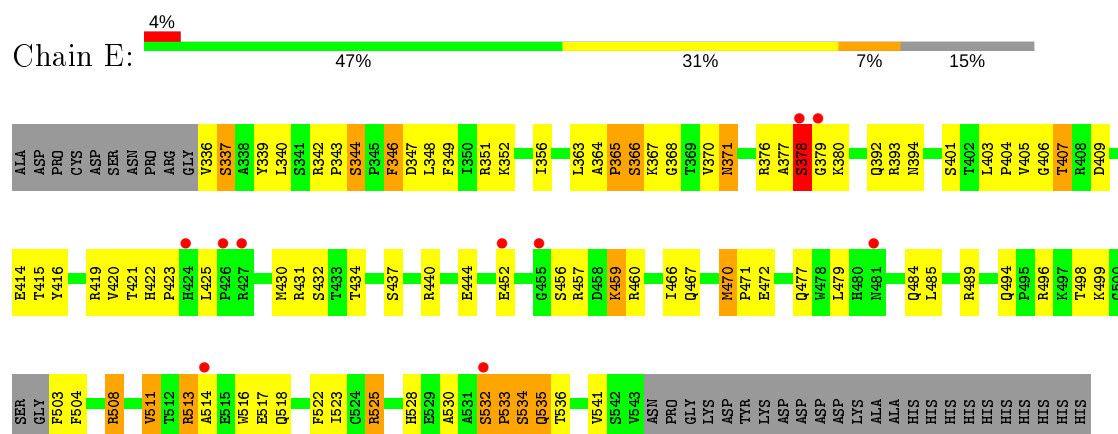
• Molecule 2: IMMUNOGLOBULIN G



• Molecule 2: IMMUNOGLOBULIN G

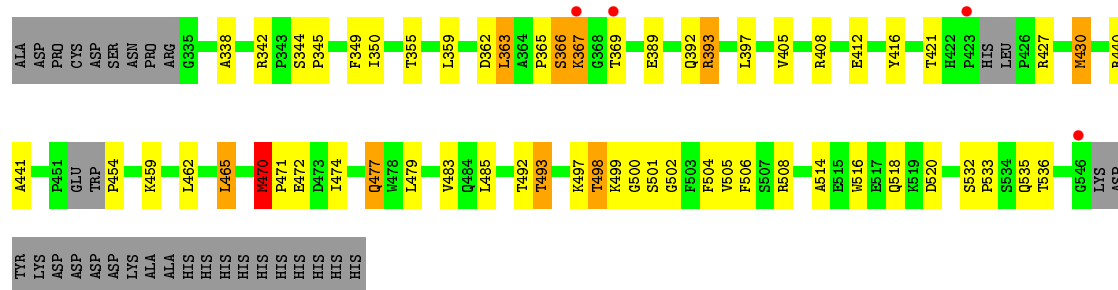


• Molecule 3: IG EPSILON CHAIN C REGION

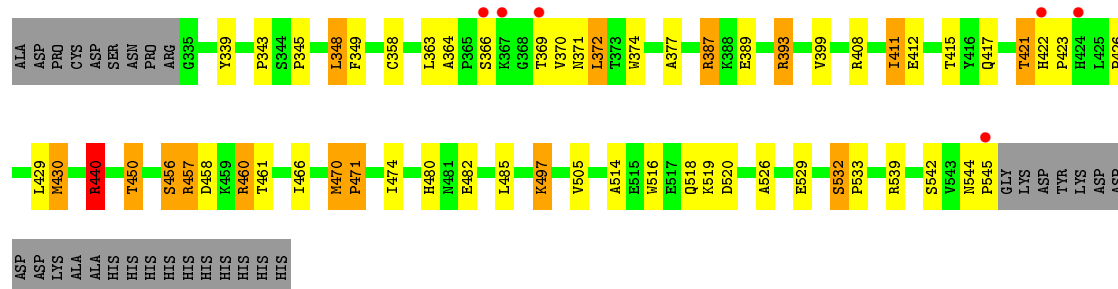


• Molecule 3: IG EPSILON CHAIN C REGION





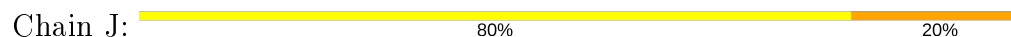
• Molecule 3: IG EPSILON CHAIN C REGION



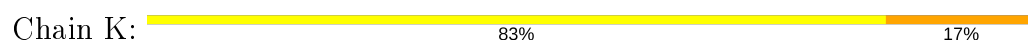
• Molecule 4: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



• Molecule 5: beta-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



• Molecule 6: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[beta-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-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	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	140.96Å 140.96Å 244.45Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	31.29 – 2.85 31.28 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.9 (31.29-2.85) 100.0 (31.28-2.85)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.28 (at 2.85Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.196 , 0.260 0.192 , 0.252	Depositor DCC
R_{free} test set	3355 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	57.9	Xtriage
Anisotropy	0.195	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 46.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14788	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% 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: BMA, NAG, MAN

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	0.64	0/1623	0.81	2/2222 (0.1%)
1	C	0.63	0/1623	0.82	2/2222 (0.1%)
1	L	0.61	3/1413 (0.2%)	0.74	1/1930 (0.1%)
2	B	0.79	0/1715	1.02	8/2332 (0.3%)
2	D	0.78	0/1706	1.11	6/2321 (0.3%)
2	H	0.54	0/1617	0.87	3/2199 (0.1%)
3	E	0.54	0/1677	0.79	2/2282 (0.1%)
3	F	0.62	0/1666	0.93	6/2262 (0.3%)
3	G	0.70	0/1708	0.90	3/2326 (0.1%)
All	All	0.66	3/14748 (0.0%)	0.90	33/20096 (0.2%)

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	1
1	C	0	2
1	L	0	1
2	B	0	4
2	D	0	1
2	H	0	2
3	E	0	3
3	F	0	1
3	G	0	2
All	All	0	17

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	14	PRO	C-O	7.56	1.38	1.23
1	L	14	PRO	C-N	6.41	1.44	1.33
1	L	13	ALA	C-N	-5.40	1.24	1.34

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	156	GLU	C-N-CD	-21.33	73.67	120.60
3	F	470	MET	C-N-CD	-16.27	84.80	120.60
2	H	156	GLU	C-N-CD	-13.12	91.74	120.60
2	D	156	GLU	C-N-CA	12.97	176.46	122.00
2	D	154	PHE	C-N-CD	-12.80	92.45	120.60
2	B	156	GLU	C-N-CD	-10.98	96.45	120.60
3	G	470	MET	C-N-CD	-9.53	99.62	120.60
2	B	78	ARG	NE-CZ-NH1	-9.25	115.68	120.30
2	B	109	ASP	CB-CG-OD1	-8.46	110.69	118.30
1	C	145	TYR	C-N-CD	-7.70	103.66	120.60
2	B	154	PHE	C-N-CD	-7.64	103.78	120.60
3	E	470	MET	C-N-CD	-7.44	104.22	120.60
3	F	470	MET	C-N-CA	7.30	152.64	122.00
2	D	155	PRO	N-CA-C	-7.28	93.18	112.10
1	A	145	TYR	C-N-CD	-6.89	105.44	120.60
2	B	146	LEU	CA-CB-CG	6.34	129.89	115.30
3	E	337	SER	N-CA-C	6.15	127.60	111.00
3	G	440	ARG	NE-CZ-NH1	6.12	123.36	120.30
3	F	342	ARG	NE-CZ-NH1	-6.11	117.24	120.30
2	H	86	LEU	CA-CB-CG	6.11	129.35	115.30
2	D	154	PHE	C-N-CA	6.06	147.45	122.00
2	B	100	MET	CB-CG-SD	-6.03	94.31	112.40
2	B	157	PRO	N-CA-C	-5.95	96.64	112.10
2	D	157	PRO	CA-N-CD	-5.73	103.48	111.50
2	B	101	GLY	N-CA-C	-5.61	99.08	113.10
3	F	430	MET	CG-SD-CE	-5.47	91.45	100.20
3	F	470	MET	N-CA-C	5.42	125.63	111.00
2	H	186	LEU	CA-CB-CG	5.29	127.46	115.30
1	C	104	THR	N-CA-CB	-5.14	100.54	110.30
1	A	49	LEU	CA-CB-CG	5.12	127.07	115.30
3	G	440	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	L	98	LEU	CA-CB-CG	5.03	126.88	115.30
3	F	440	ARG	NE-CZ-NH2	-5.02	117.79	120.30

There are no chirality outliers.

All (17) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	145	TYR	Peptide
2	B	154	PHE	Mainchain,Peptide
2	B	156	GLU	Mainchain,Peptide
1	C	145	TYR	Peptide
1	C	160	VAL	Peptide
2	D	154	PHE	Peptide
3	E	336	VAL	Peptide
3	E	366	SER	Peptide
3	E	470	MET	Peptide
3	F	470	MET	Peptide
3	G	470	MET	Peptide
3	G	532	SER	Peptide
2	H	154	PHE	Peptide
2	H	156	GLU	Peptide
1	L	114	PRO	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	1582	0	1528	32	0
1	C	1582	0	1528	23	0
1	L	1383	0	1339	34	0
2	B	1673	0	1642	26	0
2	D	1664	0	1629	26	0
2	H	1580	0	1537	64	0
3	E	1636	0	1622	62	0
3	F	1628	0	1616	33	0
3	G	1665	0	1647	38	0
4	I	72	0	61	0	0
5	J	61	0	52	1	0
6	K	72	0	61	1	0
7	A	14	0	0	0	0
7	B	60	0	0	2	0
7	C	16	0	0	0	0
7	D	60	0	0	2	0
7	E	6	0	0	1	0
7	F	10	0	0	0	0
7	G	22	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	H	2	0	0	0	0
All	All	14788	0	14262	318	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 11.

All (318) 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)
3:G:371:ASN:HB2	3:G:421:THR:HG22	1.33	1.06
3:F:514:ALA:HB2	2:H:165:GLY:HA3	1.41	1.01
3:G:371:ASN:HB2	3:G:421:THR:CG2	1.96	0.96
3:E:508:ARG:HG3	3:E:508:ARG:HH11	1.30	0.95
2:H:30:THR:HG22	2:H:54:ASP:HA	1.48	0.92
2:H:43:LYS:HD2	2:H:44:GLY:H	1.38	0.86
2:H:192:VAL:HG22	2:H:193:PRO:HD2	1.56	0.86
2:H:6:GLN:HG2	2:H:22:CYS:HB3	1.59	0.85
2:H:62:GLU:HA	2:H:65:GLN:HG3	1.60	0.84
1:A:155:ALA:O	1:A:156:ASP:HB2	1.76	0.84
2:H:154:PHE:CD2	2:H:155:PRO:HD3	2.12	0.84
1:A:160:VAL:HG23	1:A:161:LYS:H	1.44	0.83
3:E:477:GLN:HG2	3:E:484:GLN:HE22	1.42	0.83
1:A:145:TYR:CD2	1:A:146:PRO:HD3	2.16	0.80
3:E:365:PRO:O	3:E:367:LYS:N	2.14	0.79
2:H:88:PHE:O	2:H:91:THR:HG23	1.82	0.79
1:C:161:LYS:HE2	1:C:164:VAL:HB	1.64	0.79
3:G:440:ARG:HH11	3:G:440:ARG:HG2	1.47	0.78
2:H:150:VAL:HG12	2:H:186:LEU:HD12	1.66	0.77
2:B:134:PRO:HD3	2:B:146:LEU:HB3	1.66	0.76
1:A:145:TYR:CD2	1:A:146:PRO:CD	2.68	0.75
1:A:160:VAL:HG23	1:A:161:LYS:N	2.02	0.74
2:H:150:VAL:CG1	2:H:186:LEU:HD12	2.16	0.74
2:H:99:VAL:O	2:H:100:MET:SD	2.45	0.74
3:F:497:LYS:NZ	3:F:501:SER:HA	2.02	0.73
3:G:387:ARG:NH1	3:G:389:GLU:HG3	2.04	0.72
2:H:30:THR:HG22	2:H:54:ASP:CA	2.20	0.72
3:E:367:LYS:HD2	3:E:423:PRO:HB2	1.72	0.71
3:E:525:ARG:HD3	3:E:536:THR:HG23	1.71	0.71
3:E:532:SER:N	3:E:533:PRO:HD2	2.04	0.71
2:H:30:THR:CG2	2:H:54:ASP:HA	2.19	0.71
1:C:145:TYR:CD2	1:C:146:PRO:HD3	2.26	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:38:ARG:NH1	2:D:46:GLU:OE2	2.25	0.69
3:G:417:GLN:OE1	3:G:430:MET:HE1	1.92	0.69
2:D:100:MET:HB2	3:E:394:ASN:HD22	1.58	0.69
2:B:134:PRO:HD3	2:B:146:LEU:CB	2.24	0.68
3:E:365:PRO:C	3:E:367:LYS:H	1.97	0.67
1:A:172:GLN:HG3	1:A:176:LYS:O	1.95	0.67
2:H:50:LEU:N	2:H:70:MET:HE3	2.11	0.67
3:G:456:SER:O	3:G:458:ASP:N	2.29	0.66
3:E:344:SER:HB3	3:E:347:ASP:OD2	1.96	0.66
2:H:124:THR:HA	2:H:155:PRO:HD2	1.78	0.65
1:C:167:THR:HG22	1:C:168:THR:O	1.95	0.65
2:H:91:THR:HG22	2:H:119:VAL:H	1.61	0.65
3:E:508:ARG:CG	3:E:508:ARG:HH11	2.08	0.64
1:L:124:PRO:HB2	1:L:125:PRO:CD	2.26	0.64
2:H:48:MET:CE	2:H:94:TYR:HD2	2.10	0.64
3:G:544:ASN:HB3	3:G:545:PRO:HD2	1.78	0.64
3:E:415:THR:OG1	3:E:434:THR:CG2	2.46	0.63
3:F:493:THR:HG22	3:F:505:VAL:HA	1.81	0.63
3:E:349:PHE:HB3	3:E:535:GLN:HG3	1.82	0.62
1:L:124:PRO:HB2	1:L:125:PRO:HD2	1.81	0.62
2:H:174:PHE:CE1	1:L:140:LEU:HD23	2.35	0.62
1:L:154:LYS:HD3	1:L:197:SER:HB2	1.82	0.62
1:A:49:LEU:O	1:A:57:PRO:HD2	1.98	0.62
1:A:145:TYR:CE2	1:A:146:PRO:HD3	2.34	0.61
1:L:18:VAL:HG22	1:L:80:LEU:HD11	1.82	0.61
3:F:441:ALA:HB3	3:F:470:MET:HB2	1.82	0.61
2:B:101:GLY:O	2:B:102:LYS:HB2	2.01	0.60
3:G:372:LEU:HD23	3:G:372:LEU:C	2.22	0.60
3:E:430:MET:O	3:E:431:ARG:NH1	2.34	0.60
2:H:48:MET:HE1	2:H:94:TYR:HD2	1.65	0.60
3:E:508:ARG:HG3	3:E:508:ARG:NH1	2.08	0.59
2:H:91:THR:CG2	2:H:119:VAL:H	2.15	0.59
2:H:154:PHE:CD2	2:H:155:PRO:CD	2.84	0.59
1:A:52:ASP:O	1:A:53:ASN:HB2	2.03	0.58
3:E:364:ALA:N	3:E:365:PRO:CD	2.67	0.58
3:E:337:SER:HB2	7:E:2001:HOH:O	2.02	0.58
1:L:81:GLN:HB3	1:L:83:GLU:HG2	1.86	0.58
3:F:365:PRO:C	3:F:367:LYS:N	2.55	0.58
3:F:516:TRP:O	3:F:520:ASP:HB2	2.03	0.58
1:A:197:SER:OG	1:A:210:THR:HB	2.03	0.58
1:C:167:THR:CG2	1:C:168:THR:O	2.51	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ARG:HG3	1:A:78:THR:HG22	1.84	0.57
3:F:498:THR:HB	3:F:502:GLY:O	2.05	0.57
3:F:365:PRO:C	3:F:367:LYS:H	2.06	0.57
2:H:11:VAL:O	2:H:11:VAL:HG12	2.03	0.57
1:C:137:LEU:HD22	1:C:183:LEU:HD23	1.87	0.57
1:C:161:LYS:O	1:C:162:ALA:HB3	2.05	0.57
3:E:432:SER:H	2:H:74:THR:HB	1.70	0.57
3:F:465:LEU:HG	3:F:506:PHE:CE1	2.40	0.56
2:D:34:ILE:HD12	2:D:79:ALA:HB2	1.87	0.56
1:C:167:THR:HB	1:C:180:SER:H	1.69	0.56
2:H:24:VAL:HG11	2:H:27:TYR:CE1	2.39	0.56
3:E:414:GLU:O	3:E:434:THR:HG22	2.06	0.56
2:B:125:LYS:HA	1:C:128:GLU:HG2	1.87	0.56
2:B:102:LYS:HA	7:B:2046:HOH:O	2.07	0.55
3:G:514:ALA:O	3:G:518:GLN:HB2	2.06	0.55
2:D:131:PRO:HD3	2:D:217:LYS:HE2	1.88	0.55
3:E:376:ARG:NH2	3:E:409:ASP:OD2	2.40	0.55
2:D:16:ALA:O	2:D:86:LEU:HB2	2.06	0.55
1:A:153:TRP:HB2	1:A:160:VAL:CG2	2.37	0.55
3:G:349:PHE:CE1	3:G:411:ILE:HD11	2.42	0.55
1:C:80:LEU:HD11	1:C:108:LEU:HD21	1.89	0.54
1:L:41:LEU:HB3	1:L:42:PRO:HD3	1.90	0.54
3:G:417:GLN:OE1	3:G:430:MET:CE	2.54	0.54
2:H:150:VAL:HG12	2:H:186:LEU:O	2.07	0.54
2:H:17:THR:OG1	2:H:84:SER:HA	2.08	0.54
2:H:8:GLY:O	2:H:9:ALA:O	2.26	0.54
1:L:94:ASP:O	1:L:96:PRO:HD2	2.08	0.54
3:E:479:LEU:HB2	3:E:523:ILE:HB	1.89	0.53
2:H:48:MET:CE	2:H:94:TYR:CD2	2.91	0.53
1:L:94:ASP:C	1:L:96:PRO:HD2	2.28	0.53
1:C:129:GLU:OE2	2:D:151:LYS:NZ	2.40	0.53
2:B:101:GLY:O	2:B:102:LYS:CB	2.56	0.53
3:F:408:ARG:O	3:F:412:GLU:HG3	2.08	0.53
2:H:2:VAL:HB	2:H:110:TYR:CE1	2.43	0.53
3:G:440:ARG:HH11	3:G:440:ARG:CG	2.19	0.53
3:E:431:ARG:HD2	2:H:74:THR:HG21	1.91	0.53
3:F:393:ARG:HH21	2:H:107:GLY:H	1.56	0.53
2:D:63:LYS:HB3	7:D:2028:HOH:O	2.08	0.52
3:G:387:ARG:HH12	3:G:389:GLU:HG3	1.74	0.52
1:C:13:ALA:O	1:C:16:GLN:HB2	2.10	0.52
3:E:343:PRO:HD3	3:E:356:ILE:HG22	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:508:ARG:CG	3:E:508:ARG:NH1	2.69	0.52
3:E:377:ALA:O	3:E:378:SER:HB2	2.10	0.52
1:A:160:VAL:CG2	1:A:161:LYS:H	2.19	0.52
1:C:17:ARG:HG3	1:C:78:THR:HG22	1.91	0.52
3:E:467:GLN:OE1	3:F:508:ARG:NE	2.43	0.52
3:E:430:MET:CG	3:E:431:ARG:N	2.73	0.52
1:L:124:PRO:CB	1:L:125:PRO:CD	2.87	0.52
3:G:339:TYR:CE2	6:K:3:BMA:H3	2.45	0.52
1:L:81:GLN:HB3	1:L:83:GLU:H	1.75	0.52
2:D:24:VAL:HG13	2:D:27:TYR:CE1	2.45	0.51
2:H:32:TYR:CE2	2:H:100:MET:HG3	2.45	0.51
1:L:6:GLN:HE22	1:L:90:CYS:H	1.58	0.51
3:F:366:SER:O	3:F:367:LYS:HG3	2.11	0.51
1:A:128:GLU:HG2	2:D:125:LYS:HA	1.92	0.51
3:G:480:HIS:HE1	3:G:519:LYS:O	1.94	0.51
3:E:496:ARG:HB2	3:E:504:PHE:CE1	2.46	0.51
2:H:15:GLY:O	2:H:85:SER:HA	2.10	0.51
1:C:89:TYR:CD2	2:D:45:LEU:HD22	2.46	0.51
2:H:24:VAL:CG1	2:H:27:TYR:CE1	2.93	0.51
3:F:389:GLU:HG3	3:F:397:LEU:HD11	1.93	0.51
1:L:93:TYR:CE2	1:L:98:LEU:HA	2.45	0.51
1:A:64:PHE:CD1	1:A:77:ILE:HG12	2.46	0.50
2:B:50:LEU:C	2:B:50:LEU:HD12	2.31	0.50
3:E:534:SER:C	3:E:535:GLN:HG2	2.31	0.50
3:E:405:VAL:HB	3:E:416:TYR:CZ	2.46	0.50
2:D:157:PRO:O	2:D:157:PRO:HD2	2.12	0.50
1:L:4:LEU:HD22	1:L:29:ILE:HD11	1.92	0.50
3:E:351:ARG:O	3:E:352:LYS:HB2	2.11	0.50
2:H:154:PHE:O	2:H:208:HIS:CE1	2.65	0.49
3:F:345:PRO:HG2	3:F:474:ILE:HA	1.95	0.49
2:D:198:GLY:HA3	3:G:532:SER:HB3	1.94	0.49
1:A:46:PRO:HG2	2:B:111:TRP:CZ3	2.48	0.49
1:C:190:TRP:CZ2	1:C:213:PRO:HA	2.47	0.49
1:L:95:SER:HB3	1:L:96:PRO:CD	2.42	0.49
2:H:145:ALA:HB2	2:H:191:THR:HG22	1.93	0.49
3:G:440:ARG:HG2	3:G:440:ARG:NH1	2.20	0.49
1:A:172:GLN:OE1	1:A:178:ALA:HB2	2.13	0.49
3:E:459:LYS:O	3:E:460:ARG:HG3	2.12	0.49
1:L:77:ILE:HD11	1:L:80:LEU:HG	1.92	0.49
3:E:511:VAL:HG21	3:E:522:PHE:CE2	2.48	0.49
1:L:94:ASP:OD2	1:L:99:THR:HG22	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:134:PRO:HG2	2:B:221:PRO:HA	1.94	0.49
3:G:466:ILE:HD13	3:G:526:ALA:HB2	1.94	0.49
2:B:60:TYR:CE2	2:B:70:MET:HE2	2.48	0.48
2:D:19:LYS:NZ	7:D:2013:HOH:O	2.46	0.48
3:F:392:GLN:OE1	3:F:392:GLN:HA	2.12	0.48
2:D:100:MET:HB2	3:E:394:ASN:ND2	2.26	0.48
3:F:497:LYS:HZ2	3:F:501:SER:HA	1.77	0.48
3:G:440:ARG:HE	3:G:529:GLU:CD	2.15	0.48
2:B:11:VAL:O	2:B:12:LYS:HD3	2.13	0.48
3:G:364:ALA:C	3:G:366:SER:H	2.17	0.48
1:L:24:GLY:HA3	1:L:29:ILE:HD13	1.94	0.48
1:C:152:ALA:O	1:C:199:GLN:HB2	2.13	0.48
3:G:371:ASN:CB	3:G:421:THR:HG22	2.24	0.48
3:G:460:ARG:HD3	3:G:460:ARG:HA	1.59	0.48
2:D:155:PRO:HD2	2:D:208:HIS:NE2	2.28	0.48
2:H:91:THR:HG22	2:H:119:VAL:HB	1.95	0.48
2:B:135:SER:HB2	7:B:2053:HOH:O	2.12	0.48
2:B:146:LEU:HD22	2:B:190:VAL:CG1	2.43	0.48
2:B:30:THR:HA	2:B:53:PRO:HB2	1.95	0.48
3:E:422:HIS:HB3	3:E:425:LEU:HB2	1.96	0.48
1:A:13:ALA:H	1:A:111:LEU:H	1.62	0.48
2:D:134:PRO:HB3	2:D:146:LEU:HB3	1.95	0.48
1:C:145:TYR:CD2	1:C:146:PRO:CD	2.97	0.47
3:G:497:LYS:HB2	3:G:497:LYS:HE3	1.68	0.47
1:L:41:LEU:HB3	1:L:42:PRO:CD	2.45	0.47
1:C:52:ASP:O	1:C:53:ASN:HB2	2.15	0.47
2:B:105:LYS:HD3	2:B:105:LYS:HA	1.73	0.46
3:E:508:ARG:HD2	3:F:504:PHE:CE1	2.50	0.46
2:H:133:ALA:HB1	2:H:134:PRO:HD2	1.97	0.46
1:L:37:TRP:CZ3	1:L:90:CYS:HB3	2.50	0.46
3:F:349:PHE:CG	3:F:535:GLN:HG2	2.49	0.46
3:G:544:ASN:O	3:G:545:PRO:C	2.54	0.46
1:A:137:LEU:HD11	1:A:185:LEU:HD11	1.96	0.46
1:A:95:SER:N	1:A:96:PRO:CD	2.79	0.46
3:E:376:ARG:HD2	3:E:416:TYR:CE1	2.49	0.46
1:C:64:PHE:CD1	1:C:77:ILE:HG12	2.51	0.46
2:H:192:VAL:HG22	2:H:193:PRO:CD	2.37	0.46
1:L:93:TYR:HE2	1:L:98:LEU:HA	1.81	0.46
1:C:129:GLU:OE2	1:C:136:THR:OG1	2.32	0.46
3:E:403:LEU:HD12	3:E:404:PRO:HD2	1.98	0.46
3:E:430:MET:HG2	3:E:431:ARG:N	2.30	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:154:PHE:CG	2:H:155:PRO:N	2.84	0.46
2:H:192:VAL:HG21	2:H:202:TYR:CE2	2.51	0.46
3:F:362:ASP:OD2	3:F:365:PRO:HA	2.16	0.45
1:L:97:THR:O	1:L:98:LEU:HB3	2.16	0.45
2:D:12:LYS:HA	2:D:12:LYS:HD2	1.66	0.45
1:L:96:PRO:O	1:L:97:THR:OG1	2.25	0.45
2:B:209:LYS:N	2:B:210:PRO:CD	2.79	0.45
1:A:113:GLN:HA	1:A:114:PRO:HD3	1.80	0.45
2:D:6:GLN:OE1	2:D:112:GLY:HA3	2.17	0.45
3:E:378:SER:O	3:E:380:LYS:N	2.39	0.45
3:E:514:ALA:HA	3:E:517:GLU:HB2	1.99	0.45
1:L:154:LYS:HB2	1:L:197:SER:HB2	1.99	0.45
2:H:146:LEU:HD13	2:H:219:VAL:HG21	1.99	0.45
2:H:132:LEU:HD13	2:H:149:LEU:HB2	1.99	0.45
1:A:153:TRP:HB2	1:A:160:VAL:HG21	1.99	0.45
2:B:142:GLY:O	2:B:194:SER:HB3	2.16	0.45
1:L:154:LYS:HD2	1:L:199:GLN:NE2	2.31	0.45
1:L:99:THR:O	1:L:101:PRO:HD3	2.17	0.45
2:D:60:TYR:CE1	2:D:70:MET:HE2	2.52	0.45
1:L:153:TRP:CD2	1:L:183:LEU:HD12	2.52	0.45
3:E:479:LEU:HD21	3:E:525:ARG:HB2	1.98	0.44
3:E:499:LYS:N	3:E:499:LYS:HD2	2.33	0.44
3:G:389:GLU:HG2	3:G:399:VAL:HG22	1.98	0.44
3:E:343:PRO:HD3	3:E:356:ILE:CG2	2.47	0.44
3:G:345:PRO:HG2	3:G:474:ILE:HA	1.99	0.44
1:L:161:LYS:O	1:L:162:ALA:HB3	2.17	0.44
3:F:498:THR:HG23	3:F:499:LYS:N	2.32	0.44
2:H:154:PHE:O	2:H:208:HIS:NE2	2.50	0.44
1:L:6:GLN:HA	1:L:7:PRO:HD2	1.76	0.44
3:G:358:CYS:HB2	3:G:374:TRP:CH2	2.52	0.44
2:H:49:GLY:HA3	2:H:70:MET:HE1	2.00	0.44
3:E:364:ALA:N	3:E:365:PRO:HD2	2.32	0.44
3:E:528:HIS:CD2	3:E:530:ALA:H	2.36	0.44
3:E:532:SER:N	3:E:533:PRO:CD	2.76	0.44
3:F:363:LEU:HD22	5:J:1:NAG:H83	2.00	0.44
2:B:100:MET:H	2:B:100:MET:HG2	1.05	0.44
3:G:450:THR:HB	3:G:461:THR:O	2.18	0.44
3:G:440:ARG:CZ	3:G:471:PRO:HG3	2.48	0.44
1:L:6:GLN:NE2	1:L:90:CYS:SG	2.91	0.43
1:A:13:ALA:O	1:A:16:GLN:HB2	2.18	0.43
1:A:144:PHE:CE1	1:A:177:TYR:HB2	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:190:VAL:HG22	2:D:192:VAL:HG13	1.99	0.43
2:H:39:GLN:O	2:H:39:GLN:HG2	2.18	0.43
3:E:525:ARG:HH11	3:E:536:THR:HG21	1.82	0.43
2:H:48:MET:HE2	2:H:94:TYR:CD2	2.53	0.43
2:D:192:VAL:HG11	2:D:202:TYR:CE1	2.54	0.43
3:E:472:GLU:OE1	2:H:78:ARG:NH2	2.51	0.43
2:H:43:LYS:CD	2:H:44:GLY:H	2.20	0.43
2:H:6:GLN:OE1	2:H:95:TYR:HA	2.18	0.43
2:B:55:ASN:CG	2:B:57:GLU:HG3	2.39	0.43
2:D:30:THR:HA	2:D:53:PRO:HB2	2.01	0.43
1:A:188:GLU:H	1:A:188:GLU:CD	2.22	0.43
1:L:37:TRP:CH2	1:L:90:CYS:HB3	2.54	0.43
1:A:124:PRO:HA	1:A:137:LEU:HD23	2.00	0.42
1:A:12:GLY:HA3	1:A:80:LEU:CD2	2.49	0.42
2:B:103:TRP:HA	2:B:103:TRP:CE3	2.53	0.42
3:E:525:ARG:HD3	3:E:536:THR:CG2	2.44	0.42
3:E:444:GLU:OE2	3:F:454:PRO:HG2	2.18	0.42
3:G:393:ARG:C	3:G:393:ARG:HD3	2.39	0.42
3:E:415:THR:OG1	3:E:434:THR:HG21	2.20	0.42
2:B:200:GLN:HG2	2:B:202:TYR:CZ	2.54	0.42
3:F:405:VAL:HG12	3:F:416:TYR:CE2	2.55	0.42
3:F:493:THR:HG22	3:F:505:VAL:CA	2.48	0.42
3:E:342:ARG:NH2	3:E:434:THR:O	2.44	0.42
2:D:124:THR:HA	2:D:154:PHE:O	2.19	0.42
3:E:365:PRO:C	3:E:367:LYS:N	2.64	0.42
1:C:161:LYS:O	1:C:162:ALA:CB	2.66	0.42
3:E:498:THR:HG23	3:E:503:PHE:N	2.33	0.42
3:F:338:ALA:HA	3:F:359:LEU:O	2.20	0.42
3:G:520:ASP:O	3:G:542:SER:OG	2.29	0.42
2:H:18:VAL:O	2:H:82:GLU:HA	2.18	0.42
3:G:387:ARG:HH12	3:G:389:GLU:CG	2.33	0.42
3:E:430:MET:HB3	2:H:55:ASN:HA	2.02	0.42
2:B:32:TYR:CE1	2:B:101:GLY:O	2.73	0.42
2:B:174:PHE:CD1	2:B:174:PHE:N	2.88	0.42
3:E:431:ARG:CD	2:H:74:THR:HG21	2.49	0.42
2:H:192:VAL:HG21	2:H:202:TYR:HE2	1.84	0.42
2:H:23:LYS:HD3	2:H:25:TYR:CE1	2.55	0.42
1:A:108:LEU:HD13	1:A:109:THR:N	2.36	0.41
2:B:19:LYS:HE2	2:B:80:TYR:CG	2.54	0.41
2:D:24:VAL:CG1	2:D:27:TYR:CE1	3.03	0.41
3:E:343:PRO:HB2	3:E:348:LEU:HD11	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:477:GLN:H	3:F:477:GLN:HG2	1.75	0.41
2:H:12:LYS:HB3	2:H:86:LEU:HD23	2.02	0.41
2:H:134:PRO:HB3	2:H:146:LEU:HB3	2.01	0.41
3:G:345:PRO:HG2	3:G:474:ILE:CA	2.50	0.41
3:E:479:LEU:CD1	3:E:484:GLN:HA	2.51	0.41
3:E:513:ARG:HD3	3:E:516:TRP:CE2	2.55	0.41
3:F:350:ILE:HD13	3:F:350:ILE:N	2.36	0.41
2:H:162:TRP:HB2	2:H:167:LEU:HB3	2.03	0.41
2:D:170:GLY:O	2:D:190:VAL:HA	2.20	0.41
3:G:377:ALA:HB2	3:G:415:THR:HB	2.03	0.41
3:G:408:ARG:O	3:G:412:GLU:HG3	2.21	0.41
2:H:49:GLY:HA3	2:H:70:MET:CE	2.50	0.41
2:H:55:ASN:OD1	2:H:57:GLU:HB2	2.21	0.41
2:B:55:ASN:OD1	2:B:57:GLU:HG3	2.21	0.41
3:E:466:ILE:O	3:E:504:PHE:HA	2.19	0.41
3:F:493:THR:CG2	3:F:506:PHE:H	2.33	0.41
1:L:94:ASP:O	1:L:95:SER:CB	2.69	0.41
1:A:153:TRP:HB2	1:A:160:VAL:HG22	2.01	0.41
3:E:371:ASN:O	3:E:420:VAL:HA	2.20	0.41
3:F:497:LYS:HZ1	3:F:501:SER:HA	1.82	0.41
1:L:15:GLY:C	1:L:79:GLY:HA2	2.41	0.41
1:L:63:ARG:CG	1:L:78:THR:O	2.69	0.41
1:C:140:LEU:HD13	2:D:189:VAL:HG21	2.03	0.41
3:G:544:ASN:CB	3:G:545:PRO:HD2	2.48	0.41
2:H:14:PRO:HA	2:H:86:LEU:O	2.21	0.41
1:A:12:GLY:O	1:A:13:ALA:HB3	2.21	0.41
2:B:154:PHE:CG	2:B:155:PRO:N	2.89	0.41
1:C:201:THR:HG23	1:C:206:THR:OG1	2.20	0.41
1:C:172:GLN:O	1:C:174:ASN:O	2.38	0.40
3:F:344:SER:O	3:F:345:PRO:C	2.55	0.40
2:H:87:ARG:NH2	2:H:89:GLU:OE1	2.54	0.40
3:E:344:SER:O	3:E:346:PHE:N	2.51	0.40
3:F:465:LEU:HD23	3:F:465:LEU:HA	1.94	0.40
3:G:343:PRO:HB2	3:G:348:LEU:HD22	2.02	0.40
2:H:99:VAL:C	2:H:100:MET:SD	3.00	0.40
1:A:160:VAL:CG2	1:A:161:LYS:N	2.73	0.40
1:A:161:LYS:HG2	1:A:161:LYS:O	2.22	0.40
3:F:393:ARG:HD3	3:F:393:ARG:C	2.42	0.40
3:G:371:ASN:HB2	3:G:421:THR:HG21	1.91	0.40
2:H:150:VAL:CG1	2:H:186:LEU:CD1	2.96	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	211/217 (97%)	189 (90%)	16 (8%)	6 (3%)	5	16
1	C	211/217 (97%)	189 (90%)	15 (7%)	7 (3%)	4	13
1	L	177/217 (82%)	146 (82%)	21 (12%)	10 (6%)	2	4
2	B	216/229 (94%)	204 (94%)	7 (3%)	5 (2%)	6	20
2	D	215/229 (94%)	200 (93%)	12 (6%)	3 (1%)	11	31
2	H	200/229 (87%)	174 (87%)	16 (8%)	10 (5%)	2	6
3	E	202/242 (84%)	175 (87%)	14 (7%)	13 (6%)	1	3
3	F	202/242 (84%)	181 (90%)	16 (8%)	5 (2%)	5	18
3	G	209/242 (86%)	191 (91%)	10 (5%)	8 (4%)	3	10
All	All	1843/2064 (89%)	1649 (90%)	127 (7%)	67 (4%)	3	11

All (67) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	113	GLN
1	A	146	PRO
1	A	213	PRO
2	B	155	PRO
2	B	157	PRO
1	C	113	GLN
1	C	146	PRO
1	C	174	ASN
2	D	155	PRO
2	D	157	PRO
3	E	365	PRO
3	E	366	SER
3	E	471	PRO
3	F	471	PRO
3	G	422	HIS
3	G	423	PRO

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Mol	Chain	Res	Type
3	G	471	PRO
3	G	533	PRO
2	H	9	ALA
2	H	63	LYS
2	H	65	GLN
2	H	155	PRO
2	H	157	PRO
1	L	95	SER
1	C	111	LEU
1	C	112	GLY
3	E	368	GLY
3	E	379	GLY
3	E	407	THR
3	F	533	PRO
3	G	457	ARG
2	H	166	ALA
1	L	97	THR
1	A	157	SER
2	B	102	LYS
2	B	141	GLY
2	D	199	THR
3	E	456	SER
3	E	459	LYS
3	E	533	PRO
3	G	426	PRO
3	G	482	GLU
2	H	16	ALA
1	A	13	ALA
1	C	156	ASP
3	E	346	PHE
3	E	378	SER
2	H	41	PRO
1	L	43	GLY
1	L	124	PRO
1	L	125	PRO
1	L	156	ASP
1	A	156	ASP
2	B	135	SER
1	C	162	ALA
3	F	366	SER
2	H	99	VAL
1	L	7	PRO

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Mol	Chain	Res	Type
1	L	26	SER
1	L	96	PRO
3	F	472	GLU
3	G	456	SER
3	E	406	GLY
3	F	500	GLY
2	H	131	PRO
1	L	213	PRO
3	E	532	SER

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	179/183 (98%)	165 (92%)	14 (8%)	12	32
1	C	179/183 (98%)	166 (93%)	13 (7%)	14	35
1	L	158/183 (86%)	135 (85%)	23 (15%)	3	8
2	B	184/193 (95%)	170 (92%)	14 (8%)	13	33
2	D	183/193 (95%)	172 (94%)	11 (6%)	19	45
2	H	173/193 (90%)	148 (86%)	25 (14%)	3	8
3	E	183/213 (86%)	155 (85%)	28 (15%)	2	7
3	F	182/213 (85%)	161 (88%)	21 (12%)	5	15
3	G	186/213 (87%)	166 (89%)	20 (11%)	6	17
All	All	1607/1767 (91%)	1438 (90%)	169 (10%)	7	18

All (169) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	THR
1	A	10	VAL
1	A	69	SER
1	A	74	SER
1	A	108	LEU

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Mol	Chain	Res	Type
1	A	127	SER
1	A	146	PRO
1	A	150	THR
1	A	168	THR
1	A	172	GLN
1	A	180	SER
1	A	185	LEU
1	A	203	GLU
1	A	210	THR
2	B	13	LYS
2	B	21	SER
2	B	100	MET
2	B	102	LYS
2	B	103	TRP
2	B	132	LEU
2	B	146	LEU
2	B	157	PRO
2	B	159	THR
2	B	168	THR
2	B	169	SER
2	B	172	HIS
2	B	186	LEU
2	B	209	LYS
1	C	19	THR
1	C	44	THR
1	C	56	ARG
1	C	58	SER
1	C	74	SER
1	C	127	SER
1	C	134	LYS
1	C	146	PRO
1	C	150	THR
1	C	158	SER
1	C	180	SER
1	C	191	LYS
1	C	205	SER
2	D	1	GLU
2	D	45	LEU
2	D	63	LYS
2	D	86	LEU
2	D	99	VAL
2	D	113	ARG

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Mol	Chain	Res	Type
2	D	156	GLU
2	D	168	THR
2	D	190	VAL
2	D	199	THR
2	D	223	SER
3	E	339	TYR
3	E	340	LEU
3	E	344	SER
3	E	363	LEU
3	E	370	VAL
3	E	371	ASN
3	E	378	SER
3	E	392	GLN
3	E	393	ARG
3	E	401	SER
3	E	407	THR
3	E	419	ARG
3	E	421	THR
3	E	437	SER
3	E	440	ARG
3	E	452	GLU
3	E	457	ARG
3	E	485	LEU
3	E	489	ARG
3	E	494	GLN
3	E	508	ARG
3	E	511	VAL
3	E	513	ARG
3	E	518	GLN
3	E	525	ARG
3	E	534	SER
3	E	535	GLN
3	E	541	VAL
3	F	355	THR
3	F	363	LEU
3	F	367	LYS
3	F	369	THR
3	F	393	ARG
3	F	421	THR
3	F	427	ARG
3	F	430	MET
3	F	459	LYS

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Mol	Chain	Res	Type
3	F	462	LEU
3	F	465	LEU
3	F	477	GLN
3	F	479	LEU
3	F	483	VAL
3	F	485	LEU
3	F	492	THR
3	F	493	THR
3	F	498	THR
3	F	518	GLN
3	F	532	SER
3	F	536	THR
3	G	348	LEU
3	G	363	LEU
3	G	369	THR
3	G	370	VAL
3	G	372	LEU
3	G	387	ARG
3	G	393	ARG
3	G	411	ILE
3	G	421	THR
3	G	429	LEU
3	G	430	MET
3	G	440	ARG
3	G	450	THR
3	G	457	ARG
3	G	460	ARG
3	G	485	LEU
3	G	497	LYS
3	G	505	VAL
3	G	516	TRP
3	G	539	ARG
2	H	17	THR
2	H	18	VAL
2	H	30	THR
2	H	43	LYS
2	H	45	LEU
2	H	55	ASN
2	H	65	GLN
2	H	74	THR
2	H	78	ARG
2	H	83	LEU

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Mol	Chain	Res	Type
2	H	87	ARG
2	H	89	GLU
2	H	91	THR
2	H	99	VAL
2	H	113	ARG
2	H	117	VAL
2	H	148	CYS
2	H	159	THR
2	H	167	LEU
2	H	186	LEU
2	H	199	THR
2	H	201	THR
2	H	204	CYS
2	H	206	VAL
2	H	217	LYS
1	L	5	THR
1	L	9	SER
1	L	21	SER
1	L	23	THR
1	L	34	ASP
1	L	67	SER
1	L	77	ILE
1	L	87	ASP
1	L	93	TYR
1	L	99	THR
1	L	109	THR
1	L	110	VAL
1	L	122	LEU
1	L	141	ILE
1	L	142	SER
1	L	168	THR
1	L	171	LYS
1	L	180	SER
1	L	184	SER
1	L	188	GLU
1	L	194	ARG
1	L	208	GLU
1	L	214	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	GLN
3	E	535	GLN
3	F	477	GLN
3	G	392	GLN
3	G	480	HIS
3	G	535	GLN
1	L	6	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 ⓘ

17 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	$\# Z > 2$
4	NAG	I	1	3,4	14,14,15	0.86	1 (7%)	17,19,21	1.05	0
4	NAG	I	2	4	14,14,15	0.71	0	17,19,21	1.45	2 (11%)
4	BMA	I	3	4	11,11,12	0.75	0	15,15,17	1.68	5 (33%)
4	MAN	I	4	4	11,11,12	0.71	0	15,15,17	1.69	4 (26%)
4	MAN	I	5	4	11,11,12	0.61	0	15,15,17	0.96	0
4	MAN	I	6	4	11,11,12	0.89	1 (9%)	15,15,17	1.70	3 (20%)
5	NAG	J	1	3,5	14,14,15	0.99	1 (7%)	17,19,21	1.39	2 (11%)
5	NAG	J	2	5	14,14,15	0.84	1 (7%)	17,19,21	2.61	9 (52%)
5	BMA	J	3	5	11,11,12	0.85	0	15,15,17	1.67	3 (20%)
5	BMA	J	4	5	11,11,12	1.02	1 (9%)	15,15,17	1.91	4 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	MAN	J	5	5	11,11,12	0.99	1 (9%)	15,15,17	1.78	3 (20%)
6	NAG	K	1	3,6	14,14,15	1.38	2 (14%)	17,19,21	2.17	4 (23%)
6	NAG	K	2	6	14,14,15	0.89	1 (7%)	17,19,21	2.14	4 (23%)
6	BMA	K	3	6	11,11,12	1.13	0	15,15,17	2.62	5 (33%)
6	MAN	K	4	6	11,11,12	0.72	0	15,15,17	2.93	4 (26%)
6	MAN	K	5	6	11,11,12	0.82	0	15,15,17	3.24	6 (40%)
6	BMA	K	6	6	11,11,12	0.60	0	15,15,17	1.54	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
4	NAG	I	1	3,4	-	0/6/23/26	0/1/1/1
4	NAG	I	2	4	-	0/6/23/26	0/1/1/1
4	BMA	I	3	4	-	2/2/19/22	0/1/1/1
4	MAN	I	4	4	-	0/2/19/22	0/1/1/1
4	MAN	I	5	4	-	0/2/19/22	0/1/1/1
4	MAN	I	6	4	-	0/2/19/22	0/1/1/1
5	NAG	J	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	J	2	5	-	2/6/23/26	0/1/1/1
5	BMA	J	3	5	-	2/2/19/22	0/1/1/1
5	BMA	J	4	5	-	2/2/19/22	0/1/1/1
5	MAN	J	5	5	-	0/2/19/22	0/1/1/1
6	NAG	K	1	3,6	-	2/6/23/26	0/1/1/1
6	NAG	K	2	6	-	0/6/23/26	0/1/1/1
6	BMA	K	3	6	-	0/2/19/22	0/1/1/1
6	MAN	K	4	6	-	2/2/19/22	0/1/1/1
6	MAN	K	5	6	-	2/2/19/22	1/1/1/1
6	BMA	K	6	6	-	2/2/19/22	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	K	1	NAG	O5-C1	-3.79	1.37	1.43
6	K	2	NAG	O5-C1	-2.54	1.39	1.43
5	J	4	BMA	C2-C3	2.51	1.56	1.52
5	J	5	MAN	C2-C3	2.30	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	K	1	NAG	O5-C5	-2.22	1.39	1.43
5	J	2	NAG	C2-N2	-2.18	1.42	1.46
4	I	1	NAG	O5-C1	-2.04	1.40	1.43
5	J	1	NAG	O5-C1	-2.03	1.40	1.43
4	I	6	MAN	C2-C3	2.01	1.55	1.52

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	K	5	MAN	C1-C2-C3	-10.52	96.73	109.67
6	K	4	MAN	C1-O5-C5	7.85	122.83	112.19
6	K	2	NAG	O5-C1-C2	-6.91	100.37	111.29
6	K	4	MAN	C1-C2-C3	-6.21	102.03	109.67
6	K	3	BMA	O3-C3-C2	5.94	121.37	109.99
5	J	2	NAG	C2-N2-C7	4.97	129.97	122.90
5	J	4	BMA	C1-C2-C3	4.94	115.74	109.67
5	J	2	NAG	C1-O5-C5	4.81	118.71	112.19
6	K	3	BMA	O5-C5-C6	4.59	114.39	107.20
6	K	1	NAG	C1-C2-N2	-4.54	102.73	110.49
6	K	3	BMA	C3-C4-C5	-4.37	102.44	110.24
5	J	5	MAN	C1-C2-C3	4.32	114.97	109.67
6	K	1	NAG	C3-C4-C5	-4.28	102.60	110.24
6	K	6	BMA	C3-C4-C5	4.10	117.56	110.24
4	I	6	MAN	C3-C4-C5	4.08	117.52	110.24
6	K	1	NAG	C6-C5-C4	3.80	121.91	113.00
5	J	2	NAG	O5-C1-C2	-3.73	105.40	111.29
4	I	2	NAG	C1-O5-C5	3.69	117.20	112.19
4	I	4	MAN	C1-C2-C3	3.56	114.04	109.67
5	J	3	BMA	C1-C2-C3	3.55	114.03	109.67
4	I	6	MAN	C2-C3-C4	3.44	116.84	110.89
5	J	4	BMA	O5-C5-C6	3.36	112.48	107.20
5	J	2	NAG	C1-C2-N2	3.30	116.13	110.49
5	J	2	NAG	C8-C7-N2	3.19	121.50	116.10
6	K	5	MAN	C1-O5-C5	-3.16	107.91	112.19
4	I	3	BMA	C1-O5-C5	3.16	116.47	112.19
5	J	3	BMA	C1-O5-C5	3.12	116.41	112.19
6	K	5	MAN	O5-C5-C6	3.08	112.03	107.20
6	K	3	BMA	C1-O5-C5	-3.07	108.04	112.19
6	K	2	NAG	O5-C5-C6	3.05	111.98	107.20
4	I	4	MAN	C2-C3-C4	3.01	116.10	110.89
5	J	5	MAN	O5-C5-C6	2.99	111.89	107.20
5	J	1	NAG	O5-C1-C2	-2.95	106.62	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	J	2	NAG	O7-C7-N2	-2.82	116.77	121.95
4	I	3	BMA	C1-C2-C3	2.81	113.12	109.67
6	K	5	MAN	O2-C2-C3	2.80	115.75	110.14
5	J	2	NAG	O4-C4-C3	-2.78	103.93	110.35
4	I	2	NAG	O4-C4-C3	-2.76	103.96	110.35
4	I	3	BMA	O3-C3-C4	-2.72	104.07	110.35
6	K	1	NAG	O5-C1-C2	-2.68	107.06	111.29
5	J	5	MAN	C2-C3-C4	2.63	115.44	110.89
6	K	4	MAN	O5-C5-C4	2.49	116.89	110.83
6	K	4	MAN	O3-C3-C4	2.49	116.10	110.35
5	J	2	NAG	C3-C4-C5	-2.47	105.83	110.24
4	I	4	MAN	C1-O5-C5	2.45	115.51	112.19
6	K	2	NAG	O7-C7-N2	2.43	126.41	121.95
6	K	5	MAN	O5-C1-C2	-2.38	107.09	110.77
5	J	1	NAG	C3-C4-C5	-2.37	106.02	110.24
5	J	3	BMA	O5-C5-C6	2.34	110.87	107.20
6	K	2	NAG	O7-C7-C8	-2.25	117.87	122.06
5	J	4	BMA	O5-C5-C4	-2.24	105.38	110.83
4	I	4	MAN	O5-C5-C6	2.21	110.68	107.20
5	J	4	BMA	C2-C3-C4	2.18	114.67	110.89
5	J	2	NAG	O4-C4-C5	2.17	114.69	109.30
6	K	5	MAN	C6-C5-C4	2.17	118.08	113.00
4	I	6	MAN	O2-C2-C3	2.17	114.47	110.14
6	K	3	BMA	O5-C5-C4	-2.12	105.68	110.83
4	I	3	BMA	O4-C4-C5	2.10	114.50	109.30
6	K	6	BMA	C2-C3-C4	2.05	114.45	110.89
4	I	3	BMA	O4-C4-C3	-2.04	105.64	110.35

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	J	1	NAG	C4-C5-C6-O6
6	K	4	MAN	C4-C5-C6-O6
6	K	6	BMA	O5-C5-C6-O6
6	K	4	MAN	O5-C5-C6-O6
6	K	5	MAN	C4-C5-C6-O6
5	J	1	NAG	O5-C5-C6-O6
5	J	4	BMA	O5-C5-C6-O6
6	K	1	NAG	C4-C5-C6-O6
4	I	3	BMA	O5-C5-C6-O6
5	J	4	BMA	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
5	J	2	NAG	C8-C7-N2-C2
5	J	2	NAG	O7-C7-N2-C2
6	K	5	MAN	O5-C5-C6-O6
5	J	3	BMA	O5-C5-C6-O6
4	I	3	BMA	C4-C5-C6-O6
5	J	3	BMA	C4-C5-C6-O6
6	K	1	NAG	O5-C5-C6-O6
6	K	6	BMA	C4-C5-C6-O6

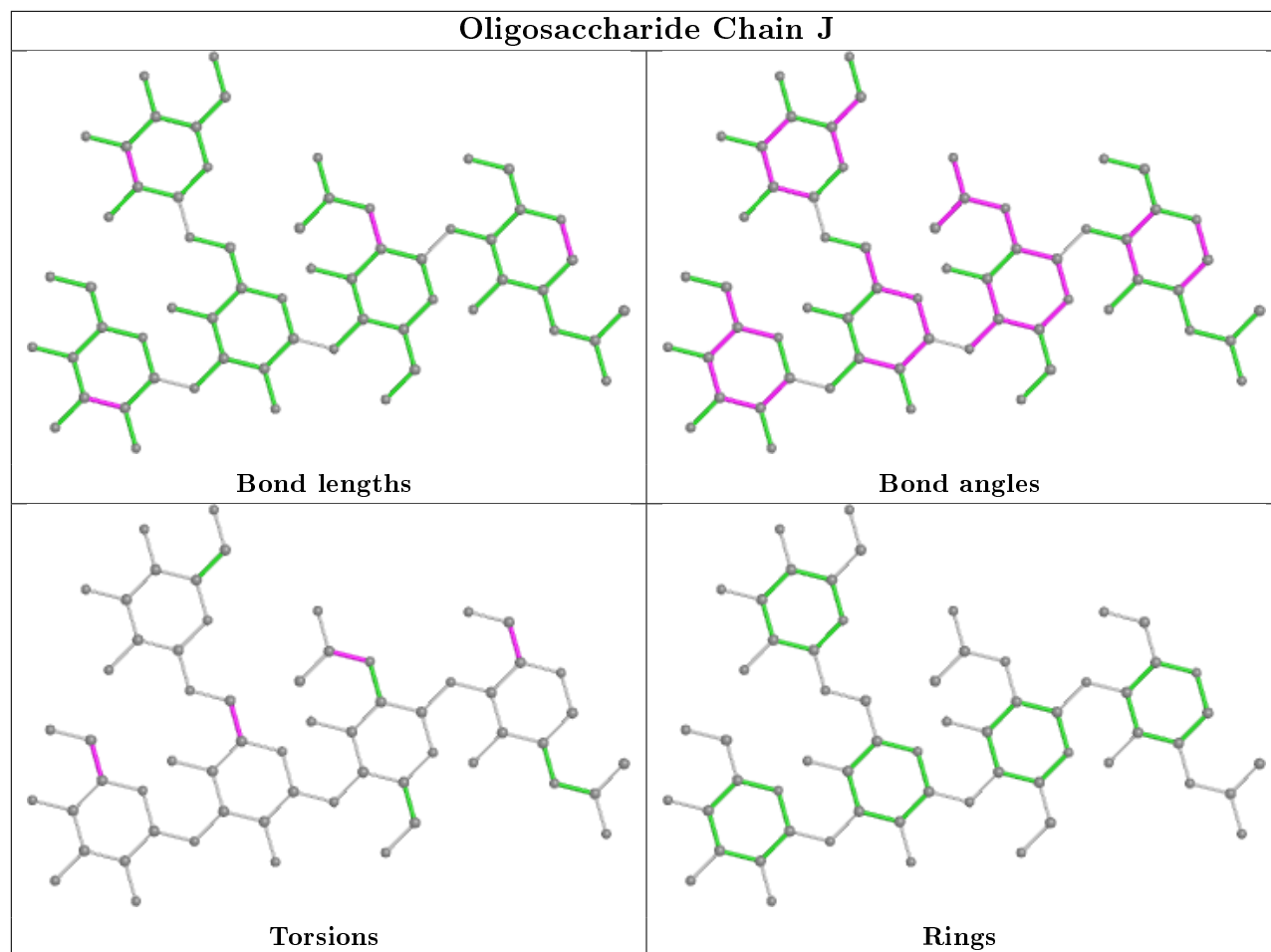
All (1) ring outliers are listed below:

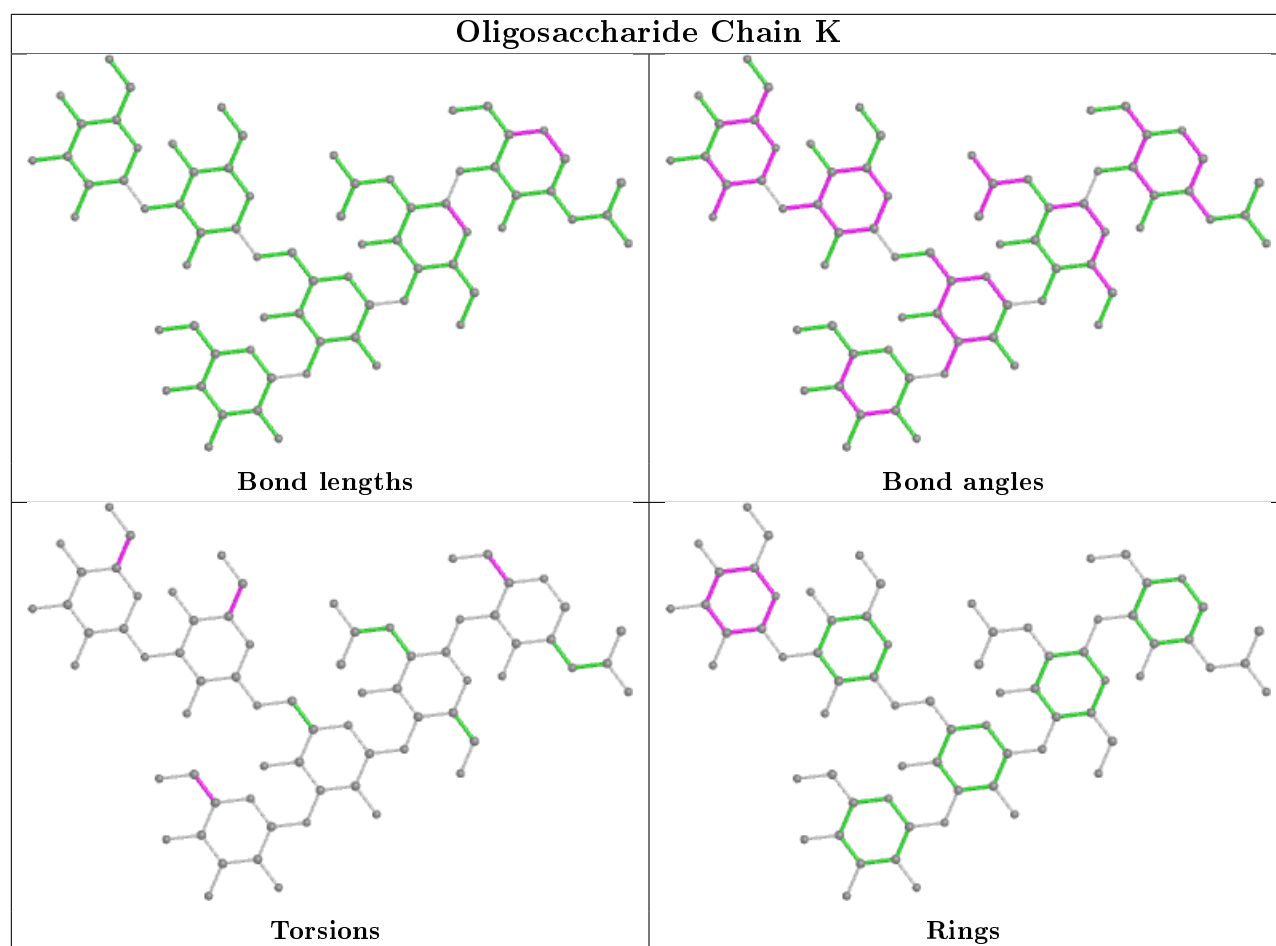
Mol	Chain	Res	Type	Atoms
6	K	5	MAN	C1-C2-C3-C4-C5-O5

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	J	1	NAG	1	0
6	K	3	BMA	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.





5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	213/217 (98%)	-0.40	0 100 100	30, 56, 83, 105	0
1	C	213/217 (98%)	-0.43	0 100 100	31, 56, 90, 108	0
1	L	189/217 (87%)	0.23	8 (4%) 36 31	79, 118, 147, 159	0
2	B	220/229 (96%)	-0.50	3 (1%) 75 74	25, 35, 92, 148	0
2	D	219/229 (95%)	-0.50	1 (0%) 91 90	25, 37, 69, 110	0
2	H	208/229 (90%)	0.14	9 (4%) 35 30	60, 89, 127, 157	0
3	E	206/242 (85%)	0.15	10 (4%) 29 25	54, 92, 137, 165	0
3	F	208/242 (85%)	-0.17	4 (1%) 66 64	31, 67, 126, 153	0
3	G	211/242 (87%)	-0.33	6 (2%) 53 48	28, 47, 98, 161	0
All	All	1887/2064 (91%)	-0.21	41 (2%) 62 59	25, 62, 131, 165	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	546	GLY	7.0
3	E	426	PRO	5.4
1	L	159	PRO	4.7
3	G	424	HIS	4.6
3	G	545	PRO	4.2
2	H	218	LYS	4.2
3	E	532	SER	4.1
1	L	82	ALA	3.7
3	E	424	HIS	3.5
2	H	202	TYR	3.5
2	B	141	GLY	3.5
2	D	103	TRP	3.4
2	H	220	GLU	3.3
2	H	201	THR	3.2
3	G	366	SER	2.9

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Mol	Chain	Res	Type	RSRZ
1	L	186	THR	2.8
3	F	369	THR	2.7
2	B	136	SER	2.6
1	L	193	HIS	2.6
1	L	68	LYS	2.6
3	E	452	GLU	2.6
1	L	154	LYS	2.5
1	L	25	SER	2.4
3	E	481	ASN	2.4
2	H	166	ALA	2.4
3	E	514	ALA	2.4
3	E	378	SER	2.3
2	H	192	VAL	2.3
3	G	369	THR	2.3
3	G	367	LYS	2.3
3	F	367	LYS	2.3
2	H	191	THR	2.3
2	B	137	LYS	2.3
2	H	62	GLU	2.2
2	H	200	GLN	2.2
3	E	427	ARG	2.2
3	G	422	HIS	2.1
3	E	379	GLY	2.1
3	E	455	GLY	2.1
1	L	32	GLY	2.1
3	F	423	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

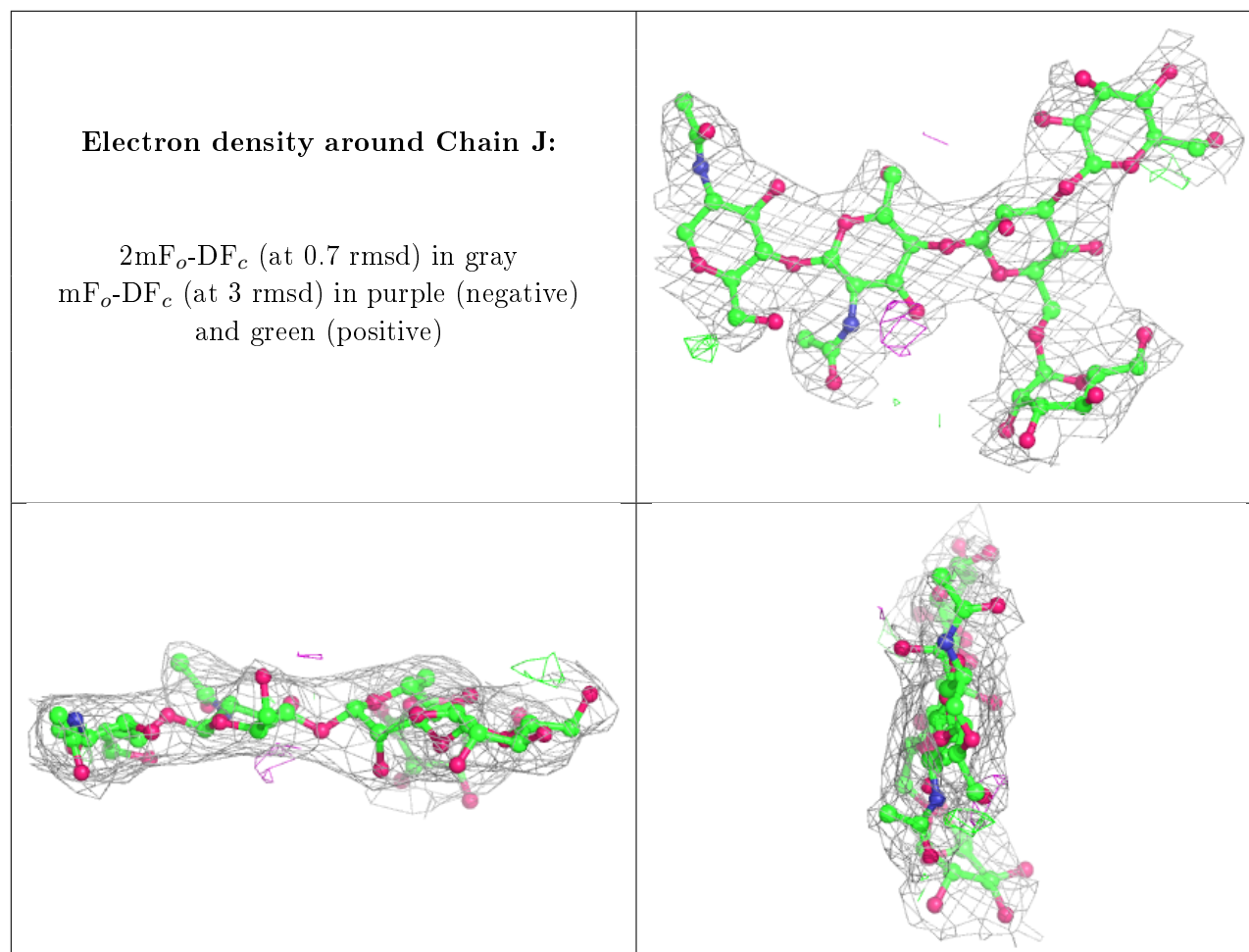
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MAN	I	5	11/12	0.73	0.37	141,147,150,151	0
5	BMA	J	4	11/12	0.73	0.25	85,103,107,114	0

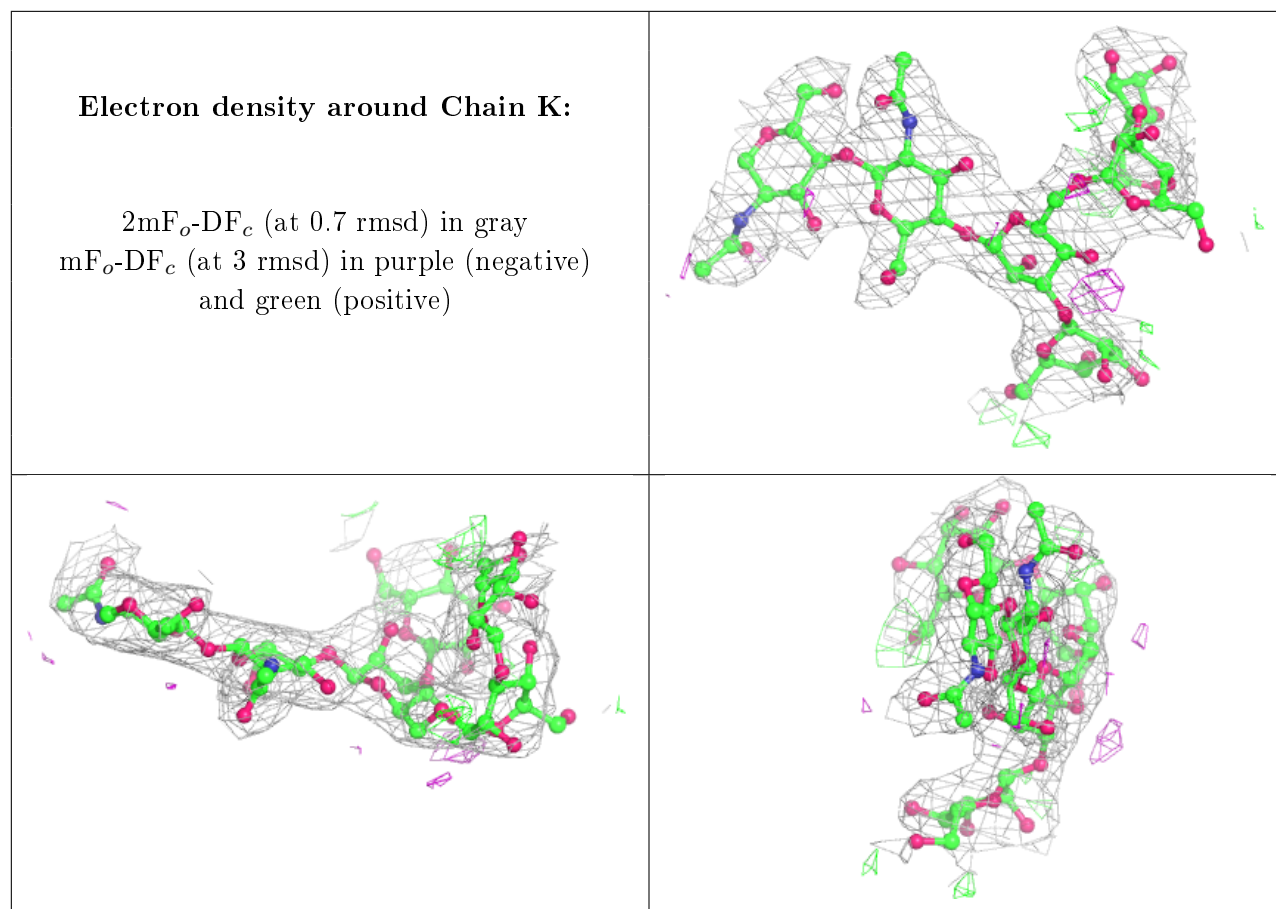
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	BMA	K	6	11/12	0.80	0.18	99,111,121,134	0
5	MAN	J	5	11/12	0.83	0.18	91,95,102,106	0
6	MAN	K	5	11/12	0.84	0.23	83,100,111,119	0
4	MAN	I	6	11/12	0.86	0.18	89,103,107,109	0
4	MAN	I	4	11/12	0.86	0.16	101,117,126,143	0
6	MAN	K	4	11/12	0.93	0.14	81,83,94,97	0
5	BMA	J	3	11/12	0.95	0.10	64,72,81,87	0
4	NAG	I	2	14/15	0.96	0.12	59,62,69,71	0
4	BMA	I	3	11/12	0.96	0.09	74,85,96,107	0
5	NAG	J	1	14/15	0.96	0.17	60,62,68,76	0
5	NAG	J	2	14/15	0.97	0.13	56,63,68,70	0
6	BMA	K	3	11/12	0.97	0.12	51,57,70,76	0
4	NAG	I	1	14/15	0.98	0.11	47,50,54,55	0
6	NAG	K	2	14/15	0.98	0.14	37,41,44,46	0
6	NAG	K	1	14/15	0.98	0.15	34,37,38,39	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 ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.